## GNU Octave

A high-level interactive language for numerical computations
Edition 7 for Octave version 7.3.0
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Free Your Numbers

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## Table of Contents

Preface ..... 1
Acknowledgements ..... 1
Citing Octave in Publications ..... 5
How You Can Contribute to Octave ..... 6
Distribution ..... 6
1 A Brief Introduction to Octave ..... 7
1.1 Running Octave ..... 7
1.2 Simple Examples ..... 7
1.2.1 Elementary Calculations ..... 7
1.2.2 Creating a Matrix ..... 8
1.2.3 Matrix Arithmetic ..... 8
1.2.4 Solving Systems of Linear Equations ..... 8
1.2.5 Integrating Differential Equations ..... 9
1.2.6 Producing Graphical Output ..... 10
1.2.7 Help and Documentation ..... 10
1.2.8 Editing What You Have Typed ..... 10
1.3 Conventions ..... 11
1.3.1 Fonts ..... 11
1.3.2 Evaluation Notation ..... 11
1.3.3 Printing Notation ..... 11
1.3.4 Error Messages ..... 12
1.3.5 Format of Descriptions ..... 12
1.3.5.1 A Sample Function Description ..... 12
1.3.5.2 A Sample Command Description ..... 13
2 Getting Started ..... 15
2.1 Invoking Octave from the Command Line ..... 15
2.1.1 Command Line Options ..... 15
2.1.2 Startup Files ..... 18
2.2 Quitting Octave ..... 19
2.3 Commands for Getting Help ..... 21
2.4 Command Line Editing ..... 25
2.4.1 Cursor Motion ..... 26
2.4.2 Killing and Yanking ..... 27
2.4.3 Commands for Changing Text ..... 27
2.4.4 Letting Readline Type for You ..... 28
2.4.5 Commands for Manipulating the History ..... 28
2.4.6 Customizing readline ..... 32
2.4.7 Customizing the Prompt ..... 33
2.4.8 Diary and Echo Commands ..... 34
2.5 How Octave Reports Errors ..... 36
2.6 Executable Octave Programs ..... 37
2.6.1 Passing Arguments to Executable Scripts ..... 38
2.6.2 Dual-Purpose Executable Scripts and Octave Functions ..... 38
2.7 Comments in Octave Programs ..... 38
2.7.1 Single Line Comments ..... 38
2.7.2 Block Comments ..... 39
2.7.3 Comments and the Help System ..... 39
3 Data Types ..... 41
3.1 Built-in Data Types ..... 41
3.1.1 Numeric Objects ..... 44
3.1.2 Missing Data ..... 45
3.1.3 String Objects ..... 45
3.1.4 Data Structure Objects ..... 46
3.1.5 Cell Array Objects ..... 46
3.2 User-defined Data Types ..... 46
3.3 Object Sizes ..... 46
4 Numeric Data Types ..... 51
4.1 Matrices ..... 52
4.1.1 Empty Matrices ..... 55
4.2 Ranges ..... 56
4.3 Single Precision Data Types ..... 57
4.4 Integer Data Types ..... 58
4.4.1 Integer Arithmetic ..... 61
4.5 Bit Manipulations ..... 62
4.6 Logical Values. ..... 64
4.7 Automatic Conversion of Data Types ..... 65
4.8 Predicates for Numeric Objects ..... 67
5 Strings ..... 71
5.1 Escape Sequences in String Constants ..... 71
5.2 Character Arrays ..... 72
5.3 String Operations ..... 74
5.3.1 Common String Operations ..... 75
5.3.2 Concatenating Strings ..... 77
5.3.3 Splitting and Joining Strings ..... 80
5.3.4 Searching in Strings ..... 84
5.3.5 Searching and Replacing in Strings ..... 90
5.4 Converting Strings ..... 94
5.4.1 String encoding ..... 94
5.4.2 Numerical Data and Strings ..... 95
5.4.3 JSON data encoding/decoding ..... 104
5.5 Character Class Functions ..... 108
6 Data Containers ..... 111
6.1 Structures ..... 111
6.1.1 Basic Usage and Examples ..... 111
6.1.2 Structure Arrays ..... 115
6.1.3 Creating Structures ..... 116
6.1.4 Manipulating Structures ..... 119
6.1.5 Processing Data in Structures ..... 123
6.2 containers.Map. ..... 124
6.3 Cell Arrays ..... 124
6.3.1 Basic Usage of Cell Arrays ..... 125
6.3.2 Creating Cell Arrays ..... 126
6.3.3 Indexing Cell Arrays ..... 130
6.3.4 Cell Arrays of Strings ..... 133
6.3.5 Processing Data in Cell Arrays ..... 134
6.4 Comma Separated Lists ..... 134
6.4.1 Comma Separated Lists Generated from Cell Arrays ..... 135
6.4.2 Comma Separated Lists Generated from Structure Arrays ..... 136
7 Variables ..... 137
7.1 Global Variables ..... 139
7.2 Persistent Variables ..... 140
7.3 Status of Variables ..... 142
8 Expressions ..... 151
8.1 Index Expressions ..... 151
8.1.1 Advanced Indexing ..... 153
8.2 Calling Functions ..... 157
8.2.1 Call by Value ..... 158
8.2.2 Recursion ..... 159
8.2.3 Access via Handle ..... 160
8.3 Arithmetic Operators ..... 161
8.4 Comparison Operators ..... 164
8.5 Boolean Expressions. ..... 166
8.5.1 Element-by-element Boolean Operators ..... 166
8.5.2 Short-circuit Boolean Operators ..... 167
8.6 Assignment Expressions ..... 169
8.7 Increment Operators ..... 171
8.8 Operator Precedence ..... 172
9 Evaluation ..... 175
9.1 Calling a Function by its Name ..... 176
9.2 Evaluation in a Different Context ..... 177
10 Statements ..... 179
10.1 The if Statement ..... 179
10.2 The switch Statement ..... 181
10.2.1 Notes for the C Programmer ..... 182
10.3 The while Statement ..... 183
10.4 The do-until Statement ..... 184
10.5 The for Statement ..... 184
10.5.1 Looping Over Structure Elements ..... 185
10.6 The break Statement ..... 186
10.7 The continue Statement ..... 187
10.8 The unwind_protect Statement ..... 188
10.9 The try Statement ..... 188
10.10 Continuation Lines ..... 189
11 Functions and Scripts ..... 191
11.1 Introduction to Function and Script Files ..... 191
11.2 Defining Functions ..... 191
11.3 Returning from a Function ..... 194
11.4 Multiple Return Values ..... 195
11.5 Variable-length Return Lists ..... 198
11.6 Variable-length Argument Lists ..... 199
11.7 Ignoring Arguments ..... 200
11.8 Default Arguments ..... 202
11.9 Validating Arguments ..... 202
11.9.1 Validating the number of Arguments ..... 202
11.9.2 Validating the type of Arguments ..... 203
11.9.3 Parsing Arguments ..... 208
11.10 Function Files ..... 210
11.10.1 Manipulating the Load Path ..... 213
11.10.2 Subfunctions ..... 217
11.10.3 Private Functions ..... 217
11.10.4 Nested Functions ..... 218
11.10.5 Overloading and Autoloading ..... 220
11.10.6 Function Locking ..... 221
11.10.7 Function Precedence ..... 223
11.11 Script Files ..... 223
11.11.1 Publish Octave Script Files ..... 225
11.11.2 Publishing Markup ..... 227
11.11.2.1 Using Publishing Markup in Script Files ..... 227
11.11.2.2 Text Formatting ..... 228
11.11.2.3 Sections ..... 228
11.11.2.4 Preformatted Code ..... 229
11.11.2.5 Preformatted Text ..... 229
11.11.2.6 Bulleted Lists ..... 230
11.11.2.7 Numbered Lists ..... 230
11.11.2.8 Including File Content ..... 230
11.11.2.9 Including Graphics ..... 230
11.11.2.10 Including URLs ..... 231
11.11.2.11 Mathematical Equations ..... 231
11.11.2.12 HTML Markup ..... 231
11.11.2.13 LaTeX Markup ..... 232
11.11.3 Jupyter Notebooks ..... 232
11.12 Function Handles and Anonymous Functions ..... 233
11.12.1 Function Handles ..... 233
11.12.2 Anonymous Functions ..... 235
11.13 Command Syntax and Function Syntax ..... 236
11.14 Organization of Functions Distributed with Octave ..... 236
12 Errors and Warnings ..... 239
12.1 Handling Errors ..... 239
12.1.1 Raising Errors ..... 239
12.1.2 Catching Errors ..... 242
12.1.3 Recovering From Errors ..... 245
12.2 Handling Warnings ..... 245
12.2.1 Issuing Warnings ..... 246
12.2.2 Enabling and Disabling Warnings ..... 254
13 Debugging ..... 257
13.1 Entering Debug Mode ..... 257
13.2 Leaving Debug Mode. ..... 258
13.3 Breakpoints ..... 258
13.4 Debug Mode ..... 262
13.5 Call Stack ..... 263
13.6 Profiling ..... 264
13.7 Profiler Example ..... 266
14 Input and Output ..... 271
14.1 Basic Input and Output ..... 271
14.1.1 Terminal Output ..... 271
14.1.1.1 Paging Screen Output ..... 275
14.1.2 Terminal Input ..... 276
14.1.3 Simple File I/O ..... 278
14.1.3.1 Saving Data on Unexpected Exits ..... 292
14.2 C-Style I/O Functions ..... 294
14.2.1 Opening and Closing Files ..... 294
14.2.2 Simple Output ..... 297
14.2.3 Line-Oriented Input ..... 297
14.2.4 Formatted Output ..... 298
14.2.5 Output Conversion for Matrices ..... 300
14.2.6 Output Conversion Syntax ..... 300
14.2.7 Table of Output Conversions ..... 301
14.2.8 Integer Conversions ..... 302
14.2.9 Floating-Point Conversions ..... 303
14.2.10 Other Output Conversions ..... 303
14.2.11 Formatted Input ..... 304
14.2.12 Input Conversion Syntax ..... 306
14.2.13 Table of Input Conversions ..... 306
14.2.14 Numeric Input Conversions ..... 307
14.2.15 String Input Conversions ..... 308
14.2.16 Binary I/O ..... 308
14.2.17 Temporary Files ..... 311
14.2.18 End of File and Errors ..... 312
14.2.19 File Positioning ..... 313
15 Plotting ..... 315
15.1 Introduction to Plotting ..... 315
15.2 High-Level Plotting ..... 315
15.2.1 Two-Dimensional Plots ..... 315
15.2.1.1 Axis Configuration ..... 348
15.2.1.2 Two-dimensional Function Plotting ..... 353
15.2.1.3 Two-dimensional Geometric Shapes ..... 356
15.2.2 Three-Dimensional Plots ..... 357
15.2.2.1 Aspect Ratio ..... 383
15.2.2.2 Three-dimensional Function Plotting. ..... 384
15.2.2.3 Three-dimensional Geometric Shapes ..... 388
15.2.3 Plot Annotations ..... 389
15.2.4 Multiple Plots on One Page ..... 398
15.2.5 Multiple Plot Windows ..... 399
15.2.6 Manipulation of Plot Objects ..... 400
15.2.7 Manipulation of Plot Windows ..... 401
15.2.8 Use of the interpreter Property ..... 406
15.2.8.1 none interpreter ..... 406
15.2.8.2 tex interpreter ..... 406
15.2.8.3 latex interpreter ..... 408
15.2.9 Printing and Saving Plots ..... 409
15.2.10 Interacting with Plots ..... 419
15.2.11 Test Plotting Functions ..... 420
15.3 Graphics Data Structures ..... 421
15.3.1 Introduction to Graphics Structures ..... 421
15.3.2 Graphics Objects ..... 423
15.3.2.1 Creating Graphics Objects ..... 423
15.3.2.2 Handle Functions ..... 427
15.3.3 Graphics Object Properties ..... 433
15.3.3.1 Root Properties ..... 433
15.3.3.2 Figure Properties ..... 434
15.3.3.3 Axes Properties ..... 441
15.3.3.4 Legend Properties ..... 449
15.3.3.5 Line Properties ..... 450
15.3.3.6 Text Properties ..... 453
15.3.3.7 Image Properties ..... 455
15.3.3.8 Patch Properties ..... 457
15.3.3.9 Scatter Properties ..... 461
15.3.3.10 Surface Properties ..... 465
15.3.3.11 Light Properties ..... 468
15.3.3.12 Uimenu Properties ..... 470
15.3.3.13 Uibuttongroup Properties ..... 472
15.3.3.14 Uicontextmenu Properties ..... 474
15.3.3.15 Uipanel Properties ..... 476
15.3.3.16 Uicontrol Properties ..... 478
15.3.3.17 Uitable Properties ..... 481
15.3.3.18 Uitoolbar Properties ..... 483
15.3.3.19 Uipushtool Properties ..... 485
15.3.3.20 Uitoggletool Properties ..... 486
15.3.4 Searching Properties ..... 488
15.3.5 Managing Default Properties ..... 489
15.4 Advanced Plotting ..... 491
15.4.1 Colors ..... 491
15.4.2 Line Styles ..... 492
15.4.3 Marker Styles ..... 492
15.4.4 Callbacks ..... 492
15.4.5 Application-defined Data ..... 494
15.4.6 Object Groups ..... 495
15.4.6.1 Data Sources in Object Groups ..... 500
15.4.6.2 Area Series ..... 500
15.4.6.3 Bar Series ..... 501
15.4.6.4 Contour Groups ..... 502
15.4.6.5 Error Bar Series ..... 503
15.4.6.6 Line Series ..... 504
15.4.6.7 Quiver Group ..... 504
15.4.6.8 Stair Group ..... 505
15.4.6.9 Stem Series ..... 506
15.4.6.10 Surface Group ..... 506
15.4.7 Transform Groups ..... 507
15.4.8 Graphics Toolkits ..... 507
15.4.8.1 Customizing Toolkit Behavior ..... 508
15.4.8.2 Hardware vs. Software Rendering ..... 508
15.4.8.3 Precision issues ..... 509
16 Matrix Manipulation ..... 511
16.1 Finding Elements and Checking Conditions ..... 511
16.2 Rearranging Matrices ..... 515
16.3 Special Utility Matrices ..... 525
16.4 Famous Matrices ..... 537
17 Arithmetic ..... 547
17.1 Exponents and Logarithms ..... 547
17.2 Complex Arithmetic ..... 549
17.3 Trigonometry ..... 550
17.4 Sums and Products ..... 555
17.5 Utility Functions ..... 557
17.6 Special Functions ..... 565
17.7 Rational Approximations ..... 577
17.8 Coordinate Transformations ..... 578
17.9 Mathematical Constants ..... 580
18 Linear Algebra ..... 585
18.1 Techniques Used for Linear Algebra ..... 585
18.2 Basic Matrix Functions ..... 585
18.3 Matrix Factorizations ..... 595
18.4 Functions of a Matrix ..... 609
18.5 Specialized Solvers ..... 611
19 Vectorization and Faster Code Execution ..... 625
19.1 Basic Vectorization. ..... 625
19.2 Broadcasting ..... 627
19.2.1 Broadcasting and Legacy Code ..... 630
19.3 Function Application ..... 631
19.4 Accumulation ..... 636
19.5 Miscellaneous Techniques ..... 638
19.6 Examples ..... 640
20 Nonlinear Equations ..... 641
20.1 Solvers ..... 641
20.2 Minimizers ..... 645
21 Diagonal and Permutation Matrices ..... 651
21.1 Creating and Manipulating Diagonal/Permutation Matrices. ..... 651
21.1.1 Creating Diagonal Matrices. ..... 652
21.1.2 Creating Permutation Matrices ..... 652
21.1.3 Explicit and Implicit Conversions ..... 653
21.2 Linear Algebra with Diagonal/Permutation Matrices ..... 654
21.2.1 Expressions Involving Diagonal Matrices ..... 654
21.2.2 Expressions Involving Permutation Matrices ..... 655
21.3 Functions That Are Aware of These Matrices ..... 656
21.3.1 Diagonal Matrix Functions ..... 656
21.3.2 Permutation Matrix Functions ..... 656
21.4 Examples of Usage ..... 656
21.5 Differences in Treatment of Zero Elements ..... 657
22 Sparse Matrices ..... 659
22.1 Creation and Manipulation of Sparse Matrices ..... 659
22.1.1 Storage of Sparse Matrices ..... 659
22.1.2 Creating Sparse Matrices ..... 660
22.1.3 Finding Information about Sparse Matrices ..... 666
22.1.4 Basic Operators and Functions on Sparse Matrices ..... 670
22.1.4.1 Sparse Functions ..... 670
22.1.4.2 Return Types of Operators and Functions ..... 670
22.1.4.3 Mathematical Considerations ..... 672
22.2 Linear Algebra on Sparse Matrices ..... 680
22.3 Iterative Techniques Applied to Sparse Matrices ..... 690
22.4 Real Life Example using Sparse Matrices ..... 698
23 Numerical Integration ..... 703
23.1 Functions of One Variable ..... 703
23.2 Orthogonal Collocation ..... 712
23.3 Functions of Multiple Variables ..... 712
24 Differential Equations ..... 719
24.1 Ordinary Differential Equations ..... 719
24.2 Differential-Algebraic Equations ..... 722
24.3 Matlab-compatible solvers ..... 730
25 Optimization ..... 741
25.1 Linear Programming ..... 741
25.2 Quadratic Programming. ..... 747
25.3 Nonlinear Programming ..... 749
25.4 Linear Least Squares ..... 751
26 Statistics ..... 757
26.1 Descriptive Statistics ..... 757
26.2 Statistics on Sliding Windows of Data ..... 766
26.3 Basic Statistical Functions ..... 783
26.4 Correlation and Regression Analysis ..... 786
26.5 Distributions ..... 788
26.6 Random Number Generation ..... 789
27 Sets ..... 791
27.1 Set Operations ..... 793
28 Polynomial Manipulations ..... 797
28.1 Evaluating Polynomials ..... 797
28.2 Finding Roots ..... 798
28.3 Products of Polynomials ..... 799
28.4 Derivatives / Integrals / Transforms ..... 802
28.5 Polynomial Interpolation ..... 803
28.6 Miscellaneous Functions ..... 813
29 Interpolation ..... 815
29.1 One-dimensional Interpolation ..... 815
29.2 Multi-dimensional Interpolation ..... 819
30 Geometry ..... 825
30.1 Delaunay Triangulation ..... 825
30.1.1 Plotting the Triangulation ..... 827
30.1.2 Identifying Points in Triangulation ..... 830
30.2 Voronoi Diagrams ..... 832
30.3 Convex Hull ..... 836
30.4 Interpolation on Scattered Data ..... 838
30.5 Vector Rotation Matrices ..... 841
31 Signal Processing ..... 845
32 Image Processing ..... 859
32.1 Loading and Saving Images ..... 859
32.2 Displaying Images ..... 865
32.3 Representing Images ..... 867
32.4 Plotting on top of Images ..... 878
32.5 Color Conversion ..... 879
33 Audio Processing ..... 881
33.1 Audio File Utilities. ..... 881
33.2 Audio Device Information ..... 882
33.3 Audio Player ..... 883
33.3.1 Playback ..... 883
33.3.2 Properties ..... 884
33.4 Audio Recorder ..... 884
33.4.1 Recording ..... 885
33.4.2 Data Retrieval ..... 885
33.4.3 Properties ..... 886
33.5 Audio Data Processing ..... 886
34 Object Oriented Programming ..... 889
34.1 Creating a Class ..... 889
34.2 Class Methods ..... 891
34.3 Indexing Objects ..... 894
34.3.1 Defining Indexing And Indexed Assignment ..... 894
34.3.2 Indexed Assignment Optimization ..... 898
34.4 Overloading Objects ..... 899
34.4.1 Function Overloading ..... 899
34.4.2 Operator Overloading ..... 900
34.4.3 Precedence of Objects ..... 901
34.5 Inheritance and Aggregation ..... 902
34.6 classdef Classes ..... 906
34.6.1 Creating a classdef Class ..... 906
34.6.2 Properties ..... 908
34.6.3 Methods ..... 909
34.6.4 Inheritance ..... 911
34.6.5 Value Classes vs. Handle Classes ..... 911
35 GUI Development ..... 915
35.1 I/O Dialogs ..... 915
35.2 Progress Bar ..... 923
35.3 UI Elements ..... 924
35.4 GUI Utility Functions ..... 933
35.5 User-Defined Preferences ..... 937
35.6 Octave Workspace Windows ..... 939
36 System Utilities ..... 941
36.1 Timing Utilities ..... 941
36.2 Filesystem Utilities ..... 953
36.3 File Archiving Utilities ..... 963
36.4 Networking Utilities ..... 966
36.4.1 FTP Objects ..... 966
36.4.2 WWW Access ..... 968
36.4.3 Base64 and Binary Data Transmission ..... 971
36.5 Controlling Subprocesses ..... 972
36.6 Process, Group, and User IDs ..... 980
36.7 Environment Variables ..... 981
36.8 Current Working Directory ..... 982
36.9 Password Database Functions ..... 983
36.10 Group Database Functions ..... 984
36.11 System Information ..... 985
36.12 Hashing Functions ..... 994
37 Packages ..... 997
37.1 Installing and Removing Packages ..... 997
37.2 Using Packages ..... 1002
37.3 Administrating Packages ..... 1003
37.4 Creating Packages ..... 1003
37.4.1 The DESCRIPTION File ..... 1005
37.4.2 The INDEX File ..... 1007
37.4.3 PKG_ADD and PKG_DEL Directives ..... 1008
37.4.4 Missing Components ..... 1008
Appendix A External Code Interface ..... 1009
A. 1 Oct-Files ..... 1010
A.1.1 Getting Started with Oct-Files ..... 1010
A.1.2 Matrices and Arrays in Oct-Files ..... 1014
A.1.3 Character Strings in Oct-Files ..... 1017
A.1.4 Cell Arrays in Oct-Files ..... 1019
A.1.5 Structures in Oct-Files ..... 1019
A.1.6 Sparse Matrices in Oct-Files ..... 1021
A.1.6.1 Array and Sparse Class Differences ..... 1021
A.1.6.2 Creating Sparse Matrices in Oct-Files ..... 1022
A.1.6.3 Using Sparse Matrices in Oct-Files ..... 1025
A.1.7 Accessing Global Variables in Oct-Files ..... 1026
A.1.8 Calling Octave Functions from Oct-Files ..... 1027
A.1.9 Calling External Code from Oct-Files ..... 1028
A.1.10 Allocating Local Memory in Oct-Files ..... 1030
A.1.11 Input Parameter Checking in Oct-Files ..... 1030
A.1.12 Exception and Error Handling in Oct-Files ..... 1031
A.1.13 Documentation and Testing of Oct-Files ..... 1033
A. 2 Mex-Files ..... 1034
A.2.1 Getting Started with Mex-Files ..... 1034
A.2.2 Working with Matrices and Arrays in Mex-Files ..... 1036
A.2.3 Character Strings in Mex-Files ..... 1038
A.2.4 Cell Arrays with Mex-Files ..... 1039
A.2.5 Structures with Mex-Files ..... 1040
A.2.6 Sparse Matrices with Mex-Files ..... 1042
A.2.7 Calling Other Functions in Mex-Files ..... 1045
A. 3 Standalone Programs ..... 1046
A. 4 Java Interface ..... 1049
A.4.1 Making Java Classes Available ..... 1050
A.4.2 How to use Java from within Octave ..... 1051
A.4.3 Set up the JVM ..... 1053
A.4.4 Java Interface Functions ..... 1054
Appendix B Test and Demo Functions ..... 1061
B. 1 Test Functions ..... 1061
B. 2 Demonstration Functions ..... 1069
Appendix C Obsolete Functions ..... 1073
Appendix D Known Causes of Trouble ..... 1079
D. 1 Actual Bugs We Haven't Fixed Yet ..... 1079
D. 2 Reporting Bugs... ..... 1079
D.2.1 Have You Found a Bug? ..... 1079
D.2.2 Where to Report Bugs ..... 1080
D.2.3 How to Report Bugs ..... 1080
D.2.4 Sending Patches for Octave ..... 1081
D. 3 How To Get Help with Octave ..... 1082
D. 4 How to Distinguish Between Octave and Matlab ..... 1082
Appendix E Installing Octave ..... 1085
E. 1 Build Dependencies ..... 1085
E.1.1 Obtaining the Dependencies Automatically ..... 1085
E.1.2 Build Tools ..... 1085
E.1.3 External Packages ..... 1086
E. 2 Running Configure and Make ..... 1089
E. 3 Compiling Octave with 64 -bit Indexing ..... 1093
E. 4 Installation Problems ..... 1096
Appendix F Grammar and Parser ..... 1099
F. 1 Keywords ..... 1099
F. 2 Parser ..... 1099
Appendix G GNU GENERAL PUBLIC LICENSE ..... 1101
Concept Index ..... 1113
Function Index ..... 1119
Operator Index ..... 1133
Graphics Properties Index ..... 1135

## Preface

Octave was originally intended to be companion software for an undergraduate-level textbook on chemical reactor design being written by James B. Rawlings of the University of Wisconsin-Madison and John G. Ekerdt of the University of Texas.

Clearly, Octave is now much more than just another 'courseware' package with limited utility beyond the classroom. Although our initial goals were somewhat vague, we knew that we wanted to create something that would enable students to solve realistic problems, and that they could use for many things other than chemical reactor design problems. We find that most students pick up the basics of Octave quickly, and are using it confidently in just a few hours.

Although it was originally intended to be used to teach reactor design, it has been used in several other undergraduate and graduate courses in the Chemical Engineering Department at the University of Texas, and the math department at the University of Texas has been using it for teaching differential equations and linear algebra as well. More recently, Octave has been used as the primary computational tool for teaching Stanford's online Machine Learning class (http://ml-class.org) taught by Andrew Ng. Tens of thousands of students participated in the course.

If you find Octave useful, please let us know. We are always interested to find out how Octave is being used.

Virtually everyone thinks that the name Octave has something to do with music, but it is actually the name of one of John W. Eaton's former professors who wrote a famous textbook on chemical reaction engineering, and who was also well known for his ability to do quick 'back of the envelope' calculations. We hope that this software will make it possible for many people to do more ambitious computations just as easily.

Everyone is encouraged to share this software with others under the terms of the GNU General Public License (see Appendix G [Copying], page 1101). You are also encouraged to help make Octave more useful by writing and contributing additional functions for it, and by reporting any problems you may have.

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## Citing Octave in Publications

In view of the many contributions made by numerous developers over many years it is common courtesy to cite Octave in publications when it has been used during the course of research or the preparation of figures. The citation function can automatically generate a recommended citation text for Octave or any of its packages. See the help text below on how to use citation.

```
citation
```

citation package
Display instructions for citing GNU Octave or its packages in publications.

When called without an argument, display information on how to cite the core GNU Octave system.
When given a package name package, display information on citing the specific named package. Note that some packages may not yet have instructions on how to cite them.
The GNU Octave developers and its active community of package authors have invested a lot of time and effort in creating GNU Octave as it is today. Please give credit where credit is due and cite GNU Octave and its packages when you use them.

## How You Can Contribute to Octave

There are a number of ways that you can contribute to help make Octave a better system. Perhaps the most important way to contribute is to write high-quality code for solving new problems, and to make your code freely available for others to use. See https://www. octave.org/get-involved.html for detailed information.

If you find Octave useful, consider providing additional funding to continue its development. Even a modest amount of additional funding could make a significant difference in the amount of time that is available for development and support.

Donations supporting Octave development may be made on the web at https://my. fsf.org/donate/working-together/octave. These donations also help to support the Free Software Foundation

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If you pay by check, please be sure to write "GNU Octave" in the memo field of your check.
If you cannot provide funding or contribute code, you can still help make Octave better and more reliable by reporting any bugs you find and by offering suggestions for ways to improve Octave. See Appendix D [Trouble], page 1079, for tips on how to write useful bug reports.

## Distribution

Octave is free software. This means that everyone is free to use it and free to redistribute it on certain conditions. Octave is not, however, in the public domain. It is copyrighted and there are restrictions on its distribution, but the restrictions are designed to ensure that others will have the same freedom to use and redistribute Octave that you have. The precise conditions can be found in the GNU General Public License that comes with Octave and that also appears in Appendix G [Copying], page 1101.

To download a copy of Octave, please visit https://www.octave.org/download.html.

## 1 A Brief Introduction to Octave

GNU Octave is a high-level language primarily intended for numerical computations. It is typically used for such problems as solving linear and nonlinear equations, numerical linear algebra, statistical analysis, and for performing other numerical experiments. It may also be used as a batch-oriented language for automated data processing.

The current version of Octave executes in a graphical user interface (GUI). The GUI hosts an Integrated Development Environment (IDE) which includes a code editor with syntax highlighting, built-in debugger, documentation browser, as well as the interpreter for the language itself. A command-line interface for Octave is also available.

GNU Octave is freely redistributable software. You may redistribute it and/or modify it under the terms of the GNU General Public License as published by the Free Software Foundation. The GPL is included in this manual, see Appendix G [Copying], page 1101.

This manual provides comprehensive documentation on how to install, run, use, and extend GNU Octave. Additional chapters describe how to report bugs and help contribute code.

This document corresponds to Octave version 7.3.0.

### 1.1 Running Octave

On most systems, Octave is started with the shell command 'octave'. This starts the graphical user interface. The central window in the GUI is the Octave command-line interface. In this window Octave displays an initial message and then a prompt indicating it is ready to accept input. If you have chosen the traditional command-line interface then only the command prompt appears in the same window that was running a shell. In either case, you can immediately begin typing Octave commands.

If you get into trouble, you can usually interrupt Octave by typing Control-C (written $C-c$ for short). $C-c$ gets its name from the fact that you type it by holding down CTRL and then pressing c. Doing this will normally return you to Octave's prompt.

To exit Octave, type quit or exit at the Octave prompt.
On systems that support job control, you can suspend Octave by sending it a SIGTSTP signal, usually by typing $C-z$.

### 1.2 Simple Examples

The following chapters describe all of Octave's features in detail, but before doing that, it might be helpful to give a sampling of some of its capabilities.

If you are new to Octave, we recommend that you try these examples to begin learning Octave by using it. Lines marked like so, 'octave:13>', are lines you type, ending each with a carriage return. Octave will respond with an answer, or by displaying a graph.

### 1.2.1 Elementary Calculations

Octave can easily be used for basic numerical calculations. Octave knows about arithmetic operations ( $+,-,{ }^{*}, /$ ), exponentiation ( ${ }^{\wedge}$ ), natural logarithms/exponents (log, exp), and the trigonometric functions (sin, cos, ...). Moreover, Octave calculations work on real or imaginary numbers ( $\mathrm{i}, \mathrm{j}$ ). In addition, some mathematical constants such as the base of
the natural logarithm (e) and the ratio of a circle's circumference to its diameter (pi) are pre-defined.
For example, to verify Euler's Identity,

$$
e^{2 \pi}=-1
$$

type the following which will evaluate to -1 within the tolerance of the calculation.

```
octave:1> exp (i*pi)
```


### 1.2.2 Creating a Matrix

Vectors and matrices are the basic building blocks for numerical analysis. To create a new matrix and store it in a variable so that you can refer to it later, type the command
octave:1> $A=[1,1,2 ; 3,5,8 ; 13,21,34]$
Octave will respond by printing the matrix in neatly aligned columns. Octave uses a comma or space to separate entries in a row, and a semicolon or carriage return to separate one row from the next. Ending a command with a semicolon tells Octave not to print the result of the command. For example,
octave:2> $B=$ rand (3, 2);
will create a 3 row, 2 column matrix with each element set to a random value between zero and one.

To display the value of a variable, simply type the name of the variable at the prompt. For example, to display the value stored in the matrix B, type the command

```
octave:3> B
```


### 1.2.3 Matrix Arithmetic

Octave uses standard mathematical notation with the advantage over low-level languages that operators may act on scalars, vector, matrices, or N-dimensional arrays. For example, to multiply the matrix A by a scalar value, type the command

```
octave:4> 2 * A
```

To multiply the two matrices A and B, type the command

```
octave:5> A * B
```

and to form the matrix product $\mathrm{A}^{T} \mathrm{~A}$, type the command
octave:6> A' * A

### 1.2.4 Solving Systems of Linear Equations

Systems of linear equations are ubiquitous in numerical analysis. To solve the set of linear equations $\mathrm{Ax}=\mathrm{b}$, use the left division operator, ' $\backslash$ ':

```
x = A \ b
```

This is conceptually equivalent to $\mathrm{A}^{-1} \mathrm{~b}$, but avoids computing the inverse of a matrix directly.

If the coefficient matrix is singular, Octave will print a warning message and compute a minimum norm solution.

A simple example comes from chemistry and the need to obtain balanced chemical equations. Consider the burning of hydrogen and oxygen to produce water.

$$
\mathrm{H}_{2}+\mathrm{O}_{2} \rightarrow \mathrm{H}_{2} \mathrm{O}
$$

The equation above is not accurate. The Law of Conservation of Mass requires that the number of molecules of each type balance on the left- and right-hand sides of the equation. Writing the variable overall reaction with individual equations for hydrogen and oxygen one finds:

$$
\begin{gathered}
x_{1} \mathrm{H}_{2}+x_{2} \mathrm{O}_{2} \rightarrow \mathrm{H}_{2} \mathrm{O} \\
\mathrm{H}: \quad 2 x_{1}+0 x_{2} \rightarrow 2 \\
\mathrm{O}: \quad 0 x_{1}+2 x_{2} \rightarrow 1
\end{gathered}
$$

The solution in Octave is found in just three steps.

```
octave:1> A = [ 2, 0; 0, 2 ];
octave:2> b = [ 2; 1 ];
octave:3> x = A \ b
```


### 1.2.5 Integrating Differential Equations

Octave has built-in functions for solving nonlinear differential equations of the form

$$
\frac{d x}{d t}=f(x, t), \quad x\left(t=t_{0}\right)=x_{0}
$$

For Octave to integrate equations of this form, you must first provide a definition of the function $f(x, t)$. This is straightforward, and may be accomplished by entering the function body directly on the command line. For example, the following commands define the righthand side function for an interesting pair of nonlinear differential equations. Note that while you are entering a function, Octave responds with a different prompt, to indicate that it is waiting for you to complete your input.

```
octave:1> function xdot = f (x, t)
>
> r = 0.25;
> k = 1.4;
> a = 1.5;
> b = 0.16;
> c = 0.9;
> d = 0.8;
>
> xdot(1) = r*x(1)*(1 - x(1)/k) - a*x(1)*x(2)/(1 + b*x(1));
> xdot(2) = c*a*x(1)*x(2)/(1 + b*x(1)) - d*x(2);
>
> endfunction
```

Given the initial condition
octave:2> x0 = [1; 2];
and the set of output times as a column vector (note that the first output time corresponds to the initial condition given above)
octave:3> t = linspace ( $0,50,200$ )';
it is easy to integrate the set of differential equations:
octave:4> $x=1$ sode ("f", x0, t);
The function 1sode uses the Livermore Solver for Ordinary Differential Equations, described in A. C. Hindmarsh, ODEPACK, a Systematized Collection of ODE Solvers, in: Scientific Computing, R. S. Stepleman et al. (Eds.), North-Holland, Amsterdam, 1983, pages 55-64.

### 1.2.6 Producing Graphical Output

To display the solution of the previous example graphically, use the command
octave:1> plot ( $t, x$ )
Octave will automatically create a separate window to display the plot.
To save a plot once it has been displayed on the screen, use the print command. For example,
print foo.pdf
will create a file called foo.pdf that contains a rendering of the current plot in Portable Document Format. The command
help print
explains more options for the print command and provides a list of additional output file formats.

### 1.2.7 Help and Documentation

Octave has an extensive help facility. The same documentation that is available in printed form is also available from the Octave prompt, because both forms of the documentation are created from the same input file.

In order to get good help you first need to know the name of the command that you want to use. The name of this function may not always be obvious, but a good place to start is to type help --list. This will show you all the operators, keywords, built-in functions, and loadable functions available in the current session of Octave. An alternative is to search the documentation using the lookfor function (described in Section 2.3 [Getting Help], page 21).

Once you know the name of the function you wish to use, you can get more help on the function by simply including the name as an argument to help. For example,
help plot
will display the help text for the plot function.
The part of Octave's help facility that allows you to read the complete text of the printed manual from within Octave normally uses a separate program called Info. When you invoke Info you will be put into a menu driven program that contains the entire Octave manual. Help for using Info is provided in this manual, see Section 2.3 [Getting Help], page 21.

### 1.2.8 Editing What You Have Typed

At the Octave prompt, you can recall, edit, and reissue previous commands using Emacsor vi-style editing commands. The default keybindings use Emacs-style commands. For example, to recall the previous command, press Control-p (written C-p for short). Doing this will normally bring back the previous line of input. $C-n$ will bring up the next line of
input, $C-b$ will move the cursor backward on the line, $C-f$ will move the cursor forward on the line, etc.

A complete description of the command line editing capability is given in this manual, see Section 2.4 [Command Line Editing], page 25.

### 1.3 Conventions

This section explains the notation conventions that are used in this manual. You may want to skip this section and refer back to it later.

### 1.3.1 Fonts

Examples of Octave code appear in this font or form: svd (a). Names that represent variables or function arguments appear in this font or form: first-number. Commands that you type at the shell prompt appear in this font or form: 'octave --no-init-file'. Commands that you type at the Octave prompt sometimes appear in this font or form: foo --bar --baz. Specific keys on your keyboard appear in this font or form: RET.

### 1.3.2 Evaluation Notation

In the examples in this manual, results from expressions that you evaluate are indicated with ' $\Rightarrow$ '. For example:
sqrt (2)

$$
\Rightarrow 1.4142
$$

You can read this as "sqrt (2) evaluates to 1.4142 ".
In some cases, matrix values that are returned by expressions are displayed like this

$$
\begin{aligned}
& {[1,2 ; 3,4]==[1,3 ; 2,4]} \\
& \quad \Rightarrow[1,0 ; 0,1]
\end{aligned}
$$

and in other cases, they are displayed like this

```
eye (3)
    # 1 0 0
    0}1
    0
```

in order to clearly show the structure of the result.
Sometimes to help describe one expression, another expression is shown that produces identical results. The exact equivalence of expressions is indicated with ' $\equiv$ '. For example:

```
rot90 ([1, 2; 3, 4], -1)
\equiv
rot90 ([1, 2; 3, 4], 3)
\equiv
rot90 ([1, 2; 3, 4], 7)
```


### 1.3.3 Printing Notation

Many of the examples in this manual print text when they are evaluated. In this manual the printed text resulting from an example is indicated by ' $\dashv$ '. The value that is returned
by evaluating the expression is displayed with ' $\Rightarrow$ ' ( 1 in the next example) and follows on a separate line.

```
printf ("foo %s\n", "bar")
    foo bar
    => 1
```


### 1.3.4 Error Messages

Some examples signal errors. This normally displays an error message on your terminal. Error messages are shown on a line beginning with error:.

```
fieldnames ([1, 2; 3, 4])
error: fieldnames: Invalid input argument
```


### 1.3.5 Format of Descriptions

Functions and commands are described in this manual in a uniform format. The first line of a description contains the name of the item followed by its arguments, if any. If there are multiple ways to invoke the function then each allowable form is listed.

The description follows on succeeding lines, sometimes with examples.

### 1.3.5.1 A Sample Function Description

In a function description, the name of the function being described appears first. It is followed on the same line by a list of parameters. The names used for the parameters are also used in the body of the description.

After all of the calling forms have been enumerated, the next line is a concise one-sentence summary of the function.

After the summary there may be documentation on the inputs and outputs, examples of function usage, notes about the algorithm used, and references to related functions.

Here is a description of an imaginary function foo:

```
foo (x)
foo (x, y)
foo (x, y, ...)
    Subtract }x\mathrm{ from y, then add any remaining arguments to the result.
```

The input $x$ must be a numeric scalar, vector, or array.

The optional input $y$ defaults to 19 if it is not supplied.

Example:

```
foo (1, [3, 5], 3, 9)
            =>[ 14, 16 ]
foo (5)
            => 14
```

More generally,

```
foo (w, x, y, ...)
\equiv
x - w + y + ...
```

See also: bar

Any parameter whose name contains the name of a type (e.g., integer or matrix) is expected to be of that type. Parameters named object may be of any type. Parameters with other sorts of names (e.g., new_file) are discussed specifically in the description of the function. In some sections, features common to parameters of several functions are described at the beginning.

### 1.3.5.2 A Sample Command Description

Commands are functions that may be called without surrounding their arguments in parentheses. Command descriptions have a format similar to function descriptions. For example, here is the description for Octave's diary command:

```
diary
diary on
diary off
diary filename
[status, diaryfile] = diary
```

Record a list of all commands and the output they produce, mixed together just as they appear on the terminal.
Valid options are:
on Start recording a session in a file called diary in the current working directory.
off Stop recording the session in the diary file.
filename Record the session in the file named filename.
With no input or output arguments, diary toggles the current diary state.
If output arguments are requested, diary ignores inputs and returns the current status. The boolean status indicates whether recording is on or off, and diaryfile is the name of the file where the session is stored.
See also: history, evalc.

## 2 Getting Started

This chapter explains some of Octave's basic features, including how to start an Octave session, get help at the command prompt, edit the command line, and write Octave programs that can be executed as commands from your shell.

### 2.1 Invoking Octave from the Command Line

Normally, Octave is used interactively by running the program 'octave' without any arguments. Once started, Octave reads commands from the terminal until you tell it to exit.

You can also specify the name of a file on the command line, and Octave will read and execute the commands from the named file and then exit when it is finished.

You can further control how Octave starts by using the command-line options described in the next section, and Octave itself can remind you of the options available. Type 'octave --help' to display all available options and briefly describe their use ('octave -h' is a shorter equivalent).

### 2.1.1 Command Line Options

Here is a complete list of the command line options that Octave accepts.
--built-in-docstrings-file filename
Specify the name of the file containing documentation strings for the built-in functions of Octave. This value is normally correct and should only need to specified in extraordinary situations.
--debug
-d Enter parser debugging mode. Using this option will cause Octave's parser to print a lot of information about the commands it reads, and is probably only useful if you are actually trying to debug the parser.
--doc-cache-file filename
Specify the name of the doc cache file to use. The value of filename specified on the command line will override any value of OCTAVE_DOC_CACHE_FILE found in the environment, but not any commands in the system or user startup files that use the doc_cache_file function.
--echo-commands
$-\mathrm{x} \quad$ Echo commands as they are executed.
--eval code
Evaluate code and exit when finished unless --persist is also specified.
--exec-path path
Specify the path to search for programs to run. The value of path specified on the command line will override any value of OCTAVE_EXEC_PATH found in the environment, but not any commands in the system or user startup files that call the EXEC_PATH function.
--gui Start the graphical user interface (GUI).

```
--help
-h Print short help message and exit.
--image-path path
```

Add path to the head of the search path for images. The value of path specified on the command line will override any value of OCTAVE_IMAGE_PATH found in the environment, but not any commands in the system or user startup files that call the IMAGE_PATH function.

## --info-file filename

Specify the name of the info file to use. The value of filename specified on the command line will override any value of OCTAVE_INFO_FILE found in the environment, but not any commands in the system or user startup files that use the info_file function.

```
--info-program program
```

Specify the name of the info program to use. The value of program specified on the command line will override any value of OCTAVE_INFO_PROGRAM found in the environment, but not any commands in the system or user startup files that use the info_program function.

## --interactive

-i Force interactive behavior. This can be useful for running Octave via a remote shell command or inside an Emacs shell buffer.

## --line-editing

Force readline use for command-line editing.
--no-gui Disable the graphical user interface (GUI) and use the command line interface (CLI) instead. This is the default behavior, but this option may be useful to override a previous --gui.

```
--no-history
```

-H Disable recording of command-line history.
--no-init-file
Don't read the initialization files ~/.octaverc and .octaverc.
--no-init-path
Don't initialize the search path for function files to include default locations.
--no-line-editing
Disable command-line editing.

```
--no-site-file
```

Don't read the site-wide octaverc initialization files.
--no-window-system
-W Disable use of a windowing system including graphics. This forces a strictly terminal-only environment.
--norc
-f Don't read any of the system or user initialization files at startup. This is equivalent to using both of the options --no-init-file and --no-site-file.
--path path
-p path Add path to the head of the search path for function files. The value of path specified on the command line will override any value of OCTAVE_PATH found in the environment, but not any commands in the system or user startup files that set the internal load path through one of the path functions.
--persist
Go to interactive mode after --eval or reading from a file named on the command line.
--silent
--quiet
-q Don't print the usual greeting and version message at startup.
--texi-macros-file filename
Specify the name of the file containing Texinfo macros for use by makeinfo.
--traditional
--braindead
For compatibility with matlab, set initial values for user preferences to the following values

```
PS1 = ">> "
PS2 = ""
beep_on_error = true
confirm_recursive_rmdir = false
crash_dumps_octave_core = false
optimize_diagonal_matrix = false
optimize_permutation_matrix = false
optimize_range = false
fixed_point_format = true
history_timestamp_format_string = "%%-- %D %I:%M %p --%%"
print_empty_dimensions = false
print_struct_array_contents = true
save_default_options = "-mat-binary"
struct_levels_to_print = 0
```

and disable the following warnings

```
Octave:abbreviated-property-match
Octave:colon-nonscalar-argument
Octave:data-file-in-path
Octave:empty-index
Octave:function-name-clash
Octave:possible-matlab-short-circuit-operator
```

Note that this does not enable the Octave:language-extension warning, which you might want if you want to be told about writing code that works in Octave but not matlab (see [warning], page 246, [warning_ids], page 248).
--verbose
-V Turn on verbose output.
--version
-v Print the program version number and exit.
file Execute commands from file. Exit when done unless --persist is also specified.
Octave also includes several functions which return information about the command line, including the number of arguments and all of the options.
argv ()
Return the command line arguments passed to Octave.
For example, if you invoked Octave using the command

```
            octave --no-line-editing --silent
```

        argv would return a cell array of strings with the elements --no-line-editing and
        --silent.
    If you write an executable Octave script, argv will return the list of arguments passed to the script. See Section 2.6 [Executable Octave Programs], page 37, for an example of how to create an executable Octave script.
program_name ()
Return the last component of the value returned by program_invocation_name.
See also: [program_invocation_name], page 18.
program_invocation_name ()
Return the name that was typed at the shell prompt to run Octave.
If executing a script from the command line (e.g., octave foo.m) or using an executable Octave script, the program name is set to the name of the script. See Section 2.6 [Executable Octave Programs], page 37, for an example of how to create an executable Octave script.
See also: [program_name], page 18.
Here is an example of using these functions to reproduce the command line which invoked Octave.

```
printf ("%s", program_name ());
```

arg_list = argv () ;
for $i=1$ :nargin
printf (" \%s", arg_list\{i\});
endfor
printf ("\n");

See Section 6.3.3 [Indexing Cell Arrays], page 130, for an explanation of how to retrieve objects from cell arrays, and Section 11.2 [Defining Functions], page 191, for information about the variable nargin.

### 2.1.2 Startup Files

When Octave starts, it looks for commands to execute from the files in the following list. These files may contain any valid Octave commands, including function definitions.
octave-home/share/octave/site/m/startup/octaverc
where octave-home is the directory in which Octave is installed (the default is /usr/local). This file is provided so that changes to the default Octave
environment can be made globally for all users at your site for all versions of Octave you have installed. Care should be taken when making changes to this file since all users of Octave at your site will be affected. The default file may be overridden by the environment variable OCTAVE_SITE_INITFILE.
octave-home/share/octave/version/m/startup/octaverc
where octave-home is the directory in which Octave is installed (the default is /usr/local), and version is the version number of Octave. This file is provided so that changes to the default Octave environment can be made globally for all users of a particular version of Octave. Care should be taken when making changes to this file since all users of Octave at your site will be affected. The default file may be overridden by the environment variable OCTAVE_VERSION_INITFILE.
config-dir/octave/octaverc
where config-dir is the platform-dependent location for user local configuration files (e.g., \$XDG_CONFIG_HOME on many Unix-like operating systems or \%APPDATA\% on Windows).
~/.octaverc
This file is used to make personal changes to the default Octave environment.
.octaverc
This file can be used to make changes to the default Octave environment for a particular project. Octave searches for this file in the current directory after it reads ~/.octaverc. Any use of the cd command in the ~/.octaverc file will affect the directory where Octave searches for .octaverc.
If you start Octave in your home directory, commands from the file ~/.octaverc will only be executed once.
startup.m
This file is used to make personal changes to the default Octave environment. It is executed for mATLAB compatibility, but $\sim /$. octaverc is the preferred location for configuration changes.
A message will be displayed as each of the startup files is read if you invoke Octave with the --verbose option but without the --silent option.

The startup files are always processed in the system's locale charset (independent of the m-file encoding that is set, for example, in the GUI properties). In other words, the system's locale charset is in effect until a user manually sets the m-file encoding (e.g., in one of the startup files) and triggers re-parsing of any relevant m-files. Octave can be forced to use a new encoding with the undocumented function __mfile_encoding__:

```
__mfile_encoding__ ("utf-8"); # set new encoding
clear ("functions"); # re-parse all .m files in the new encoding
```

This changes the encoding that is used to interpret all subsequently run startup and m -files (not including the currently executing file).

### 2.2 Quitting Octave

Shutdown is initiated with the exit or quit commands (they are equivalent). Similar to startup, Octave has a shutdown process that can be customized by user script files.

During shutdown Octave will search for the script file finish.m in the function load path. Commands to save all workspace variables or cleanup temporary files may be placed there. Additional functions to execute on shutdown may be registered with atexit.

Quit the current Octave session.
The exit function is an alias for quit.
If the optional integer value status is supplied, pass that value to the operating system as Octave's exit status. The default value is zero.
When exiting, Octave will attempt to run the m-file finish.m if it exists. User commands to save the workspace or clean up temporary files may be placed in that file. Alternatively, another m-file may be scheduled to run using atexit. If an error occurs while executing the finish.m file, Octave does not exit and control is returned to the command prompt.
If the optional argument "cancel" is provided, Octave does not exit and control is returned to the command prompt. This feature allows the finish.m file to cancel the quit process.
If the user preference to request confirmation before exiting, Octave will display a dialog and give the user an option to cancel the exit process.
If the optional argument "force" is provided, no confirmation is requested, and the execution of the finish.m file is skipped.
See also: [atexit], page 20.
atexit (fcn)
atexit (fcn, flag)
Register a function to be called when Octave exits.
For example,

```
function last_words ()
    disp ("Bye bye");
endfunction
atexit ("last_words");
```

will print the message "Bye bye" when Octave exits.
The additional argument flag will register or unregister fcn from the list of functions to be called when Octave exits. If flag is true, the function is registered, and if flag is false, it is unregistered. For example, after registering the function last_words above,

```
atexit ("last_words", false);
```

will remove the function from the list and Octave will not call last_words when it exits.

Note that atexit only removes the first occurrence of a function from the list, so if a function was placed in the list multiple times with atexit, it must also be removed from the list multiple times.

See also: [quit], page 20.

### 2.3 Commands for Getting Help

The entire text of this manual is available from the Octave prompt via the command doc. In addition, the documentation for individual user-written functions and variables is also available via the help command. This section describes the commands used for reading the manual and the documentation strings for user-supplied functions and variables. See Section 11.10 [Function Files], page 210, for more information about how to document the functions you write.

```
help name
help -list
help.
help
```

Display the help text for name.
For example, the command help help prints a short message describing the help command.
Given the single argument --list, list all operators, keywords, built-in functions, and loadable functions available in the current session of Octave.
Given the single argument ., list all operators available in the current session of Octave.
If invoked without any arguments, help displays instructions on how to access help from the command line.
The help command can provide information about most operators, but name must be enclosed by single or double quotes to prevent the Octave interpreter from acting on name. For example, help "+" displays help on the addition operator.
See also: [doc], page 21, [lookfor], page 21, [which], page 148, [info], page 22.
doc function_name
doc
Display documentation for the function function_name directly from an online version of the printed manual, using the GNU Info browser.
If invoked without an argument, the manual is shown from the beginning.
For example, the command doc rand starts the GNU Info browser at the rand node in the online version of the manual.
Once the GNU Info browser is running, help for using it is available using the command $C$-h.

See also: [help], page 21.
lookfor str
lookfor -all str
[fcn, help1str] = lookfor (str)
[fcn, help1str] = lookfor ("-all", str)
Search for the string str in the documentation of all functions in the current function search path.
By default, lookfor looks for str in just the first sentence of the help string for each function found. The entire help text of each function can be searched by using the "-all" argument. All searches are case insensitive.
When called with no output arguments, lookfor prints the list of matching functions to the terminal. Otherwise, the output argument fcns contains the function names and help1str contains the first sentence from the help string of each function.
Programming Note: The ability of lookfor to correctly identify the first sentence of the help text is dependent on the format of the function's help. All Octave core functions are correctly formatted, but the same can not be guaranteed for external packages and user-supplied functions. Therefore, the use of the "-all" argument may be necessary to find related functions that are not a part of Octave.
The speed of lookup is greatly enhanced by having a cached documentation file. For more information, see [doc_cache_create], page 24 .
See also: [help], page 21, [doc], page 21, [which], page 148, [path], page 214, [doc_cache_create], page 24.

To see what is new in the current release of Octave, use the news function.
news
news package
Display the current NEWS file for Octave or an installed package.
When called without an argument, display the NEWS file for Octave.
When given a package name package, display the current NEWS file for that package.
See also: [ver], page 988, [pkg], page 997.
info ()
Display contact information for the GNU Octave community.
warranty ()
Describe the conditions for copying and distributing Octave.
The following functions can be used to change which programs are used for displaying the documentation, and where the documentation can be found.

```
val = info_file ()
old_val = info_file (new_val)
info_file (new_val, "local")
```

Query or set the internal variable that specifies the name of the Octave info file.
The default value is octave-home/share/info/octave.info, in which octave-home is the root directory of the Octave installation. The default value may be overridden by the environment variable OCTAVE_INFO_FILE, or the command line argument --info-file FNAME.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [info_program], page 23, [doc], page 21, [help], page 21, [makeinfo_program], page 23 .
val = info_program ()
old_val = info_program (new_val)
info_program (new_val, "local")
Query or set the internal variable that specifies the name of the info program to run.
The default value is info. The default value may be overridden by the environment variable OCTAVE_INFO_PROGRAM, or the command line argument --info-program NAME.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [info_file], page 22, [doc], page 21, [help], page 21, [makeinfo_program], page 23 .
val = makeinfo_program ()
old_val = makeinfo_program (new_val)
makeinfo_program (new_val, "local")
Query or set the internal variable that specifies the name of the program that Octave runs to format help text containing Texinfo markup commands.
The default value is makeinfo.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [texi_macros_file], page 23, [info_file], page 22, [info_program], page 23, [doc], page 21 , [help], page 21 .

```
val = texi_macros_file ()
```

old_val = texi_macros_file (new_val)
texi_macros_file (new_val, "local")

Query or set the internal variable that specifies the name of the file containing Texinfo macros that are prepended to documentation strings before they are passed to makeinfo.
The default value is octave-home/share/octave/version/etc/macros.texi, in which octave-home is the root directory of the Octave installation, and version is the Octave version number. The default value may be overridden by the environment variable OCTAVE_TEXI_MACROS_FILE, or the command line argument --texi-macros-file FNAME.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [makeinfo_program], page 23.

```
val = doc_cache_file ()
```

old_val = doc_cache_file (new_val)
doc_cache_file (new_val, "local")

Query or set the internal variable that specifies the name of the Octave documentation cache file.
A cache file significantly improves the performance of the lookfor command. The default value is octave-home/share/octave/version/etc/doc-cache, in which octave-home is the root directory of the Octave installation, and version is the Octave version number. The default value may be overridden by the environment variable OCTAVE_DOC_CACHE_FILE, or the command line argument --doc-cache-file FNAME.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [doc_cache_create], page 24, [lookfor], page 21, [info_program], page 23, [doc], page 21, [help], page 21, [makeinfo_program], page 23.

See also: [lookfor], page 21.
val = built_in_docstrings_file ()
old_val = built_in_docstrings_file (new_val)
built_in_docstrings_file (new_val, "local")
Query or set the internal variable that specifies the name of the file containing docstrings for built-in Octave functions.
The default value is octave-home/share/octave/version/etc/built-indocstrings, in which octave-home is the root directory of the Octave installation, and version is the Octave version number. The default value may be overridden by the environment variable OCTAVE_BUILT_IN_DOCSTRINGS_FILE, or the command line argument --built-in-docstrings-file FNAME.
Note: This variable is only used when Octave is initializing itself. Modifying it during a running session of Octave will have no effect.

```
val = suppress_verbose_help_message ()
old_val = suppress_verbose_help_message (new_val)
suppress_verbose_help_message (new_val, "local")
```

Query or set the internal variable that controls whether Octave will add additional help information to the end of the output from the help command and usage messages for built-in commands.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

The following functions are principally used internally by Octave for generating the documentation. They are documented here for completeness and because they may occasionally be useful for users.

```
doc_cache_create (out_file, directory)
doc_cache_create (out_file)
```

doc_cache_create ()
Generate documentation cache for all functions in directory.
A documentation cache is generated for all functions in directory which may be a single string or a cell array of strings. The cache is used to speed up the function lookfor.
The cache is saved in the file out_file which defaults to the value doc-cache if not given.
If no directory is given (or it is the empty matrix), a cache for built-in functions, operators, and keywords is generated.
See also: [doc_cache_file], page 24, [lookfor], page 21, [path], page 214.

```
[text, format] = get_help_text (name)
```

Return the raw help text of function name.
The raw help text is returned in text and the format in format. The format is a string which is one of "texinfo", "html", or "plain text".
See also: [get_help_text_from_file], page 25.
[text, format] = get_help_text_from_file (fname)
Return the raw help text from the file fname.
The raw help text is returned in text and the format in format. The format is a string which is one of "texinfo", "html", or "plain text".
See also: [get_help_text], page 25.
text = get_first_help_sentence (name)
text = get_first_help_sentence (name, max_len)
[text, status] = get_first_help_sentence (...)
Return the first sentence of a function's help text.
The first sentence is defined as the text after the function declaration until either the first period (".") or the first appearance of two consecutive newlines (" $\mathrm{n} \backslash \mathrm{n}$ "). The text is truncated to a maximum length of max_len, which defaults to 80 . If the text must be truncated the last three characters of the text are replaced with "..." to indicate that more text was available.
The optional output argument status returns the status reported by makeinfo. If only one output argument is requested, and status is nonzero, a warning is displayed.
As an example, the first sentence of this help text is

```
get_first_help_sentence ("get_first_help_sentence")
\dashvans = Return the first sentence of a function's help text.
```


### 2.4 Command Line Editing

Octave uses the GNU Readline library to provide an extensive set of command-line editing and history features. Only the most common features are described in this manual. In addition, all of the editing functions can be bound to different key strokes at the user's discretion. This manual assumes no changes from the default Emacs bindings. See the

GNU Readline Library manual for more information on customizing Readline and for a complete feature list.

To insert printing characters (letters, digits, symbols, etc.), simply type the character. Octave will insert the character at the cursor and advance the cursor forward.

Many of the command-line editing functions operate using control characters. For example, the character Control-a moves the cursor to the beginning of the line. To type $C-$ a, hold down CTRL and then press a. In the following sections, control characters such as Control-a are written as C -a.

Another set of command-line editing functions use Meta characters. To type $M-u$, hold down the META key and press $u$. Depending on the keyboard, the META key may be labeled ALT or even WINDOWS. If your terminal does not have a META key, you can still type Meta characters using two-character sequences starting with ESC. Thus, to enter M-u, you would type ESC u. The ESC character sequences are also allowed on terminals with real Meta keys. In the following sections, Meta characters such as Meta-u are written as M-u.

### 2.4.1 Cursor Motion

The following commands allow you to position the cursor.
C-b Move back one character.
$C-f \quad$ Move forward one character.
BACKSPACE
Delete the character to the left of the cursor.
DEL Delete the character underneath the cursor.
C-d Delete the character underneath the cursor.
M-f Move forward a word.
M-b Move backward a word.
C-a $\quad$ Move to the start of the line.
C-e $\quad$ Move to the end of the line.
C-1 Clear the screen, reprinting the current line at the top.
$\begin{array}{ll}C- \\ C- & \text { Undo the last action. You can undo all the way back to an empty line. }\end{array}$
$M-r \quad U n d o$ all changes made to this line. This is like typing the 'undo' command enough times to get back to the beginning.

The above table describes the most basic possible keystrokes that you need in order to do editing of the input line. On most terminals, you can also use the left and right arrow keys in place of $C-f$ and $C-b$ to move forward and backward.

Notice how $C-f$ moves forward a character, while $M-f$ moves forward a word. It is a loose convention that control keystrokes operate on characters while meta keystrokes operate on words.

The function clc will allow you to clear the screen from within Octave programs.
clc ()
home ()
Clear the terminal screen and move the cursor to the upper left corner.

### 2.4.2 Killing and Yanking

Killing text means to delete the text from the line, but to save it away for later use, usually by yanking it back into the line. If the description for a command says that it 'kills' text, then you can be sure that you can get the text back in a different (or the same) place later.

Here is the list of commands for killing text.
$C-k \quad$ Kill the text from the current cursor position to the end of the line.
$M-d \quad$ Kill from the cursor to the end of the current word, or if between words, to the end of the next word.

M-DEL Kill from the cursor to the start of the previous word, or if between words, to the start of the previous word.
$C-$ K Kill from the cursor to the previous whitespace. This is different than M-DEL because the word boundaries differ.

And, here is how to yank the text back into the line. Yanking means to copy the most-recently-killed text from the kill buffer.

C-y Yank the most recently killed text back into the buffer at the cursor.
$M-y \quad$ Rotate the kill-ring, and yank the new top. You can only do this if the prior command is $C-y$ or $M-y$.

When you use a kill command, the text is saved in a kill-ring. Any number of consecutive kills save all of the killed text together, so that when you yank it back, you get it in one clean sweep. The kill ring is not line specific; the text that you killed on a previously typed line is available to be yanked back later, when you are typing another line.

### 2.4.3 Commands for Changing Text

The following commands can be used for entering characters that would otherwise have a special meaning (e.g., TAB, $C-q$, etc.), or for quickly correcting typing mistakes.
$C-q$
$C-v \quad$ Add the next character that you type to the line verbatim. This is how to insert things like $C$-q for example.
$M-T A B \quad$ Insert a tab character.
$C-t \quad$ Drag the character before the cursor forward over the character at the cursor, also moving the cursor forward. If the cursor is at the end of the line, then transpose the two characters before it.
$M-t \quad$ Drag the word behind the cursor past the word in front of the cursor moving the cursor over that word as well.

M-u Uppercase the characters following the cursor to the end of the current (or following) word, moving the cursor to the end of the word.

M-1 Lowercase the characters following the cursor to the end of the current (or following) word, moving the cursor to the end of the word.
$M-c \quad$ Uppercase the character following the cursor (or the beginning of the next word if the cursor is between words), moving the cursor to the end of the word.

### 2.4.4 Letting Readline Type for You

The following commands allow Octave to complete command and variable names for you.
TAB Attempt to do completion on the text before the cursor. Octave can complete the names of commands and variables.
$M-$ ? List the possible completions of the text before the cursor.
val = completion_append_char ()
old_val = completion_append_char (new_val)
completion_append_char (new_val, "local")
Query or set the internal character variable that is appended to successful commandline completion attempts.
The default value is " " (a single space).
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

```
completion_matches (hint)
```

Generate possible completions given hint.
This function is provided for the benefit of programs like Emacs which might be controlling Octave and handling user input. The current command number is not incremented when this function is called. This is a feature, not a bug.

### 2.4.5 Commands for Manipulating the History

Octave normally keeps track of the commands you type so that you can recall previous commands to edit or execute them again. When you exit Octave, the most recent commands you have typed, up to the number specified by the variable history_size, are saved in a file. When Octave starts, it loads an initial list of commands from the file named by the variable history_file.

Here are the commands for simple browsing and searching the history list.
LFD
RET Accept the current line regardless of where the cursor is. If the line is nonempty, add it to the history list. If the line was a history line, then restore the history line to its original state.
$c-p \quad$ Move 'up' through the history list.
C-n Move 'down' through the history list.
$M-<\quad$ Move to the first line in the history.
$M \rightarrow \quad$ Move to the end of the input history, i.e., the line you are entering!
$C-r \quad$ Search backward starting at the current line and moving 'up' through the history as necessary. This is an incremental search.

C-s Search forward starting at the current line and moving 'down' through the history as necessary.

On most terminals, you can also use the up and down arrow keys in place of $C-p$ and $C-n$ to move through the history list.

In addition to the keyboard commands for moving through the history list, Octave provides three functions for viewing, editing, and re-running chunks of commands from the history list.

```
history
history opt1...
h = history ()
h = history (opt1,...)
```

If invoked with no arguments, history displays a list of commands that you have executed.

Valid options are:
n
-n Display only the most recent $n$ lines of history.
-c Clear the history list.
-q Don't number the displayed lines of history. This is useful for cutting and pasting commands using the X Window System.
-r file Read the file file, appending its contents to the current history list. If the name is omitted, use the default history file (normally ~/.octave_hist).
${ }^{-w}$ file Write the current history to the file file. If the name is omitted, use the default history file (normally ~/.octave_hist).

For example, to display the five most recent commands that you have typed without displaying line numbers, use the command history -q 5 .
If invoked with a single output argument, the history will be saved to that argument as a cell string and will not be output to screen.
See also: [edit_history], page 29, [run_history], page 30.

```
edit_history
edit_history cmd_number
edit_history first last
```

Edit the history list using the editor named by the variable EDITOR.
The commands to be edited are first copied to a temporary file. When you exit the editor, Octave executes the commands that remain in the file. It is often more convenient to use edit_history to define functions rather than attempting to enter them directly on the command line. The block of commands is executed as soon as you exit the editor. To avoid executing any commands, simply delete all the lines from the buffer before leaving the editor.

When invoked with no arguments, edit the previously executed command; With one argument, edit the specified command cmd_number; With two arguments, edit the list of commands between first and last. Command number specifiers may also be negative where -1 refers to the most recently executed command. The following are equivalent and edit the most recently executed command.

```
edit_history
edit_history -1
```

When using ranges, specifying a larger number for the first command than the last command reverses the list of commands before they are placed in the buffer to be edited.

See also: [run_history], page 30, [history], page 29.
run_history
run_history cmd_number
run_history first last
Run commands from the history list.
When invoked with no arguments, run the previously executed command;
With one argument, run the specified command cmd_number;
With two arguments, run the list of commands between first and last. Command number specifiers may also be negative where -1 refers to the most recently executed command. For example, the command

```
run_history
    OR
run_history -1
```

executes the most recent command again. The command

```
run_history 13 169
```

executes commands 13 through 169.
Specifying a larger number for the first command than the last command reverses the list of commands before executing them. For example:

```
disp (1)
disp (2)
run_history -1 -2
#
    2
    1
```

See also: [edit_history], page 29, [history], page 29.
Octave also allows you customize the details of when, where, and how history is saved.

```
val = history_save ()
old_val = history_save (new_val)
history_save (new_val, "local")
```

Query or set the internal variable that controls whether commands entered on the command line are saved in the history file.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [history_control], page 31, [history_file], page 31, [history_size], page 32, [history_timestamp_format_string], page 32.
val = history_control ()
old_val = history_control (new_val)
Query or set the internal variable that specifies how commands are saved to the history list.
The default value is an empty character string, but may be overridden by the environment variable OCTAVE_HISTCONTROL.
The value of history_control is a colon-separated list of values controlling how commands are saved on the history list. If the list of values includes ignorespace, lines which begin with a space character are not saved in the history list. A value of ignoredups causes lines matching the previous history entry to not be saved. A value of ignoreboth is shorthand for ignorespace and ignoredups. A value of erasedups causes all previous lines matching the current line to be removed from the history list before that line is saved. Any value not in the above list is ignored. If history_ control is the empty string, all commands are saved on the history list, subject to the value of history_save.
See also: [history_file], page 31, [history_size], page 32, [history_timestamp_format_string],
page 32 , [history_save], page 30 .
val = history_file ()
old_val = history_file (new_val)
Query or set the internal variable that specifies the name of the file used to store command history.
All future commands issued during the current Octave session will be written to this new file (if the current setting of history_save allows for this).
The default value is \$DATA/octave/history, where \$DATA is the platform-specific location for (roaming) user data files (e.g., \$XDG_DATA_HOME or, if that is not set, ~/.local/share on Unix-like operating systems or \%APPDATA\% on Windows). The default value may be overridden by the environment variable OCTAVE_HISTFILE.
Programming Notes:
If you want to permanently change the location of Octave's history file you need to issue the history_file command in every new Octave session. This can be achieved by using Octave's .octaverc startup file.
If you also want to read the saved history commands of past Octave sessions from this different history file, then you need to use the additional command history -r after setting the new value of the history file. Example code in Octave's startup file to do this might look like this:

```
history_file ("~/new/.octave_hist");
if (exist (history_file ()))
    history ("-r", history_file());
endif
```

See also: [history], page 29, [history_control], page 31, [history_save], page 30, [history_size], page 32, [history_timestamp_format_string], page 32.
val = history_size ()
old_val = history_size (new_val)
Query or set the internal variable that specifies how many entries to store in the history file.
The default value is 1000 , but may be overridden by the environment variable OCTAVE_HISTSIZE.

See also: [history_file], page 31, [history_timestamp_format_string], page 32, [history_save], page 30.
val = history_timestamp_format_string ()
old_val = history_timestamp_format_string (new_val)
history_timestamp_format_string (new_val, "local")
Query or set the internal variable that specifies the format string for the comment line that is written to the history file when Octave exits.
The format string is passed to strftime. The default value is
"\# Octave VERSION, \%a \%b \%d \%H: \%M:\%S \%Y \%Z [USER@HOST](mailto:USER@HOST)"
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [strftime], page 943, [history_file], page 31, [history_size], page 32, [history_save], page 30.
val = EDITOR ()
old_val = EDITOR (new_val)
EDITOR (new_val, "local")
Query or set the internal variable that specifies the default text editor.
The default value is taken from the environment variable EDITOR when Octave starts. If the environment variable is not initialized, EDITOR will be set to "emacs".
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [edit], page 211, [edit_history], page 29.

### 2.4.6 Customizing readline

Octave uses the GNU Readline library for command-line editing and history features. Readline is very flexible and can be modified through a configuration file of commands (See the GNU Readline library for the exact command syntax). The default configuration file is normally ~/.inputrc.

Octave provides two commands for initializing Readline and thereby changing the command line behavior.
readline_read_init_file (file)
Read the readline library initialization file file.
If file is omitted, read the default initialization file (normally $\sim /$. inputrc).
See Section "Readline Init File" in GNU Readline Library, for details.
See also: [readline_re_read_init_file], page 33.
readline_re_read_init_file ()
Re-read the last readline library initialization file that was read.
See Section "Readline Init File" in GNU Readline Library, for details.
See also: [readline_read_init_file], page 32.

### 2.4.7 Customizing the Prompt

The following variables are available for customizing the appearance of the command-line prompts. Octave allows the prompt to be customized by inserting a number of backslashescaped special characters that are decoded as follows:
' t ’ The time.
'\d' The date.
' n ' Begins a new line by printing the equivalent of a carriage return followed by a line feed.
'\s' The name of the program (usually just 'octave').
‘\w’ The current working directory.
' W' The basename of the current working directory.
'\u' The username of the current user.
' h ' The hostname, up to the first ' $\quad$ '.
‘ H ’ The hostname.
' $\ \#$ ' The command number of this command, counting from when Octave starts.
' $\$ ' The history number of this command. This differs from ' $\backslash$ ' by the number of commands in the history list when Octave starts.
' $\$$ ' If the effective UID is 0 , a ' $\#$ ', otherwise a ' $\$$ '.
' $\backslash n n n$ ' The character whose character code in octal is nnn.
‘<br>' A backslash.
val = PS1 ()
old_val = PS1 (new_val)
PS1 (new_val, "local")
Query or set the primary prompt string.
When executing interactively, Octave displays the primary prompt when it is ready to read a command.

The default value of the primary prompt string is 'octave:<br>\#> '. To change it, use a command like

PS1 ('\u@\H> ')
which will result in the prompt 'boris@kremvax>' for the user 'boris' logged in on the host 'kremvax.kgb.su'. Note that two backslashes are required to enter a backslash into a double-quoted character string. See Chapter 5 [Strings], page 71.
You can also use ANSI escape sequences if your terminal supports them. This can be useful for coloring the prompt. For example,

PS1 ('$$
\033[01;31m
$$\s:<br>\#> $$
\033[0m
$$')

will give the default Octave prompt a red coloring.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [PS2], page 34, [PS4], page 34.
val = PS2 ()
old_val = PS2 (new_val)
PS2 (new_val, "local")
Query or set the secondary prompt string.
The secondary prompt is printed when Octave is expecting additional input to complete a command. For example, if you are typing a for loop that spans several lines, Octave will print the secondary prompt at the beginning of each line after the first. The default value of the secondary prompt string is "> ".
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [PS1], page 33, [PS4], page 34.
val = PS4 ()
old_val = PS4 (new_val)
PS4 (new_val, "local")
Query or set the character string used to prefix output produced when echoing commands is enabled.
The default value is "+ ". See Section 2.4.8 [Diary and Echo Commands], page 34, for a description of echoing commands.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [echo], page 35, [PS1], page 33, [PS2], page 34.

### 2.4.8 Diary and Echo Commands

Octave's diary feature allows you to keep a $\log$ of all or part of an interactive session by recording the input you type and the output that Octave produces in a separate file.
diary
diary on
diary off
diary filename
[status, diaryfile] = diary
Record a list of all commands and the output they produce, mixed together just as they appear on the terminal.
Valid options are:
on $\quad$ Start recording a session in a file called diary in the current working directory.
off Stop recording the session in the diary file.
filename Record the session in the file named filename.
With no input or output arguments, diary toggles the current diary state.
If output arguments are requested, diary ignores inputs and returns the current status. The boolean status indicates whether recording is on or off, and diaryfile is the name of the file where the session is stored.
See also: [history], page 29, [evalc], page 175.
Sometimes it is useful to see the commands in a function or script as they are being evaluated. This can be especially helpful for debugging some kinds of problems.
echo
echo on
echo off
echo on all
echo off all
echo function on
echo function off
Control whether commands are displayed as they are executed.
Valid options are:
on Enable echoing of commands as they are executed in script files.
off Disable echoing of commands as they are executed in script files.
on all Enable echoing of commands as they are executed in script files and functions.
off all Disable echoing of commands as they are executed in script files and functions.
function on
Enable echoing of commands as they are executed in the named function.
function off
Disable echoing of commands as they are executed in the named function.
With no arguments, echo toggles the current echo state.
See also: [PS4], page 34.

### 2.5 How Octave Reports Errors

Octave reports two kinds of errors for invalid programs.
A parse error occurs if Octave cannot understand something you have typed. For example, if you misspell a keyword,

```
octave:13> function z = f (x, y) z = x ||| 2; endfunction
```

Octave will respond immediately with a message like this:

```
parse error:
    syntax error
>>> function z = f (x, y) z = x ||| y; endfunction
```

For most parse errors, Octave uses a caret ( $\left.{ }^{(\sim `}\right)$ to mark the point on the line where it was unable to make sense of your input. In this case, Octave generated an error message because the keyword for the logical or operator (II) was misspelled. It marked the error at the third ' $\mid$ ' because the code leading up to this was correct but the final ' $\mid$ ' was not understood.

Another class of error message occurs at evaluation time. These errors are called run-time errors, or sometimes evaluation errors, because they occur when your program is being run, or evaluated. For example, if after correcting the mistake in the previous function definition, you type
octave:13> f ()
Octave will respond with

```
error: 'x' undefined near line 1 column 24
error: called from:
error: f at line 1, column 22
```

This error message has several parts, and gives quite a bit of information to help you locate the source of the error. The messages are generated from the point of the innermost error, and provide a traceback of enclosing expressions and function calls.

In the example above, the first line indicates that a variable named ' $x$ ' was found to be undefined near line 1 and column 24 of some function or expression. For errors occurring within functions, lines are counted from the beginning of the file containing the function definition. For errors occurring outside of an enclosing function, the line number indicates the input line number, which is usually displayed in the primary prompt string.

The second and third lines of the error message indicate that the error occurred within the function $f$. If the function $f$ had been called from within another function, for example, g , the list of errors would have ended with one more line:

```
error: g at line 1, column 17
```

These lists of function calls make it fairly easy to trace the path your program took before the error occurred, and to correct the error before trying again.

### 2.6 Executable Octave Programs

Once you have learned Octave, you may want to write self-contained Octave scripts, using the ' $\#$ !' script mechanism. You can do this on GNU systems and on many Unix systems ${ }^{1}$.

Self-contained Octave scripts are useful when you want to write a program which users can invoke without knowing that the program is written in the Octave language. Octave scripts are also used for batch processing of data files. Once an algorithm has been developed and tested in the interactive portion of Octave, it can be committed to an executable script and used again and again on new data files.

As a trivial example of an executable Octave script, you might create a text file named hello, containing the following lines:

```
#! octave-interpreter-name -qf
# a sample Octave program
printf ("Hello, world!\n");
```

(where octave-interpreter-name should be replaced with the full path and name of your Octave binary). Note that this will only work if '\#!' appears at the very beginning of the file. After making the file executable (with the chmod command on Unix systems), you can simply type:
hello
at the shell, and the system will arrange to run Octave as if you had typed:
octave hello
The line beginning with '\#!' lists the full path and filename of an interpreter to be run, and an optional initial command line argument to pass to that interpreter. The operating system then runs the interpreter with the given argument and the full argument list of the executed program. The first argument in the list is the full filename of the Octave executable. The rest of the argument list will either be options to Octave, or data files, or both. The '-qf' options are usually specified in stand-alone Octave programs to prevent them from printing the normal startup message, and to keep them from behaving differently depending on the contents of a particular user's ~/ . octaverc file. See Section 2.1 [Invoking Octave from the Command Line], page 15.

Note that some operating systems may place a limit on the number of characters that are recognized after '\#!'. Also, the arguments appearing in a '\#!' line are parsed differently by various shells/systems. The majority of them group all the arguments together in one string and pass it to the interpreter as a single argument. In this case, the following script:

```
#! octave-interpreter-name -q -f # comment
```

is equivalent to typing at the command line:
octave "-q -f \# comment"
which will produce an error message. Unfortunately, it is not possible for Octave to determine whether it has been called from the command line or from a ' $\#$ !' script, so some care is needed when using the '\#!' mechanism.

[^0]
### 2.6.1 Passing Arguments to Executable Scripts

Note that when Octave is started from an executable script, the built-in function argv returns a cell array containing the command line arguments passed to the executable Octave script, not the arguments passed to the Octave interpreter on the '\#!' line of the script. For example, the following program will reproduce the command line that was used to execute the script, not '-qf'.

```
#! /bin/octave -qf
printf ("%s", program_name ());
arg_list = argv ();
for i = 1:nargin
    printf (" %s", arg_list{i});
endfor
printf ("\n");
```


### 2.6.2 Dual-Purpose Executable Scripts and Octave Functions

To write m-files that can act as executable programs when called from the shell or as normal functions when called from within Octave, use default input arguments initialized with the argv function.

If a function is called from the shell Octave will not pass any input parameters to the function and therefore the default argument is used. But when a function is called from the interpreter any arguments are passed to the function and these override the default.

Additionally, the file must end with the extension .m so that the interpreter will recognize it as an Octave function. Finally, the output from argv is a cell array of strings. It may be necessary to convert this to a numeric value with str2double or str2num before processing.

As a complete example, consider the following code located in the file mysin.m.

```
#! /bin/octave -qf
function retval = mysin (x = str2double (argv(){end}))
    retval = sin (x)
endfunction
```

This can be called from the shell with
mysin.m 1.5
or from Octave with
mysin (1.5)

### 2.7 Comments in Octave Programs

A comment is some text that is included in a program for the sake of human readers, and which is NOT an executable part of the program. Comments can explain what the program does, and how it works. Nearly all programming languages have provisions for comments, because programs are typically hard to understand without them.

### 2.7.1 Single Line Comments

In the Octave language, a comment starts with either the sharp sign character, '\#', or the percent symbol ${ }^{\circ} \%$ and continues to the end of the line. Any text following the sharp sign
or percent symbol is ignored by the Octave interpreter and not executed. The following example shows whole line and partial line comments.

```
function countdown
    # Count down for main rocket engines
    disp (3);
    disp (2);
    disp (1);
    disp ("Blast Off!"); # Rocket leaves pad
endfunction
```


### 2.7.2 Block Comments

Entire blocks of code can be commented by enclosing the code between matching '\#\{' and '\#\}' or ' $\%\{$ ' and ' $\%\}$ ' markers. For example,

```
function quick_countdown
    # Count down for main rocket engines
    disp (3);
    #{
    disp (2);
    disp (1);
    #}
    disp ("Blast Off!"); # Rocket leaves pad
endfunction
```

will produce a very quick countdown from '3' to "Blast Off" as the lines "disp (2);" and "disp (1);" won't be executed.

The block comment markers must appear alone as the only characters on a line (excepting whitespace) in order to be parsed correctly.

### 2.7.3 Comments and the Help System

The help command (see Section 2.3 [Getting Help], page 21) is able to find the first block of comments in a function and return those as a documentation string. This means that the same commands used to get help on built-in functions are available for properly formatted user-defined functions. For example, after defining the function $f$ below,

```
function xdot = f (x, t)
# usage: f (x, t)
#
# This function defines the right-hand
# side functions for a set of nonlinear
# differential equations.
    r = 0.25;
endfunction
```

the command help $f$ produces the output

```
usage: f (x, t)
This function defines the right-hand
side functions for a set of nonlinear
differential equations.
```

Although it is possible to put comment lines into keyboard-composed, throw-away Octave programs, it usually isn't very useful because the purpose of a comment is to help you or another person understand the program at a later time.

The help parser currently only recognizes single line comments (see Section 2.7.1 [Single Line Comments], page 38) and not block comments for the initial help text.

## 3 Data Types

All versions of Octave include a number of built-in data types, including real and complex scalars and matrices, character strings, a data structure type, and an array that can contain all data types.

It is also possible to define new specialized data types by writing a small amount of C++ code. On some systems, new data types can be loaded dynamically while Octave is running, so it is not necessary to recompile all of Octave just to add a new type. See Appendix A [External Code Interface], page 1009, for more information about Octave's dynamic linking capabilities. Section 3.2 [User-defined Data Types], page 46, describes what you must do to define a new data type for Octave.
typeinfo ()
typeinfo (expr)
Return the type of the expression expr, as a string.
If expr is omitted, return a cell array of strings containing all the currently installed data types.
See also: [class], page 41, [isa], page 41.

### 3.1 Built-in Data Types

The standard built-in data types are real and complex scalars and matrices, ranges, character strings, a data structure type, and cell arrays. Additional built-in data types may be added in future versions. If you need a specialized data type that is not currently provided as a built-in type, you are encouraged to write your own user-defined data type and contribute it for distribution in a future release of Octave.

The data type of a variable can be determined and changed through the use of the following functions.

```
classname = class (obj)
class (s,id)
class (s, id, p,...)
```

Return the class of the object obj, or create a class with fields from structure $s$ and name (string) id.
Additional arguments name a list of parent classes from which the new class is derived.
See also: [typeinfo], page 41, [isa], page 41.
isa (obj, classname)
Return true if obj is an object from the class classname.
classname may also be one of the following class categories:
"float" Floating point value comprising classes "double" and "single".
"integer"
Integer value comprising classes (u)int8, (u)int16, (u)int32, (u)int64.
"numeric"
Numeric value comprising either a floating point or integer value.

If classname is a cell array of string, a logical array of the same size is returned, containing true for each class to which obj belongs to.

See also: [class], page 41, [typeinfo], page 41.
cast (val, "type")
cast (val, "like", var)
Convert val to data type type.
The input val may be a scalar, vector, or matrix of a class that is convertible to the target class (see below).
If a variable var is specified after "like", val is converted to the same data type and sparsity attribute. If var is complex, val will be complex, too.
var may be and type may name any of the following built-in numeric classes:

```
"double"
"single"
"logical"
"char"
"int8"
"int16"
"int32"
"int64"
"uint8"
"uint16"
"uint32"
"uint64"
```

The value val may be modified to fit within the range of the new type.
Examples:

```
cast (-5, "uint8")
    # 0
cast (300, "int8")
    => 127
```

Programming Note: This function relies on the object val having a conversion method named type. User-defined classes may implement only a subset of the full list of types shown above. In that case, it may be necessary to call cast twice in order to reach the desired type. For example, the conversion to double is nearly always implemented, but the conversion to uint8 might not be. In that case, the following code will work:

```
cast (cast (user_defined_val, "double"), "uint8")
```

See also: [typecast], page 42, [int8], page 58, [uint8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59, [double], page 51, [single], page 57, [logical], page 64, [char], page 78, [class], page 41, [typeinfo], page 41.
$y=$ typecast ( $x$, "class")
Return a new array $y$ resulting from interpreting the data of $x$ in memory as data of the numeric class class.

Both the class of $x$ and class must be one of the built-in numeric classes:

```
"logical"
"char"
"int8"
"int16"
"int32"
"int64"
"uint8"
"uint16"
"uint32"
"uint64"
"double"
"single"
"double complex"
"single complex"
```

the last two are only used with class; they indicate that a complex-valued result is requested. Complex arrays are stored in memory as consecutive pairs of real numbers. The sizes of integer types are given by their bit counts. Both logical and char are typically one byte wide; however, this is not guaranteed by C++. If your system is IEEE conformant, single and double will be 4 bytes and 8 bytes wide, respectively. "logical" is not allowed for class.

If the input is a row vector, the return value is a row vector, otherwise it is a column vector.

If the bit length of $x$ is not divisible by that of class, an error occurs.
An example of the use of typecast on a little-endian machine is

```
x = uint16 ([1, 65535]);
typecast (x, "uint8")
# [ 1, 0, 255, 255]
```

See also: [cast], page 42, [bitpack], page 43, [bitunpack], page 44, [swapbytes], page 43.
swapbytes (x)
Swap the byte order on values, converting from little endian to big endian and vice versa.

For example:

```
swapbytes (uint16 (1:4))
# 256 512 768 1024
```

See also: [typecast], page 42, [cast], page 42.
$\mathrm{y}=$ bitpack ( $\mathrm{x}, \mathrm{class}$ )
Return a new array $y$ resulting from interpreting the logical array x as raw bit patterns for data of the numeric class class.
class must be one of the built-in numeric classes:

```
"double"
"single"
"double complex"
"single complex"
"char"
"int8"
"int16"
"int32"
"int64"
"uint8"
"uint16"
"uint32"
"uint64"
```

The number of elements of $x$ should be divisible by the bit length of class. If it is not, excess bits are discarded. Bits come in increasing order of significance, i.e., x(1) is bit $0, x(2)$ is bit 1 , etc.
The result is a row vector if $x$ is a row vector, otherwise it is a column vector.
See also: [bitunpack], page 44, [typecast], page 42.
$y=$ bitunpack (x)
Return a logical array $y$ corresponding to the raw bit patterns of $x$.
$x$ must belong to one of the built-in numeric classes:

```
"double"
"single"
"char"
"int8"
"int16"
"int32"
"int64"
"uint8"
"uint16"
"uint32"
"uint64"
```

The result is a row vector if $x$ is a row vector; otherwise, it is a column vector.
See also: [bitpack], page 43, [typecast], page 42.

### 3.1.1 Numeric Objects

Octave's built-in numeric objects include real, complex, and integer scalars and matrices. All built-in floating point numeric data is currently stored as double precision numbers. On systems that use the IEEE floating point format, values in the range of approximately $2.2251 \times 10^{-308}$ to $1.7977 \times 10^{308}$ can be stored, and the relative precision is approximately $2.2204 \times 10^{-16}$. The exact values are given by the variables realmin, realmax, and eps, respectively.

Matrix objects can be of any size, and can be dynamically reshaped and resized. It is easy to extract individual rows, columns, or submatrices using a variety of powerful indexing features. See Section 8.1 [Index Expressions], page 151.

See Chapter 4 [Numeric Data Types], page 51, for more information.

### 3.1.2 Missing Data

It is possible to represent missing data explicitly in Octave using NA (short for "Not Available"). Missing data can only be represented when data is represented as floating point numbers. In this case missing data is represented as a special case of the representation of NaN .

```
val = NA
val = NA (n)
val = NA (n,m)
val = NA (n, m,k,\ldots)
val = NA (..., "like", var))
val = NA (..., class)
```

Return a scalar, matrix, or N -dimensional array whose elements are all equal to the special constant used to designate missing values.
Note that NA always compares not equal to NA (NA != NA). To find NA values, use the isna function.
When called with no arguments, return a scalar with the value ' NA '.
When called with a single argument, return a square matrix with the dimension specified.

When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
If a variable var is specified after "like", the output val will have the same data type, complexity, and sparsity as var.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [isna], page 45.
isna (x)
Return a logical array which is true where the elements of $x$ are NA (missing) values and false where they are not.
For example:

```
isna ([13, Inf, NA, NaN])
    # [0, 0, 1, 0]
```

See also: [isnan], page 512, [isinf], page 512, [isfinite], page 513.

### 3.1.3 String Objects

A character string in Octave consists of a sequence of characters enclosed in either doublequote or single-quote marks. Internally, Octave currently stores strings as matrices of characters. All the indexing operations that work for matrix objects also work for strings.

See Chapter 5 [Strings], page 71, for more information.

### 3.1.4 Data Structure Objects

Octave's data structure type can help you to organize related objects of different types. The current implementation uses an associative array with indices limited to strings, but the syntax is more like C-style structures.

See Section 6.1 [Structures], page 111, for more information.

### 3.1.5 Cell Array Objects

A Cell Array in Octave is general array that can hold any number of different data types.
See Section 6.3 [Cell Arrays], page 124, for more information.

### 3.2 User-defined Data Types

Someday I hope to expand this to include a complete description of Octave's mechanism for managing user-defined data types. Until this feature is documented here, you will have to make do by reading the code in the ov.h, ops.h, and related files from Octave's src directory.

### 3.3 Object Sizes

The following functions allow you to determine the size of a variable or expression. These functions are defined for all objects. They return -1 when the operation doesn't make sense. For example, Octave's data structure type doesn't have rows or columns, so the rows and columns functions return -1 for structure arguments.
ndims (a)
Return the number of dimensions of a.
For any array, the result will always be greater than or equal to 2 . Trailing singleton dimensions are not counted, i.e., tailing dimensions $d$ greater than 2 , for which size $(a, d)=1$.

```
ndims (ones (4, 1, 2, 1))
=>3
```

See also: [size], page 47.
columns (a)
Return the number of columns of $a$. This is equivalent to size ( $a, 2$ ).
See also: [rows], page 46, [size], page 47, [length], page 47, [numel], page 46, [isscalar], page 68, [isvector], page 68, [ismatrix], page 68.
rows (a)
Return the number of rows of $a$. This is equivalent to size ( $a, 1$ ).
See also: [columns], page 46, [size], page 47, [length], page 47, [numel], page 46, [isscalar], page 68, [isvector], page 68, [ismatrix], page 68.
numel (a)
numel (a, idx1, idx2,...)
Return the number of elements in the object a.

Optionally, if indices idx1, idx2, . . are supplied, return the number of elements that would result from the indexing

```
a(idx1, idx2, ...)
```

Note that the indices do not have to be scalar numbers. For example,

```
a = 1;
b = ones (2, 3);
numel (a, b)
```

will return 6 , as this is the number of ways to index with $b$. Or the index could be the string ":" which represents the colon operator. For example,

```
a = ones (5, 3);
numel (a, 2, ":")
```

will return 3 as the second row has three column entries.
This method is also called when an object appears as lvalue with cs-list indexing, i.e., object\{...\} or object(...).field.
See also: [size], page 47, [length], page 47, [ndims], page 46.

## length (a)

Return the length of the object a.
The length is 0 for empty objects, 1 for scalars, and the number of elements for vectors. For matrix or N-dimensional objects, the length is the number of elements along the largest dimension (equivalent to max (size (a))).
See also: [numel], page 46, [size], page 47.
sz = size (a)
dim_sz = size (a, dim)
dim_sz $=\operatorname{size}(a, d 1, d 2, \ldots)$
[rows, cols, ..., dim_ $\left.N_{-} s z\right]=$ size (...)
Return a row vector with the size (number of elements) of each dimension for the object a.
When given a second argument, dim, return the size of the corresponding dimension. If dim is a vector, return each of the corresponding dimensions. Multiple dimensions may also be specified as separate arguments.
With a single output argument, size returns a row vector. When called with multiple output arguments, size returns the size of dimension N in the Nth argument. The number of rows, dimension 1 , is returned in the first argument, the number of columns, dimension 2 , is returned in the second argument, etc. If there are more dimensions in a than there are output arguments, size returns the total number of elements in the remaining dimensions in the final output argument.
Example 1: single row vector output

```
size ([1, 2; 3, 4; 5, 6])
    # [3, 2 ]
```

Example 2: number of elements in 2nd dimension (columns)

```
size ([1, 2; 3, 4; 5, 6], 2)
    =>2
```

Example 3: number of output arguments $==$ number of dimensions

```
[nr, nc] = size ([1, 2; 3, 4; 5, 6])
    #nr = 3
    # nc = 2
```

Example 4: number of output arguments < number of dimensions

```
[nr, remainder] = size (ones (2, 3, 4, 5))
    # nr = 2
    => remainder = 60
```

See also: [numel], page 46, [ndims], page 46, [length], page 47, [rows], page 46, [columns], page 46, [size_equal], page 49, [common_size], page 513.
isempty (a)
Return true if a is an empty matrix (any one of its dimensions is zero).
See also: [isnull], page 48, [isa], page 41.
isnull (x)
Return true if $x$ is a special null matrix, string, or single quoted string.
Indexed assignment with such a null value on the right-hand side should delete array elements. This function is used in place of isempty when overloading the indexed assignment method (subsasgn) for user-defined classes. isnull is used to distinguish between these two cases:
$A(I)=[]$
and
$X=[] ; A(I)=X$
In the first assignment, the right-hand side is [] which is a special null value. As long as the index $I$ is not empty, this code should delete elements from $A$ rather than perform assignment.
In the second assignment, the right-hand side is empty (because $X$ is []), but it is not null. This code should assign the empty value to elements in $A$.
An example from Octave's built-in char class demonstrates the interpreter behavior when isnull is used correctly.

```
str = "Hello World";
nm = "Wally";
str}(7:end) = nm # indexed assignment
    | str = Hello Wally
str(7:end) = "" # indexed deletion
    str = Hello
```

See also: [isempty], page 48, [isindex], page 157.
sizeof (val)
Return the size of val in bytes.
See also: [whos], page 143.
size_equal ( $a, b, \ldots$ )
Return true if the dimensions of all arguments agree.
Trailing singleton dimensions are ignored. When called with a single argument, or no argument, size_equal returns true.
See also: [size], page 47, [numel], page 46, [ndims], page 46, [common_size], page 513. squeeze ( $x$ )

Remove singleton dimensions from $x$ and return the result.
Note that for compatibility with matlab, all objects have a minimum of two dimensions and row vectors are left unchanged.
See also: [reshape], page 518.

## 4 Numeric Data Types

A numeric constant may be a scalar, a vector, or a matrix, and it may contain complex values.

The simplest form of a numeric constant, a scalar, is a single number. Note that by default numeric constants are represented within Octave by IEEE 754 double precision (binary64) floating-point format (complex constants are stored as pairs of binary64 values). It is, however, possible to represent real integers as described in Section 4.4 [Integer Data Types], page 58.

If the numeric constant is a real integer, it can be defined in decimal, hexadecimal, or binary notation. Hexadecimal notation starts with ' $0 x$ ' or ' $0 X$ ', binary notation starts with ' Ob ' or ' OB ', otherwise decimal notation is assumed. As a consequence, ' Ob ' is not a hexadecimal number, in fact, it is not a valid number at all.

For better readability, digits may be partitioned by the underscore separator '_', which is ignored by the Octave interpreter. Here are some examples of real-valued integer constants, which all represent the same value and are internally stored as binary64:

```
42 # decimal notation
0x2A # hexadecimal notation
Ob101010 # binary notation
Ob10_1010 # underscore notation
round (42.1) # also binary64
```

In decimal notation, the numeric constant may be denoted as decimal fraction or even in scientific (exponential) notation. Note that this is not possible for hexadecimal or binary notation. Again, in the following example all numeric constants represent the same value:

```
. }10
1.05e-1
.00105e+2
```

Unlike most programming languages, complex numeric constants are denoted as the sum of real and imaginary parts. The imaginary part is denoted by a real-valued numeric constant followed immediately by a complex value indicator (' $i$ ', ' $j$ ', ' $I$ ', or 'J' which represents $\sqrt{-1}$ ). No spaces are allowed between the numeric constant and the complex value indicator. Some examples of complex numeric constants that all represent the same value:

```
3 + 42i
3 + 42j
3 + 42I
3 + 42J
3.0 + 42.0i
3.0 + 0x2Ai
3.0 + Ob10_1010i
0.3e1 + 420e-1i
```

double (x)

Convert $x$ to double precision type.
See also: [single], page 57 .

```
complex (x)
complex (re, im)
```

Return a complex value from real arguments.
With 1 real argument $x$, return the complex result $x+0 i$.
With 2 real arguments, return the complex result re $+i m i$. complex can often be more convenient than expressions such as $\mathrm{a}+\mathrm{b} * \mathrm{i}$. For example:

```
complex ([1, 2], [3, 4])
    [ [ 1 + 3i 2 + 4i ]
```

See also: [real], page 550, [imag], page 550, [iscomplex], page 67, [abs], page 549, [arg], page 549 .

### 4.1 Matrices

It is easy to define a matrix of values in Octave. The size of the matrix is determined automatically, so it is not necessary to explicitly state the dimensions. The expression

$$
a=[1,2 ; 3,4]
$$

results in the matrix

$$
a=\left[\begin{array}{ll}
1 & 2 \\
3 & 4
\end{array}\right]
$$

Elements of a matrix may be arbitrary expressions, provided that the dimensions all make sense when combining the various pieces. For example, given the above matrix, the expression
[ a, a ]
produces the matrix
ans $=$
$\begin{array}{llll}1 & 2 & 1 & 2\end{array}$
$\begin{array}{llll}3 & 4 & 3 & 4\end{array}$
but the expression
[ a, 1]
produces the error
error: number of rows must match (1 ! = 2) near line 13, column 6
(assuming that this expression was entered as the first thing on line 13, of course).
Inside the square brackets that delimit a matrix expression, Octave looks at the surrounding context to determine whether spaces and newline characters should be converted into element and row separators, or simply ignored, so an expression like

$$
a=\left[\begin{array}{ll}
1 & 2 \\
3 & 4
\end{array}\right]
$$

will work. However, some possible sources of confusion remain. For example, in the expression
[ 1 - 1 ]
the ' - ' is treated as a binary operator and the result is the scalar 0 , but in the expression
[ $\left.\begin{array}{lll}1 & -1\end{array}\right]$
the ' - ' is treated as a unary operator and the result is the vector [ $1,-1$ ]. Similarly, the expression

```
[ sin (pi) ]
```

will be parsed as
[ sin, (pi) ]
and will result in an error since the sin function will be called with no arguments. To get around this, you must omit the space between sin and the opening parenthesis, or enclose the expression in a set of parentheses:
[ ( $\sin (\mathrm{pi}))$ ]
Whitespace surrounding the single quote character ('"', used as a transpose operator and for delimiting character strings) can also cause confusion. Given $\mathrm{a}=1$, the expression
[ 1 a' ]
results in the single quote character being treated as a transpose operator and the result is the vector [ 1, 1], but the expression
[ 1 a ' ]
produces the error message

```
parse error:
    syntax error
>>> [llll
```

because not doing so would cause trouble when parsing the valid expression

```
[ a 'foo' ]
```

For clarity, it is probably best to always use commas and semicolons to separate matrix elements and rows.

The maximum number of elements in a matrix is fixed when Octave is compiled. The allowable number can be queried with the function sizemax. Note that other factors, such as the amount of memory available on your machine, may limit the maximum size of matrices to something smaller.

## sizemax ()

Return the largest value allowed for the size of an array.
If Octave is compiled with 64 -bit indexing, the result is of class int64, otherwise it is of class int 32 . The maximum array size is slightly smaller than the maximum value allowable for the relevant class as reported by intmax.
See also: [intmax], page 59 .
When you type a matrix or the name of a variable whose value is a matrix, Octave responds by printing the matrix in with neatly aligned rows and columns. If the rows of the matrix are too large to fit on the screen, Octave splits the matrix and displays a header before each section to indicate which columns are being displayed. You can use the following variables to control the format of the output.

```
val = output_precision ()
old_val = output_precision (new_val)
output_precision (new_val, "local")
```

Query or set the internal variable that specifies the minimum number of significant figures to display for numeric output.
Note that regardless of the value set for output_precision, the number of digits of precision displayed is limited to 16 for double precision values and 7 for single precision values.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [format], page 272, [fixed_point_format], page 54.
It is possible to achieve a wide range of output styles by using different values of output_ precision. Reasonable combinations can be set using the format function. See Section 14.1 [Basic Input and Output], page 271.

```
val = split_long_rows ()
old_val = split_long_rows (new_val)
split_long_rows (new_val, "local")
```

Query or set the internal variable that controls whether rows of a matrix may be split when displayed to a terminal window.
If the rows are split, Octave will display the matrix in a series of smaller pieces, each of which can fit within the limits of your terminal width and each set of rows is labeled so that you can easily see which columns are currently being displayed. For example:

```
octave:13> rand (2,10)
ans =
Columns 1 through 6:
                    0.75883 0.93290
                        0.75697}00.51942 0.40031 0.61784 0.92309 0.40201
Columns 7 through 10:
    0.90174
```

    When called from inside a function with the "local" option, the variable is changed
    locally for the function and any subroutines it calls. The original variable value is
    restored when exiting the function.
    See also: [format], page 272.
    Octave automatically switches to scientific notation when values become very large or very small. This guarantees that you will see several significant figures for every value in a matrix. If you would prefer to see all values in a matrix printed in a fixed point format, you can use the function fixed_point_format. But doing so is not recommended, because it can produce output that can easily be misinterpreted.

```
val = fixed_point_format ()
old_val = fixed_point_format (new_val)
fixed_point_format (new_val, "local")
```

Query or set the internal variable that controls whether Octave will use a scaled format to print matrix values.

The scaled format prints a scaling factor on the first line of output chosen such that the largest matrix element can be written with a single leading digit. For example:

```
fixed_point_format (true)
logspace (1, 7, 5)'
ans =
            1.0e+07 *
            0.00000
            0.00003
            0.00100
            0.03162
            1.00000
```

Notice that the first value appears to be 0 when it is actually 1 . Because of the possibility for confusion you should be careful about enabling fixed_point_format.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [format], page 272, [output_precision], page 53.

### 4.1.1 Empty Matrices

A matrix may have one or both dimensions zero, and operations on empty matrices are handled as described by Carl de Boor in An Empty Exercise, SIGNUM, Volume 25, pages 2-6, 1990 and C. N. Nett and W. M. Haddad, in A System-Theoretic Appropriate Realization of the Empty Matrix Concept, IEEE Transactions on Automatic Control, Volume 38, Number 5, May 1993. Briefly, given a scalar $s$, an $m \times n$ matrix $M_{m \times n}$, and an $m \times n$ empty matrix []$_{m \times n}$ (with either one or both dimensions equal to zero), the following are true:

$$
\begin{aligned}
s \cdot[]_{m \times n}=[]_{m \times n} \cdot s & =[]_{m \times n} \\
{[]_{m \times n}+[]_{m \times n} } & =[]_{m \times n} \\
{[]_{0 \times m} \cdot M_{m \times n} } & =[]_{0 \times n} \\
M_{m \times n} \cdot[]_{n \times 0} & =[]_{m \times 0} \\
{[]_{m \times 0} \cdot[]_{0 \times n} } & =0_{m \times n}
\end{aligned}
$$

By default, dimensions of the empty matrix are printed along with the empty matrix symbol, '[]'. The built-in variable print_empty_dimensions controls this behavior.

```
val = print_empty_dimensions ()
old_val = print_empty_dimensions (new_val)
```

print_empty_dimensions (new_val, "local")
Query or set the internal variable that controls whether the dimensions of empty matrices are printed along with the empty matrix symbol, '[]'.
For example, the expression
zeros (3, 0)
will print

```
ans = [] (3x0)
```

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [format], page 272.
Empty matrices may also be used in assignment statements as a convenient way to delete rows or columns of matrices. See Section 8.6 [Assignment Expressions], page 169.

When Octave parses a matrix expression, it examines the elements of the list to determine whether they are all constants. If they are, it replaces the list with a single matrix constant.

### 4.2 Ranges

A range is a convenient way to write a row vector with evenly spaced elements. A range expression is defined by the value of the first element in the range, an optional value for the increment between elements, and a maximum value which the elements of the range will not exceed. The base, increment, and limit are separated by colons (the ':' character) and may contain any arithmetic expressions and function calls. If the increment is omitted, it is assumed to be 1 . For example, the range

1 : 5
defines the set of values [ $1,2,3,4,5]$, and the range
$1: 3: 5$
defines the set of values [ 1, 4].
Although a range constant specifies a row vector, Octave does not normally convert range constants to vectors unless it is necessary to do so. This allows you to write a constant like 1 : 10000 without using 80,000 bytes of storage on a typical 32 -bit workstation.

A common example of when it does become necessary to convert ranges into vectors occurs when they appear within a vector (i.e., inside square brackets). For instance, whereas $\mathrm{x}=0$ : 0.1 : 1;
defines $x$ to be a variable of type range and occupies 24 bytes of memory, the expression
$y=[0: 0.1$ : 1];
defines $y$ to be of type matrix and occupies 88 bytes of memory.
This space saving optimization may be disabled using the function optimize_range.

```
val = optimize_range ()
old_val = optimize_range (new_val)
optimize_range (new_val, "local")
```

Query or set whether a special space-efficient format is used for storing ranges.

The default value is true. If this option is set to false, Octave will store ranges as full matrices.
When called from inside a function with the "local" option, the setting is changed locally for the function and any subroutines it calls. The original setting is restored when exiting the function.
See also: [optimize_diagonal_matrix], page 651, [optimize_permutation_matrix], page 651.

Note that the upper (or lower, if the increment is negative) bound on the range is not always included in the set of values, and that ranges defined by floating point values can produce surprising results because Octave uses floating point arithmetic to compute the values in the range. If it is important to include the endpoints of a range and the number of elements is known, you should use the linspace function instead (see Section 16.3 [Special Utility Matrices], page 525).

When adding a scalar to a range, subtracting a scalar from it (or subtracting a range from a scalar) and multiplying by scalar, Octave will attempt to avoid unpacking the range and keep the result as a range, too, if it can determine that it is safe to do so. For instance, doing

```
a = 2*(1:1e7) - 1;
```

will produce the same result as 1:2:2e7-1, but without ever forming a vector with ten million elements.

Using zero as an increment in the colon notation, as 1:0:1 is not allowed, because a division by zero would occur in determining the number of range elements. However, ranges with zero increment (i.e., all elements equal) are useful, especially in indexing, and Octave allows them to be constructed using the built-in function ones. Note that because a range must be a row vector, ones $(1,10)$ produces a range, while ones $(10,1)$ does not.

When Octave parses a range expression, it examines the elements of the expression to determine whether they are all constants. If they are, it replaces the range expression with a single range constant.

### 4.3 Single Precision Data Types

Octave includes support for single precision data types, and most of the functions in Octave accept single precision values and return single precision answers. A single precision variable is created with the single function.
single (x)
Convert $x$ to single precision type.
See also: [double], page 51.
for example:

```
sngl = single (rand (2, 2))
        \(\Rightarrow\) sngl =
            \(0.37569 \quad 0.92982\)
            \(0.11962 \quad 0.50876\)
class (sngl)
            \(\Rightarrow\) single
```

Many functions can also return single precision values directly. For example

```
ones (2, 2, "single")
zeros (2, 2, "single")
eye (2, 2, "single")
rand (2, 2, "single")
NaN (2, 2, "single")
NA (2, 2, "single")
Inf (2, 2, "single")
```

will all return single precision matrices.

### 4.4 Integer Data Types

Octave supports integer matrices as an alternative to using double precision. It is possible to use both signed and unsigned integers represented by $8,16,32$, or 64 bits. It should be noted that most computations require floating point data, meaning that integers will often change type when involved in numeric computations. For this reason integers are most often used to store data, and not for calculations.

In general most integer matrices are created by casting existing matrices to integers. The following example shows how to cast a matrix into 32 bit integers.

```
float = rand (2, 2)
    f float = 0.37569 0.92982
        0.11962 0.50876
integer = int32 (float)
    => integer = 0 1
        0 1
```

As can be seen, floating point values are rounded to the nearest integer when converted.

```
isinteger (x)
```

Return true if $x$ is an integer object (int8, uint8, int16, etc.).
Note that isinteger (14) is false because numeric constants in Octave are double precision floating point values.
See also: [isfloat], page 67, [ischar], page 73, [islogical], page 67, [isstring], page 73, [isnumeric], page 67, [isa], page 41.
int8 (x)
Convert $x$ to 8-bit integer type.
See also: [uint8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59.
uint8 (x)
Convert $x$ to unsigned 8-bit integer type.
See also: [int8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59.
int16 (x)
Convert $x$ to 16-bit integer type.

See also: [int8], page 58, [uint8], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59.
uint16 (x)
Convert $x$ to unsigned 16-bit integer type.
See also: [int8], page 58, [uint8], page 58, [int16], page 58, [int32], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59.
int32 (x)
Convert $x$ to 32 -bit integer type.
See also: [int8], page 58, [uint8], page 58, [int16], page 58, [uint16], page 59, [uint32], page 59, [int64], page 59, [uint64], page 59.
uint32 (x)
Convert x to unsigned 32-bit integer type.
See also: [int8], page 58, [uint8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [int64], page 59, [uint64], page 59.
int64 (x)
Convert $x$ to 64-bit integer type.
See also: [int8], page 58, [uint8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [uint64], page 59.
uint64 (x)
Convert $x$ to unsigned 64 -bit integer type.
See also: [int8], page 58, [uint8], page 58, [int16], page 58, [uint16], page 59, [int32], page 59, [uint32], page 59, [int64], page 59.

```
intmax ()
```

intmax ("type")
intmax (var)

Return the largest integer that can be represented by a specific integer type.
The input is either a string "type" specifying an integer type, or it is an existing integer variable var.
Possible values for type are
"int8" signed 8-bit integer.
"int16" signed 16-bit integer.
"int32" signed 32-bit integer.
"int64" signed 64-bit integer.
"uint8" unsigned 8-bit integer.
"uint16" unsigned 16-bit integer.
"uint32" unsigned 32-bit integer.
"uint64" unsigned 64-bit integer.

The default for type is "int32".
Example Code - query an existing variable

```
x = int8 (1);
intmax (x)
    => 127
```

See also: [intmin], page 60, [flintmax], page 60.

```
intmin ()
intmin ("type")
intmin (var) integer variable var.
Possible values for type are
"int8" signed 8-bit integer.
"int16" signed 16-bit integer.
"int32" signed 32-bit integer.
"int64" signed 64-bit integer.
"uint8" unsigned 8-bit integer.
"uint16" unsigned 16-bit integer.
"uint32" unsigned 32-bit integer.
"uint64" unsigned 64-bit integer.
```

Return the smallest integer that can be represented by a specific integer type.
The input is either a string "type" specifying an integer type, or it is an existing

The default for type is "int32".
Example Code - query an existing variable

```
x = int8 (1);
intmin (x)
        => -128
```

See also: [intmax], page 59, [flintmax], page 60.
flintmax ()
flintmax ("double")
flintmax ("single")
flintmax (var)
Return the largest integer that can be represented consecutively in a floating point value.

The input is either a string specifying a floating point type, or it is an existing floating point variable var.
The default type is "double", but "single" is a valid option. On IEEE 754 compatible systems, flintmax is $2^{53}$ for "double" and $2^{24}$ for "single".

Example Code - query an existing variable

```
x = single (1);
flintmax (x)
    # 16777216
```

See also: [intmax], page 59, [realmax], page 583, [realmin], page 583.

### 4.4.1 Integer Arithmetic

While many numerical computations can't be carried out in integers, Octave does support basic operations like addition and multiplication on integers. The operators,,.$+- *$, and ./ work on integers of the same type. So, it is possible to add two 32 bit integers, but not to add a 32 bit integer and a 16 bit integer.

When doing integer arithmetic one should consider the possibility of underflow and overflow. This happens when the result of the computation can't be represented using the chosen integer type. As an example it is not possible to represent the result of $10-20$ when using unsigned integers. Octave makes sure that the result of integer computations is the integer that is closest to the true result. So, the result of $10-20$ when using unsigned integers is zero.

When doing integer division Octave will round the result to the nearest integer. This is different from most programming languages, where the result is often floored to the nearest integer. So, the result of int32 (5) ./ int32 (8) is 1.
idivide ( $x, y, o p$ )
Integer division with different rounding rules.
The standard behavior of integer division such as a ./b is to round the result to the nearest integer. This is not always the desired behavior and idivide permits integer element-by-element division to be performed with different treatment for the fractional part of the division as determined by the op flag. op is a string with one of the values:
"fix" Calculate a./b with the fractional part rounded towards zero.
"round" Calculate a ./b with the fractional part rounded towards the nearest integer.
"floor" Calculate a ./ b with the fractional part rounded towards negative infinity.
"ceil" Calculate a ./ b with the fractional part rounded towards positive infinity.

If $o p$ is not given it defaults to "fix". An example demonstrating these rounding rules is

```
idivide (int8 ([-3, 3]), int8 (4), "fix")
    \(\Rightarrow \quad 0 \quad 0\)
idivide (int8 ([-3, 3]), int8 (4), "round")
    \(\Rightarrow \quad-1 \quad 1\)
idivide (int8 ([-3, 3]), int8 (4), "floor")
    \(\Rightarrow \quad-1 \quad 0\)
idivide (int8 ([-3, 3]), int8 (4), "ceil")
    \(\Rightarrow \quad 0 \quad 1\)
```

See also: [ldivide], page 162, [rdivide], page 163.

### 4.5 Bit Manipulations

Octave provides a number of functions for the manipulation of numeric values on a bit by bit basis. The basic functions to set and obtain the values of individual bits are bitset and bitget.
$C=$ bitset ( $A, n$ )
$C=$ bitset (A, n, val)
Set or reset bit(s) $n$ of the unsigned integers in $A$.
The least significant bit is $n=1$. val $=0$ resets bits and val $=1$ sets bits. If no val is specified it defaults to 1 (set bit). All inputs must be the same size or scalars.
Example 1: Set multiple bits

```
x = bitset (1, 3:5)
    | x =
    5 9 17
dec2bin (x)
            #
                0 0 1 0 1
                0 1 0 0 1
            10001
```

Example 2: Reset and set bits

```
x = bitset ([15 14], 1, [0 1])
    # x =
```

            \(14 \quad 15\)
    See also: [bitand], page 63, [bitor], page 63, [bitxor], page 63, [bitget], page 62, [bitcmp], page 63, [bitshift], page 63, [intmax], page 59, [flintmax], page 60.
$c=\operatorname{bitget}(A, n)$
Return the status of bit(s) $n$ of the unsigned integers in $A$.
The least significant bit is $n=1$.

```
bitget (100, 8:-1:1)
```



See also: [bitand], page 63, [bitor], page 63, [bitxor], page 63, [bitset], page 62, [bitcmp], page 63, [bitshift], page 63, [intmax], page 59, [flintmax], page 60.

The arguments to all of Octave's bitwise operations can be scalar or arrays, except for bitcmp, whose $k$ argument must a scalar. In the case where more than one argument is an array, then all arguments must have the same shape, and the bitwise operator is applied to each of the elements of the argument individually. If at least one argument is a scalar and one an array, then the scalar argument is duplicated. Therefore
bitget (100, 8:-1:1)
is the same as
bitget (100 * ones (1, 8), 8:-1:1)
It should be noted that all values passed to the bit manipulation functions of Octave are treated as integers. Therefore, even though the example for bitset above passes the floating point value 10 , it is treated as the bits [1, 0, 1, 0] rather than the bits of the native floating point format representation of 10 .

As the maximum value that can be represented by a number is important for bit manipulation, particularly when forming masks, Octave supplies two utility functions: flintmax for floating point integers, and intmax for integer objects (uint8, int64, etc.).

Octave also includes the basic bitwise 'and', 'or', and 'exclusive or' operators.
bitand (x, y)
Return the bitwise AND of non-negative integers.
$x, y$ must be in the range [0,intmax]
See also: [bitor], page 63, [bitxor], page 63, [bitset], page 62, [bitget], page 62, [bitcmp], page 63, [bitshift], page 63, [intmax], page 59, [flintmax], page 60.
bitor (x, y)
Return the bitwise OR of non-negative integers $x$ and $y$.
See also: [bitor], page 63, [bitxor], page 63, [bitset], page 62, [bitget], page 62, [bitcmp], page 63, [bitshift], page 63, [intmax], page 59, [flintmax], page 60.
bitxor ( $x, y$ )
Return the bitwise XOR of non-negative integers $x$ and $y$.
See also: [bitand], page 63, [bitor], page 63, [bitset], page 62, [bitget], page 62, [bitcmp], page 63, [bitshift], page 63, [intmax], page 59, [flintmax], page 60.

The bitwise 'not' operator is a unary operator that performs a logical negation of each of the bits of the value. For this to make sense, the mask against which the value is negated must be defined. Octave's bitwise 'not' operator is bitcmp.
bitcmp ( $A, k$ )
Return the $k$-bit complement of integers in $A$.
If $k$ is omitted $k=\log 2$ (flintmax) +1 is assumed.

```
bitcmp (7,4)
    # 8
dec2bin (11)
    => 1011
dec2bin (bitcmp (11, 6))
    => 110100
```

See also: [bitand], page 63, [bitor], page 63, [bitxor], page 63, [bitset], page 62, [bitget], page 62, [bitcmp], page 63, [bitshift], page 63, [flintmax], page 60.

Octave also includes the ability to left-shift and right-shift values bitwise.
bitshift (a, k)
bitshift (a, k, n)
Return a $k$ bit shift of $n$-digit unsigned integers in a.
A positive $k$ leads to a left shift; A negative value to a right shift.
If $n$ is omitted it defaults to $64 . n$ must be in the range $[1,64]$.

```
bitshift (eye (3), 1)
#
20
0 2 0
0 2
bitshift (10, [-2, -1, 0, 1, 2])
llllll
```

See also: [bitand], page 63, [bitor], page 63, [bitxor], page 63, [bitset], page 62, [bitget], page 62, [bitcmp], page 63, [intmax], page 59, [flintmax], page 60.

Bits that are shifted out of either end of the value are lost. Octave also uses arithmetic shifts, where the sign bit of the value is kept during a right shift. For example:

```
bitshift (-10, -1)
# -5
bitshift (int8 (-1), -1)
# -1
```

Note that bitshift (int8 ( -1 ), -1 ) is -1 since the bit representation of -1 in the int8 data type is $[1,1,1,1,1,1,1,1]$.

### 4.6 Logical Values

Octave has built-in support for logical values, i.e., variables that are either true or false. When comparing two variables, the result will be a logical value whose value depends on whether or not the comparison is true.

The basic logical operations are \& I, and !, which correspond to "Logical And", "Logical Or", and "Logical Negation". These operations all follow the usual rules of logic.

It is also possible to use logical values as part of standard numerical calculations. In this case true is converted to 1 , and false to 0 , both represented using double precision floating point numbers. So, the result of true $* 22-\mathrm{false} / 6$ is 22 .

Logical values can also be used to index matrices and cell arrays. When indexing with a logical array the result will be a vector containing the values corresponding to true parts of the logical array. The following example illustrates this.

```
data = [ 1, 2; 3, 4 ];
idx = (data <= 2);
data(idx)
    => ans = [ 1; 2 ]
```

Instead of creating the idx array it is possible to replace data(idx) with data( data <= 2 ) in the above code.

Logical values can also be constructed by casting numeric objects to logical values, or by using the true or false functions.

## logical (x)

Convert the numeric object $x$ to logical type.
Any nonzero values will be converted to true (1) while zero values will be converted to false (0). The non-numeric value NaN cannot be converted and will produce an error.
Compatibility Note: Octave accepts complex values as input, whereas matlab issues an error.
See also: [double], page 51, [single], page 57, [char], page 78.

```
val = true (x)
val = true (n,m)
val = true (n, m, k, ...)
val = true (..., "like", var))
```

Return a matrix or N-dimensional array whose elements are all logical 1.
If invoked with a single scalar integer argument, return a square matrix of the specified size.

If invoked with two or more scalar integer arguments, or a vector of integer values, return an array with given dimensions.
If a logical variable var is specified after "like", the output val will have the same sparsity as var.

See also: [false], page 65.
val = false (x)
val = false ( $n, m$ )
val = false (n, m, k,...)
val = false (..., "like", var))
Return a matrix or N-dimensional array whose elements are all logical 0.
If invoked with a single scalar integer argument, return a square matrix of the specified size.

If invoked with two or more scalar integer arguments, or a vector of integer values, return an array with given dimensions.
If a logical variable var is specified after "like", the output val will have the same sparsity as var.
See also: [true], page 65.

### 4.7 Automatic Conversion of Data Types

Many operators and functions can work with mixed data types. For example,

```
uint8 (1) + 1
    =>2
single (1) + 1
    => 2
```

```
min (single (1), 0)
    =0
```

where the results are respectively of types uint8, single, and single respectively. This is done for Matlab compatibility. Valid mixed operations are defined as follows:

| Mixed Operation | Result |
| :--- | :--- |
| double OP single | single |
| double OP integer | integer |
| double OP char | double |
| double OP logical | double |
| single OP integer | integer |
| single OP char | single |
| single OP logical | single |

When functions expect a double but are passed other types, automatic conversion is function-dependent:

```
a = det (int8 ([1 2; 3 4]))
    =>a = -2
class (a)
    => double
a = eig (int8 ([1 2; 3 4]))
    => error: eig: wrong type argument 'int8 matrix'
```

When two operands are both integers but of different widths, then some cases convert them to the wider bitwidth, and other cases throw an error:

```
a = min (int8 (100), int16 (200))
    # 100
class (a)
    => int16
int8 (100) + int16 (200)
    => error: binary operator '+' not implemented
    for 'int8 scalar' by 'int16 scalar' operations
```

For two integer operands, they typically need to both be signed or both be unsigned. Mixing signed and unsigned usually causes an error, even if they are of the same bitwidth.

```
min (int16 (100), uint16 (200))
    f error: min: cannot compute min (int16 scalar, uint16 scalar)
```

In the case of mixed type indexed assignments, the type is not changed. For example,

```
x = ones (2, 2);
x(1, 1) = single (2)
    m = 2 1
        1
```

where x remains of the double precision type.

### 4.8 Predicates for Numeric Objects

Since the type of a variable may change during the execution of a program, it can be necessary to do type checking at run-time. Doing this also allows you to change the behavior of a function depending on the type of the input. As an example, this naive implementation of abs returns the absolute value of the input if it is a real number, and the length of the input if it is a complex number.

```
function \(\mathrm{a}=\mathrm{abs}(\mathrm{x})\)
    if (isreal (x))
        \(\mathrm{a}=\operatorname{sign}(\mathrm{x}) . * \mathrm{x}\);
    elseif (iscomplex (x))
        \(a=\operatorname{sqrt}\left(\operatorname{real}(x) . \wedge 2+\operatorname{imag}(x) .^{\wedge} 2\right) ;\)
    endif
endfunction
```

The following functions are available for determining the type of a variable.
isnumeric (x)
Return true if $x$ is a numeric object, i.e., an integer, real, or complex array.
Logical and character arrays are not considered to be numeric.
See also: [isinteger], page 58, [isfloat], page 67, [isreal], page 67, [iscomplex], page 67, [ischar], page 73, [islogical], page 67, [isstring], page 73, [iscell], page 126, [isstruct], page 119, [isa], page 41.
islogical (x)
isbool (x)
Return true if $x$ is a logical object.
See also: [ischar], page 73, [isfloat], page 67, [isinteger], page 58, [isstring], page 73, [isnumeric], page 67, [isa], page 41.
isfloat (x)
Return true if $x$ is a floating-point numeric object.
Objects of class double or single are floating-point objects.
See also: [isinteger], page 58, [ischar], page 73, [islogical], page 67, [isnumeric], page 67, [isstring], page 73, [isa], page 41.
isreal (x)
Return true if $x$ is a non-complex matrix or scalar.
For compatibility with MATLAB, this includes logical and character matrices.
See also: [iscomplex], page 67, [isnumeric], page 67, [isa], page 41 .
iscomplex (x)
Return true if $x$ is a complex-valued numeric object.
See also: [isreal], page 67, [isnumeric], page 67, [ischar], page 73, [isfloat], page 67, [islogical], page 67, [isstring], page 73 , [isa], page 41 .
ismatrix (x)
Return true if $x$ is a 2-D array.
A matrix is an object with two dimensions (ndims $(x)==2)$ for which size ( $x$ ) returns [M, N] with non-negative M and N .
See also: [isscalar], page 68, [isvector], page 68, [iscell], page 126, [isstruct], page 119, [issparse], page 666, [isa], page 41.
isvector (x)
Return true if $x$ is a vector.
A vector is a 2-D array where one of the dimensions is equal to 1 (either 1 xN or Nx 1 ). As a consequence of this definition, a 1 x 1 array (a scalar) is also a vector.
See also: [isscalar], page 68, [ismatrix], page 68, [iscolumn], page 68, [isrow], page 68, [size], page 47.
isrow (x)
Return true if x is a row vector.
A row vector is a 2-D array for which size ( x ) returns [1, N] with non-negative N .
See also: [iscolumn], page 68, [isscalar], page 68, [isvector], page 68, [ismatrix], page 68, [size], page 47.
iscolumn (x)
Return true if $x$ is a column vector.
A column vector is a 2-D array for which size ( $x$ ) returns [ $\mathrm{N}, 1]$ with non-negative N .
See also: [isrow], page 68, [isscalar], page 68, [isvector], page 68, [ismatrix], page 68, [size], page 47.
isscalar (x)
Return true if $x$ is a scalar.
A scalar is an object with two dimensions for which size ( $x$ ) returns [1, 1].
See also: [isvector], page 68, [ismatrix], page 68, [size], page 47.
issquare (x)
Return true if $x$ is a $2-D$ square array.
A square array is a 2 - D object for which size ( $x$ ) returns [ $\mathrm{N}, \mathrm{N}$ ] where N is a nonnegative integer.

See also: [isscalar], page 68, [isvector], page 68, [ismatrix], page 68, [size], page 47.
issymmetric (A)
issymmetric (A, tol)
issymmetric (A, "skew")
issymmetric (A, "skew", tol)
Return true if $A$ is a symmetric or skew-symmetric matrix within the tolerance specified by $t o l$.
The default tolerance is zero (uses faster code).

The type of symmetry to check may be specified with the additional input "nonskew" (default) for regular symmetry or "skew" for skew-symmetry.
Background: A matrix is symmetric if the transpose of the matrix is equal to the original matrix: $A==A .^{\prime}$. If a tolerance is given then symmetry is determined by $\operatorname{norm}\left(A-A .^{\prime}, \operatorname{Inf}\right) / \operatorname{norm}(A, \operatorname{Inf})<t o l$.
A matrix is skew-symmetric if the transpose of the matrix is equal to the negative of the original matrix: $A==-A .^{\prime}$. If a tolerance is given then skew-symmetry is determined by norm $\left(A+A .^{\prime}\right.$, Inf) / norm ( $A$, Inf) < tol.
See also: [ishermitian], page 69, [isdefinite], page 69.
ishermitian (A)
ishermitian (A, tol)
ishermitian (A, "skew")
ishermitian ( $A$, "skew", tol)
Return true if $A$ is a Hermitian or skew-Hermitian matrix within the tolerance specified by tol.
The default tolerance is zero (uses faster code).
The type of symmetry to check may be specified with the additional input "nonskew" (default) for regular Hermitian or "skew" for skew-Hermitian.
Background: A matrix is Hermitian if the complex conjugate transpose of the matrix is equal to the original matrix: $A==A^{\prime}$. If a tolerance is given then the calculation is norm ( $A-A^{\prime}, \operatorname{Inf}$ ) / norm ( $A$, Inf) < tol.
A matrix is skew-Hermitian if the complex conjugate transpose of the matrix is equal to the negative of the original matrix: $A==-A^{\prime}$. If a tolerance is given then the calculation is norm $\left(A+A^{\prime}\right.$, Inf) / norm ( $A$, Inf) < tol.
See also: [issymmetric], page 68, [isdefinite], page 69 .
isdefinite (A)
isdefinite ( $A$, tol)
Return true if $A$ is symmetric positive definite matrix within the tolerance specified by tol.
If tol is omitted, use a tolerance of $100 *$ eps $*$ norm ( $A$, "fro").
Background: A positive definite matrix has eigenvalues which are all greater than zero. A positive semi-definite matrix has eigenvalues which are all greater than or equal to zero. The matrix $A$ is very likely to be positive semi-definite if the following two conditions hold for a suitably small tolerance tol.

```
isdefinite (A) = 0
isdefinite (A + 5*tol, tol) = 1
```

See also: [issymmetric], page 68, [ishermitian], page 69.
isbanded ( $A$, lower, upper)
Return true if $A$ is a matrix with entries confined between lower diagonals below the main diagonal and upper diagonals above the main diagonal.
lower and upper must be non-negative integers.
See also: [isdiag], page 70, [istril], page 70, [istriu], page 70, [bandwidth], page 586.
isdiag (A)
Return true if $A$ is a diagonal matrix.
See also: [isbanded], page 69, [istril], page 70, [istriu], page 70, [diag], page 525, [bandwidth], page 586.
istril (A)
Return true if $A$ is a lower triangular matrix.
A lower triangular matrix has nonzero entries only on the main diagonal and below.
See also: [istriu], page 70, [isbanded], page 69, [isdiag], page 70, [tril], page 523, [bandwidth], page 586.
istriu (A)
Return true if $A$ is an upper triangular matrix.
An upper triangular matrix has nonzero entries only on the main diagonal and above.
See also: [isdiag], page 70, [isbanded], page 69, [istril], page 70, [triu], page 523, [bandwidth], page 586.
isprime (x)
Return a logical array which is true where the elements of $x$ are prime numbers and false where they are not.
A prime number is conventionally defined as a positive integer greater than 1 (e.g., $2,3, \ldots$ ) which is divisible only by itself and 1 . Octave extends this definition to include both negative integers and complex values. A negative integer is prime if its positive counterpart is prime. This is equivalent to isprime (abs ( x )).
If class (x) is complex, then primality is tested in the domain of Gaussian integers (https://en.wikipedia.org/wiki/Gaussian_integer). Some non-complex integers are prime in the ordinary sense, but not in the domain of Gaussian integers. For example, $5=(1+2 i) *(1-2 i)$ shows that 5 is not prime because it has a factor other than itself and 1. Exercise caution when testing complex and real values together in the same matrix.

Examples:

```
isprime (1:6)
```



```
isprime ([i, 2, 3, 5])
    # 0 0 1 0
```

Programming Note: isprime is suitable for all $x$ in the range $\operatorname{abs}(x)<2^{64}$.
Compatibility Note: matlab does not extend the definition of prime numbers and will produce an error if given negative or complex inputs.
See also: [primes], page 564, [factor], page 563, [gcd], page 563, [lcm], page 563.
If instead of knowing properties of variables, you wish to know which variables are defined and to gather other information about the workspace itself, see Section 7.3 [Status of Variables], page 142.

## 5 Strings

A string constant consists of a sequence of characters enclosed in either double-quote or single-quote marks. For example, both of the following expressions

```
"parrot"
'parrot'
```

represent the string whose contents are 'parrot'. Strings in Octave can be of any length.
Since the single-quote mark is also used for the transpose operator (see Section 8.3 [Arithmetic Ops], page 161) but double-quote marks have no other purpose in Octave, it is best to use double-quote marks to denote strings.

Strings can be concatenated using the notation for defining matrices. For example, the expression
[ "foo" , "bar" , "baz" ]
produces the string whose contents are 'foobarbaz'. See Chapter 4 [Numeric Data Types], page 51 , for more information about creating matrices.

While strings can in principle store arbitrary content, most functions expect them to be UTF-8 encoded Unicode strings.

Furthermore, it is possible to create a string without actually writing a text. The function blanks creates a string of a given length consisting only of blank characters (ASCII code 32).
blanks (n)
Return a string of $n$ blanks.
For example:
blanks (10);
whos ans
$\Rightarrow$

| AttrName Size | Bytes Class <br> $========$ $=====$ <br> ans $1 \times 10$ | 10 | char |
| :---: | :---: | ---: | :--- |

See also: [repmat], page 527 .

### 5.1 Escape Sequences in String Constants

In double-quoted strings, the backslash character is used to introduce escape sequences that represent other characters. For example, ' n ' embeds a newline character in a double-quoted string and ' $\backslash$ "' embeds a double quote character. In single-quoted strings, backslash is not a special character. Here is an example showing the difference:

```
double ("\n")
    => 10
double ('\n')
    =>[ 92 110 ]
```

Here is a table of all the escape sequences used in Octave (within double quoted strings). They are the same as those used in the C programming language.
<br> Represents a literal backslash, ' $\backslash$ '.

```
\" Represents a literal double-quote character, '"'.
\' Represents a literal single-quote character, '''.
\0 Represents the null character, control-@, ASCII code 0.
\a Represents the "alert" character, control-g, ASCII code 7.
\b Represents a backspace, control-h, ASCII code 8.
\f Represents a formfeed, control-l, ASCII code 12.
\n Represents a newline, control-j, ASCII code 10.
\r Represents a carriage return, control-m, ASCII code 13.
\t Represents a horizontal tab, control-i, ASCII code 9.
\v Represents a vertical tab, control-k, ASCII code 11.
\(\backslash n n n \quad\) Represents the octal value nnn, where nnn are one to three digits between 0 and 7. For example, the code for the ASCII ESC (escape) character is ' \(\backslash 033\) '.
\xhh... Represents the hexadecimal value \(h h\), where hh are hexadecimal digits ('0' through ' 9 ' and either ' \(A\) ' through ' \(F\) ' or ' \(a\) ' through ' \(f\) '). Like the same construct in ANSI C, the escape sequence continues until the first non-hexadecimal digit is seen. However, using more than two hexadecimal digits produces undefined results.
```

In a single-quoted string there is only one escape sequence: you may insert a single quote character using two single quote characters in succession. For example,

```
'I can''t escape'
    | I can't escape
```

In scripts the two different string types can be distinguished if necessary by using is_ dq_string and is_sq_string.
is_dq_string (x)
Return true if $x$ is a double-quoted character string.
See also: [is_sq_string], page 72, [ischar], page 73.
is_sq_string (x)
Return true if $x$ is a single-quoted character string.
See also: [is_dq_string], page 72, [ischar], page 73.

### 5.2 Character Arrays

The string representation used by Octave is an array of characters, so internally the string "dddddddddd" is actually a row vector of length 10 containing the value 100 in all places (100 is the ASCII code of "d"). This lends itself to the obvious generalization to character matrices. Using a matrix of characters, it is possible to represent a collection of same-length strings in one variable. The convention used in Octave is that each row in a character matrix is a separate string, but letting each column represent a string is equally possible.

The easiest way to create a character matrix is to put several strings together into a matrix.

```
collection = [ "String #1"; "String #2" ];
```

This creates a 2 -by- 9 character matrix.
The function ischar can be used to test if an object is a character matrix.
ischar (x)
Return true if $x$ is a character array.
See also: [isfloat], page 67, [isinteger], page 58, [islogical], page 67, [isnumeric], page 67, [isstring], page 73, [iscellstr], page 133, [isa], page 41.
isstring (s)
Return true if $s$ is a string array.
A string array is a data type that stores strings (row vectors of characters) at each element in the array. It is distinct from character arrays which are N-dimensional arrays where each element is a single 1 x 1 character. It is also distinct from cell arrays of strings which store strings at each element, but use cell indexing ' $\}$ ' to access elements rather than string arrays which use ordinary array indexing '()'.

Programming Note: Octave does not yet implement string arrays so this function will always return false.

See also: [ischar], page 73, [iscellstr], page 133, [isfloat], page 67, [isinteger], page 58, [islogical], page 67, [isnumeric], page 67, [isa], page 41.

To test if an object is a string (i.e., a 1 xN row vector of characters and not a character matrix) you can use the ischar function in combination with the isrow function as in the following example:

```
ischar (collection)
    => 1
ischar (collection) && isrow (collection)
    => 0
ischar ("my string") && isrow ("my string")
    # 1
```

One relevant question is, what happens when a character matrix is created from strings of different length. The answer is that Octave puts blank characters at the end of strings shorter than the longest string. It is possible to use a different character than the blank character using the string_fill_char function.

```
val = string_fill_char ()
old_val = string_fill_char (new_val)
string_fill_char (new_val, "local")
```

Query or set the internal variable used to pad all rows of a character matrix to the same length.

The value must be a single character and the default is " " (a single space). For example:

```
string_fill_char ("X");
[ "these"; "are"; "strings" ]
    => "theseXX"
    "areXXXX"
            "strings"
```

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

Another useful function to control the text justification in this case is the strjust function.

```
strjust (s)
strjust (s, pos)
```

Return the text, $s$, justified according to pos, which may be "left", "center", or "right".
If pos is omitted it defaults to "right".
Null characters are replaced by spaces. All other character data are treated as nonwhite space.
Example:

```
strjust (["a"; "ab"; "abc"; "abcd"])
    #
        " a"
        " ab"
        " abc"
        "abcd"
```

See also: [deblank], page 75, [strrep], page 90, [strtrim], page 75 , [untabify], page 76.
This shows a problem with character matrices. It simply isn't possible to represent strings of different lengths. The solution is to use a cell array of strings, which is described in Section 6.3.4 [Cell Arrays of Strings], page 133.

### 5.3 String Operations

Octave supports a wide range of functions for manipulating strings. Since a string is just a matrix, simple manipulations can be accomplished using standard operators. The following example shows how to replace all blank characters with underscores.

```
quote = ...
    "First things first, but not necessarily in that order";
quote( quote == " " ) = "_"
=> quote =
        First_things_first,_but_not_necessarily_in_that_order
```

For more complex manipulations, such as searching, replacing, and general regular expressions, the following functions come with Octave.

### 5.3.1 Common String Operations

The following functions are useful to perform common String operations.
tolower (s)
lower (s)
Return a copy of the string or cell string $s$, with each uppercase character replaced by the corresponding lowercase one; non-alphabetic characters are left unchanged.
For example:

```
tolower ("MiXeD cAsE 123")
                    => "mixed case 123"
```

See also: [toupper], page 75 .
toupper (s)
upper (s)
Return a copy of the string or cell string $s$, with each lowercase character replaced by the corresponding uppercase one; non-alphabetic characters are left unchanged.
For example:

```
toupper ("MiXeD cAsE 123")
            => "MIXED CASE 123"
```

See also: [tolower], page 75 .
deblank (s)
Remove trailing whitespace and nulls from $s$.
If $s$ is a matrix, deblank trims each row to the length of longest string. If $s$ is a cell array of strings, operate recursively on each string element.
Examples:

```
deblank (" abc ")
    => " abc"
deblank ([" abc "; " def "])
            => [" abc " ; " def"]
```

See also: [strtrim], page 75 .
strtrim (s)
Remove leading and trailing whitespace from $s$.
If $s$ is a matrix, strtrim trims each row to the length of longest string. If $s$ is a cell array of strings, operate recursively on each string element.
For example:

```
strtrim (" abc ")
        => "abc"
strtrim ([" abc "; " def "])
    => ["abc " ; " def"]
```

See also: [deblank], page 75 .
strtrunc (s, n)
Truncate the character string $s$ to length $n$.
If $s$ is a character matrix, then the number of columns is adjusted.
If $s$ is a cell array of strings, then the operation is performed on each cell element and the new cell array is returned.
untabify ( $t$ )
untabify ( $t, t_{w}$ )
untabify ( $t, t w$, deblank)
Replace TAB characters in $t$ with spaces.
The input, $t$, may be either a 2-D character array, or a cell array of character strings. The output is the same class as the input.
The tab width is specified by $t w$, and defaults to eight.
If the optional argument deblank is true, then the spaces will be removed from the end of the character data.
The following example reads a file and writes an untabified version of the same file with trailing spaces stripped.

```
fid = fopen ("tabbed_script.m");
text = char (fread (fid, "uchar")');
fclose (fid);
fid = fopen ("untabified_script.m", "w");
text = untabify (strsplit (text, "\n"), 8, true);
fprintf (fid, "%s\n", text{:});
fclose (fid);
```

See also: [strjust], page 74, [strsplit], page 81, [deblank], page 75.
do_string_escapes (string)
Convert escape sequences in string to the characters they represent.
Escape sequences begin with a leading backslash (' $\backslash$ ') followed by $1-3$ characters (.e.g., " $\backslash \mathrm{n} " ~=>~ n e w l i n e) . ~$

See also: [undo_string_escapes], page 76.
undo_string_escapes (s)
Convert special characters in strings back to their escaped forms.
For example, the expression

```
bell = "\a";
```

assigns the value of the alert character (control-g, ASCII code 7) to the string variable bell. If this string is printed, the system will ring the terminal bell (if it is possible). This is normally the desired outcome. However, sometimes it is useful to be able to print the original representation of the string, with the special characters replaced by their escape sequences. For example,

```
octave:13> undo_string_escapes (bell)
ans = \a
```

replaces the unprintable alert character with its printable representation.
See also: [do_string_escapes], page 76.

### 5.3.2 Concatenating Strings

Strings can be concatenated using matrix notation (see Chapter 5 [Strings], page 71, Section 5.2 [Character Arrays], page 72) which is often the most natural method. For example:

```
fullname = [fname ".txt"];
email = ["<" user "@" domain ">"];
```

In each case it is easy to see what the final string will look like. This method is also the most efficient. When using matrix concatenation the parser immediately begins joining the strings without having to process the overhead of a function call and the input validation of the associated function.

The newline function can be used to join strings such that they appear as multiple lines of text when displayed.
newline
Return the character corresponding to a newline.
This is equivalent to " $\backslash \mathrm{n}$ ".
Example Code

```
joined_string = [newline "line1" newline "line2"]
#
line1
line2
```

See also: [strcat], page 79 , [strjoin], page 84 , [strsplit], page 81.
In addition, there are several other functions for concatenating string objects which can be useful in specific circumstances: char, strvcat, strcat, and cstrcat. Finally, the general purpose concatenation functions can be used: see [cat], page 517, [horzcat], page 517, and [vertcat], page 517.

- All string concatenation functions except cstrcat convert numerical input into character data by taking the corresponding UTF-8 character for each element (or multi-byte sequence), as in the following example:

```
char ([98, 97, 110, 97, 110, 97])
    b banana
```

For conversion between locale encodings and UTF-8, see [unicode2native], page 94, and [native2unicode], page 95.

- char and strvcat concatenate vertically, while strcat and cstrcat concatenate horizontally. For example:

```
char ("an apple", "two pears")
        a an apple
            two pears
strcat ("oc", "tave", " is", " good", " for you")
    => octave is good for you
```

- char generates an empty row in the output for each empty string in the input. strvcat, on the other hand, eliminates empty strings.

```
char ("orange", "green", "", "red")
    => orange
        green
        red
strvcat ("orange", "green", "", "red")
    => orange
        green
        red
```

- All string concatenation functions except cstrcat also accept cell array data (see Section 6.3 [Cell Arrays], page 124). char and strvcat convert cell arrays into character arrays, while strcat concatenates within the cells of the cell arrays:

```
char ({"red", "green", "", "blue"})
        => red
            green
            blue
strcat ({"abc"; "ghi"}, {"def"; "jkl"})
        #
            {
            [1,1] = abcdef
            [2,1] = ghijkl
            }
```

- strcat removes trailing white space in the arguments (except within cell arrays), while cstrcat leaves white space untouched. Both kinds of behavior can be useful as can be seen in the examples:

```
strcat (["dir1";"directory2"], ["/";"/"], ["file1";"file2"])
    => dir1/file1
        directory2/file2
cstrcat (["thirteen apples"; "a banana"], [" 5$";" 1$"])
    => thirteen apples 5$
    a banana 1$
```

Note that in the above example for cstrcat, the white space originates from the internal representation of the strings in a string array (see Section 5.2 [Character Arrays], page 72 ).
char (x)
char ( $x, . .$. )
char (s1, s2, ...)
char (cell_array)
Create a string array from one or more numeric matrices, character matrices, or cell arrays.

Arguments are concatenated vertically. The returned values are padded with blanks as needed to make each row of the string array have the same length. Empty input strings are significant and will concatenated in the output.
For numerical input, each element is converted to the corresponding ASCII character. A range error results if an input is outside the ASCII range (0-255).
For cell arrays, each element is concatenated separately. Cell arrays converted through char can mostly be converted back with cellstr. For example:

```
char ([97, 98, 99], "", {"98", "99", 100}, "str1", ["ha", "lf"])
    => ["abc "
        | I
        "98 "
        "99 "
        "d "
        "str1"
        "half"]
```

See also: [strvcat], page 79, [cellstr], page 133.
strvcat (x)
strvcat (x,...)
strvcat (s1, s2,...)
strvcat (cell_array)
Create a character array from one or more numeric matrices, character matrices, or cell arrays.
Arguments are concatenated vertically. The returned values are padded with blanks as needed to make each row of the string array have the same length. Unlike char, empty strings are removed and will not appear in the output.
For numerical input, each element is converted to the corresponding ASCII character. A range error results if an input is outside the ASCII range (0-255).
For cell arrays, each element is concatenated separately. Cell arrays converted through strvcat can mostly be converted back with cellstr. For example:

```
strvcat ([97, 98, 99], "", {"98", "99", 100}, "str1", ["ha", "lf"])
    # ["abc "
            "98 "
            "99 "
            "d "
            "str1"
            "half"]
```

See also: [char], page 78, [strcat], page 79, [cstrcat], page 80.

```
strcat (s1, s2,...)
```

Return a string containing all the arguments concatenated horizontally.
If the arguments are cell strings, strcat returns a cell string with the individual cells concatenated. For numerical input, each element is converted to the corresponding ASCII character. Trailing white space for any character string input is eliminated before the strings are concatenated. Note that cell string values do not have whitespace trimmed.

For example:

```
strcat ("|", " leading space is preserved", "|")
    | | leading space is preserved|
strcat ("|", "trailing space is eliminated ", "|")
    | |trailing space is eliminated|
strcat ("homogeneous space |", " ", "| is also eliminated")
    => homogeneous space || is also eliminated
s = [ "ab"; "cde" ];
strcat (s, s, s)
    #
        "ababab "
        "cdecdecde"
s = { "ab"; "cd " };
strcat (s, s, s)
        #
            {
            [1,1] = ababab
            [2,1] = cd cd cd
            }
```

See also: [cstrcat], page 80, [char], page 78, [strvcat], page 79.
cstrcat (s1, s2,...)
Return a string containing all the arguments concatenated horizontally with trailing white space preserved.
For example:

$$
\begin{gathered}
\text { cstrcat ("ab ", "cd") } \\
\Rightarrow \text { "ab cd" } \\
\mathrm{s}=[\text { "ab"; "cde" ]; } \\
\text { cstrcat (s, s, s) } \\
\Rightarrow \text { "ab ab ab" } \\
\text { "cdecdecde" }
\end{gathered}
$$

See also: [strcat], page 79, [char], page 78, [strvcat], page 79 .

### 5.3.3 Splitting and Joining Strings

substr (s, offset)
substr (s, offset, len)
Return the substring of $s$ which starts at character number offset and is len characters long.
Position numbering for offsets begins with 1. If offset is negative, extraction starts that far from the end of the string.
If len is omitted, the substring extends to the end of $s$. A negative value for len extracts to within len characters of the end of the string

Examples:

```
substr ("This is a test string", 6, 9)
    => "is a test"
substr ("This is a test string", -11)
    => "test string"
substr ("This is a test string", -11, -7)
        # "test"
```

This function is patterned after the equivalent function in Perl.
[tok, rem] = strtok (str)
[tok, rem] = strtok (str, delim)
Find all characters in the string str up to, but not including, the first character which is in the string delim.
str may also be a cell array of strings in which case the function executes on every individual string and returns a cell array of tokens and remainders.
Leading delimiters are ignored. If delim is not specified, whitespace is assumed.
If rem is requested, it contains the remainder of the string, starting at the first delimiter.

Examples:

```
strtok ("this is the life")
            => "this"
[tok, rem] = strtok ("14*27+31", "+-*/")
            =>
                tok = 14
                rem = *27+31
```

See also: [index], page 87, [strsplit], page 81, [strchr], page 87, [isspace], page 109.
[cstr] = strsplit (str)
[cstr] = strsplit (str, del)
[cstr] = strsplit (..., name, value)
[cstr, matches] = strsplit (...)
Split the string str using the delimiters specified by del and return a cell string array of substrings.
If a delimiter is not specified the string is split at whitespace $\{"$ ", " ff ", "\n", "\} \mathrm { r } " , "\t", "\v"\}. Otherwise, the delimiter, del must be a string or cell array of strings. By default, consecutive delimiters in the input string $s$ are collapsed into one resulting in a single split.

Supported name/value pair arguments are:

- collapsedelimiters which may take the value of true (default) or false.
- delimitertype which may take the value of "simple" (default) or "regularexpression". A simple delimiter matches the text exactly as written. Otherwise, the syntax for regular expressions outlined in regexp is used.

The optional second output, matches, returns the delimiters which were matched in the original string.
Examples with simple delimiters:

```
strsplit ("a b c")
    \(\Rightarrow\)
        \{
        \([1,1]=a\)
        \([1,2]=\mathrm{b}\)
        \([1,3]=c\)
        \}
strsplit ("a,b,c", ",")
        \(\Rightarrow\)
            \{
                \([1,1]=a\)
                \([1,2]=\mathrm{b}\)
                \([1,3]=c\)
            \}
strsplit ("a foo b,bar c", \{" ", ",", "foo", "bar"\})
    \(\Rightarrow\)
        \{
            \([1,1]=a\)
            \([1,2]=\mathrm{b}\)
            \([1,3]=c\)
            \}
strsplit ("a,,b, c", \{",", " "\}, "collapsedelimiters", false)
    \(\Rightarrow\)
        \{
            \([1,1]=a\)
        \([1,2]=\)
        \([1,3]=\mathrm{b}\)
        \([1,4]=\)
        \([1,5]=c\)
        \}
```

Examples with regularexpression delimiters:

```
strsplit ("a foo b,bar c", ',\\s|foolbar', ...
    "delimitertype", "regularexpression")
=>
{
        [1,1] = a
        [1,2] = b
        [1,3] = c
}
strsplit ("a,,b, c", '[, ]', "collapsedelimiters", false, ...
```

```
    "delimitertype", "regularexpression")
{
    [1,1] = a
    [1,2] =
    [1,3] = b
    [1,4] =
    [1,5] = c
}
strsplit ("a,\t,b, c", {',', '\s'}, "delimitertype", "regularexpression")
#
{
    [1,1] = a
    [1,2] = b
    [1,3] = c
}
strsplit ("a,\t,b, c", {',', ' ', '\t'}, "collapsedelimiters", false)
#
{
    [1,1] = a
    [1,2] =
    [1,3] =
    [1,4] = b
    [1,5] =
    [1,6] = c
}
```

See also: [ostrsplit], page 83, [strjoin], page 84, [strtok], page 81, [regexp], page 91.
[cstr] = ostrsplit (s, sep)
[cstr] = ostrsplit (s, sep, strip_empty)
Split the string $s$ using one or more separators sep and return a cell array of strings.
Consecutive separators and separators at boundaries result in empty strings, unless strip_empty is true. The default value of strip_empty is false.
2-D character arrays are split at separators and at the original column boundaries.
Example:

```
ostrsplit ("a,b,c", ",")
    \(\Rightarrow\)
            \{
            \([1,1]=a\)
                        \([1,2]=b\)
            \([1,3]=c\)
            \}
ostrsplit (["a,b" ; "cde"], ",")
        \(\Rightarrow\)
            \{
            \([1,1]=a\)
            \([1,2]=\mathrm{b}\)
            \([1,3]=\) cde
            \}
```

See also: [strsplit], page 81, [strtok], page 81.

```
str = strjoin (cstr)
str = strjoin (cstr, delimiter)
```

Join the elements of the cell string array, cstr, into a single string.
If no delimiter is specified, the elements of cstr are separated by a space.
If delimiter is specified as a string, the cell string array is joined using the string. Escape sequences are supported.
If delimiter is a cell string array whose length is one less than cstr, then the elements of cstr are joined by interleaving the cell string elements of delimiter. Escape sequences are not supported.

```
strjoin ({'Octave','Scilab','Lush','Yorick'}, '*')
    => 'Octave*Scilab*Lush*Yorick'
```

See also: [strsplit], page 81.

### 5.3.4 Searching in Strings

Since a string is a character array, comparisons between strings work element by element as the following example shows:

```
GNU = "GNU's Not UNIX";
spaces = (GNU == " ")
    # spaces =
        0
```

To determine if two strings are identical it is necessary to use the strcmp function. It compares complete strings and is case sensitive. strncmp compares only the first N characters (with N given as a parameter). strcmpi and strncmpi are the corresponding functions for case-insensitive comparison.
strcmp (s1, s2)
Return 1 if the character strings $s 1$ and $s 2$ are the same, and 0 otherwise.
If either $s 1$ or $s 2$ is a cell array of strings, then an array of the same size is returned, containing the values described above for every member of the cell array. The other argument may also be a cell array of strings (of the same size or with only one element), char matrix or character string.
Caution: For compatibility with matlab, Octave's strcmp function returns 1 if the character strings are equal, and 0 otherwise. This is just the opposite of the corresponding C library function.
See also: [strcmpi], page 85, [strncmp], page 84, [strncmpi], page 85.
strncmp (s1, s2, n)
Return 1 if the first $n$ characters of strings s1 and $s 2$ are the same, and 0 otherwise.

```
strncmp ("abce", "abcd", 3)
                    # 1
```

If either $s 1$ or $s 2$ is a cell array of strings, then an array of the same size is returned, containing the values described above for every member of the cell array. The other
argument may also be a cell array of strings (of the same size or with only one element), char matrix or character string.

```
strncmp ("abce", {"abcd", "bca", "abc"}, 3)
    # [1, 0, 1]
```

Caution: For compatibility with matlab, Octave's strnemp function returns 1 if the character strings are equal, and 0 otherwise. This is just the opposite of the corresponding C library function.
See also: [strncmpi], page 85, [strcmp], page 84, [strcmpi], page 85.
strcmpi (s1, s2)
Return 1 if the character strings $s 1$ and $s 2$ are the same, disregarding case of alphabetic characters, and 0 otherwise.
If either $s 1$ or $s 2$ is a cell array of strings, then an array of the same size is returned, containing the values described above for every member of the cell array. The other argument may also be a cell array of strings (of the same size or with only one element), char matrix or character string.
Caution: For compatibility with matlab, Octave's strcmp function returns 1 if the character strings are equal, and 0 otherwise. This is just the opposite of the corresponding C library function.
Caution: National alphabets are not supported.
See also: [strcmp], page 84, [strncmp], page 84, [strncmpi], page 85.
strncmpi (s1, s2, n)
Return 1 if the first $n$ character of $s 1$ and $s 2$ are the same, disregarding case of alphabetic characters, and 0 otherwise.

If either $s 1$ or $s 2$ is a cell array of strings, then an array of the same size is returned, containing the values described above for every member of the cell array. The other argument may also be a cell array of strings (of the same size or with only one element), char matrix or character string.
Caution: For compatibility with matlab, Octave's strncmpi function returns 1 if the character strings are equal, and 0 otherwise. This is just the opposite of the corresponding C library function.
Caution: National alphabets are not supported.
See also: [strncmp], page 84, [strcmp], page 84, [strcmpi], page 85.
Despite those comparison functions, there are more specialized function to find the index position of a search pattern within a string.
retval $=$ startsWith (str, pattern)
retval $=$ startsWith (str, pattern, "IgnoreCase", ignore_case)
Check whether string(s) start with pattern(s).
Return an array of logical values that indicates which string(s) in the input str (a single string or cell array of strings) begin with the input pattern (a single string or cell array of strings).

If the value of the parameter "IgnoreCase" is true, then the function will ignore the letter case of str and pattern. By default, the comparison is case sensitive.
Examples:

```
## one string and one pattern while considering case
startsWith ("hello", "he")
    # 1
## one string and one pattern while ignoring case
startsWith ("hello", "HE", "IgnoreCase", true)
    # 1
## multiple strings and multiple patterns while considering case
startsWith ({"lab work.pptx", "data.txt", "foundations.ppt"},
    {"lab", "data"})
    # 1 1 0
## multiple strings and one pattern while considering case
startsWith ({"DATASHEET.ods", "data.txt", "foundations.ppt"},
        "data", "IgnoreCase", false)
    m 0 1 0
## multiple strings and one pattern while ignoring case
startsWith ({"DATASHEET.ods", "data.txt", "foundations.ppt"},
        "data", "IgnoreCase", true)
    # 1 1 0
```

See also: [endsWith], page 86, [regexp], page 91, [strncmp], page 84, [strncmpi], page 85.
retval = endsWith (str, pattern)
retval = endsWith (str, pattern, "IgnoreCase", ignore_case)
Check whether string(s) end with pattern(s).
Return an array of logical values that indicates which string(s) in the input str (a single string or cell array of strings) end with the input pattern (a single string or cell array of strings).
If the value of the parameter "IgnoreCase" is true, then the function will ignore the letter case of str and pattern. By default, the comparison is case sensitive.
Examples:

```
## one string and one pattern while considering case
endsWith ("hello", "lo")
    # 1
## one string and one pattern while ignoring case
endsWith ("hello", "LO", "IgnoreCase", true)
    # 1
```

```
## multiple strings and multiple patterns while considering case
endsWith ({"tests.txt", "mydoc.odt", "myFunc.m", "results.pptx"},
        {".docx", ".odt", ".txt"})
    # 1 1 0 0
## multiple strings and one pattern while considering case
endsWith ({"TESTS.TXT", "mydoc.odt", "result.txt", "myFunc.m"},
        ".txt", "IgnoreCase", false)
    m}00001
## multiple strings and one pattern while ignoring case
endsWith ({"TESTS.TXT", "mydoc.odt", "result.txt", "myFunc.m"},
        ".txt", "IgnoreCase", true)
    m}10001
```

See also: [startsWith], page 85, [regexp], page 91, [strncmp], page 84, [strncmpi], page 85 .
findstr ( $s, t$ )
findstr (s, t, overlap)
This function is obsolete. Use strfind instead.
Return the vector of all positions in the longer of the two strings $s$ and $t$ where an occurrence of the shorter of the two starts.
If the optional argument overlap is true (default), the returned vector can include overlapping positions. For example:

```
findstr ("ababab", "a")
    \(\Rightarrow \quad[1,3,5]\);
findstr ("abababa", "aba", 0)
    \(\Rightarrow[1,5]\)
```

Caution: findstr is obsolete. Use strfind in all new code.
See also: [strfind], page 88, [strmatch], page 89, [strcmp], page 84, [strncmp], page 84, [strcmpi], page 85, [strncmpi], page 85, [find], page 513.

```
idx = strchr (str, chars)
```

idx = strchr (str, chars, n)
idx = strchr (str, chars, n, direction)
[i, j] = strchr (...)

Search through the string str for occurrences of characters from the set chars.
The return value(s), as well as the $n$ and direction arguments behave identically as in find.
This will be faster than using regexp in most cases.
See also: [find], page 513.
index ( $s, t$ )
index ( $s$, $t$, direction)
Return the position of the first occurrence of the string $t$ in the string $s$, or 0 if no occurrence is found.
$s$ may also be a string array or cell array of strings.
For example:

```
index ("Teststring", "t")
    =>4
```

If direction is "first", return the first element found. If direction is "last", return the last element found.
See also: [find], page 513, [rindex], page 88.
rindex ( $s, t$ )
Return the position of the last occurrence of the character string $t$ in the character string $s$, or 0 if no occurrence is found.
$s$ may also be a string array or cell array of strings.
For example:

```
rindex ("Teststring", "t")
    =>6
```

The rindex function is equivalent to index with direction set to "last".
See also: [find], page 513, [index], page 87.

```
idx = unicode_idx (str)
```

Return an array with the indices for each UTF-8 encoded character in str.

```
unicode_idx ("aäbc")
                # [1, 2, 2, 3, 4]
```

idx $=$ strfind (str, pattern)
idx $=$ strfind (cellstr, pattern)
idx $=$ strfind (..., "overlaps", val)
idx $=$ strfind (..., "forcecelloutput", val)

Search for pattern in the string str and return the starting index of every such occurrence in the vector idx.
If there is no such occurrence, or if pattern is longer than str, or if pattern itself is empty, then idx is the empty array [].
The optional argument "overlaps" determines whether the pattern can match at every position in str (true), or only for unique occurrences of the complete pattern (false). The default is true.
If a cell array of strings cellstr is specified then idx is a cell array of vectors, as specified above.
The optional argument "forcecelloutput" forces idx to be returned as a cell array of vectors. The default is false.
Examples:

```
strfind ("abababa", "aba")
    # [1, 3, 5]
strfind ("abababa", "aba", "overlaps", false)
        # [1, 5]
```

```
strfind ({"abababa", "bebebe", "ab"}, "aba")
    =>
            {
                [1,1] =
                    1 5
                [1,2] = [] (1x0)
                [1,3] = [] (1x0)
            }
strfind ("abababa", "aba", "forcecelloutput", true)
    =>
            {
                [1,1] =
                    1 3 5
            }
```

See also: [regexp], page 91, [regexpi], page 93, [find], page 513.
strmatch (s, A)
strmatch (s, A, "exact")
This function is obsolete. Use an alternative such as strncmp or strcmp instead.
Return indices of entries of $A$ which begin with the string $s$.
The second argument $A$ must be a string, character matrix, or a cell array of strings.
If the third argument "exact" is not given, then $s$ only needs to match $A$ up to the length of $s$. Trailing spaces and nulls in $s$ and $A$ are ignored when matching.

For example:

```
strmatch ("apple", "apple juice")
    # 1
strmatch ("apple", ["apple "; "apple juice"; "an apple"])
    # [1; 2]
strmatch ("apple", ["apple "; "apple juice"; "an apple"], "exact")
    # [1]
```

Caution: strmatch is obsolete (and can produce incorrect results in MATLAB when used with cell arrays of strings. Use strncmp (normal case) or strcmp ("exact" case) in all new code. Other replacement possibilities, depending on application, include regexp or validatestring.

See also: [strncmp], page 84, [strcmp], page 84, [regexp], page 91, [strfind], page 88, [validatestring], page 203.

### 5.3.5 Searching and Replacing in Strings

```
newstr \(=\) strrep (str, ptn, rep)
```

newstr $=$ strrep (cellstr, ptn, rep)
newstr $=$ strrep (..., "overlaps", val)

Replace all occurrences of the pattern ptn in the string str with the string rep and return the result.

The optional argument "overlaps" determines whether the pattern can match at every position in str (true), or only for unique occurrences of the complete pattern (false). The default is true.
$s$ may also be a cell array of strings, in which case the replacement is done for each element and a cell array is returned.
Example:

```
strrep ("This is a test string", "is", "&%$")
    => "Th&%$ &%$ a test string"
```

See also: [regexprep], page 93, [strfind], page 88.

```
newstr = erase (str, ptn)
```

Delete all occurrences of ptn within str.
str and ptn can be ordinary strings, cell array of strings, or character arrays.
Examples

```
## string, single pattern
erase ("Hello World!", " World")
    => "Hello!"
## cellstr, single pattern
erase ({"Hello", "World!"}, "World")
    => {"Hello", "!"}
## string, multiple patterns
erase ("The Octave interpreter is fabulous", ...
        {"interpreter ", "The "})
    => "Octave is fabulous"
## cellstr, multiple patterns
erase ({"The ", "Octave interpreter ", "is fabulous"}, ...
        {"interpreter ", "The "})
        => {"", "Octave ", "is fabulous"}
```

Programming Note: erase deletes the first instance of a pattern in a string when there are overlapping occurrences. For example:

```
erase ("abababa", "aba")
    => "b"
```

For processing overlaps, see [strrep], page 90.
See also: [strrep], page 90, [regexprep], page 93.
[s, e, te, m, t, nm, sp] = regexp (str, pat)
[...] = regexp (str, pat, "opt1",...)
Regular expression string matching.
Search for pat in UTF-8 encoded str and return the positions and substrings of any matches, or empty values if there are none.
The matched pattern pat can include any of the standard regex operators, including:
Match any character

*     + ? \{\} Repetition operators, representing
* Match zero or more times
$+\quad$ Match one or more times
? Match zero or one times
\{n\} Match exactly $n$ times
$\{n$,$\} \quad Match n$ or more times
$\{m, n\} \quad$ Match between $m$ and $n$ times
[...] [^...]
List operators. The pattern will match any character listed between " [" and "] ". If the first character is "~" then the pattern is inverted and any character except those listed between brackets will match.
Escape sequences defined below can also be used inside list operators. For example, a template for a floating point number might be $[-+. \backslash \mathrm{d}]+$.
() (?:) Grouping operator. The first form, parentheses only, also creates a token.
$1 \quad$ Alternation operator. Match one of a choice of regular expressions. The alternatives must be delimited by the grouping operator () above.
- \$ Anchoring operators. Requires pattern to occur at the start ( ${ }^{\wedge}$ ) or end (\$) of the string.

In addition, the following escaped characters have special meaning.
\d Match any digit
\D Match any non-digit
\s Match any whitespace character
\S Match any non-whitespace character
\w Match any word character
\W Match any non-word character
\< Match the beginning of a word
\> Match the end of a word
$\backslash$ B Match within a word

Implementation Note: For compatibility with matlab, escape sequences in pat (e.g., " n " $=>$ newline) are expanded even when pat has been defined with single quotes. To disable expansion use a second backslash before the escape sequence (e.g., "<br>n") or use the regexptranslate function.
The outputs of regexp default to the order given below
$s \quad$ The start indices of each matching substring
$e \quad$ The end indices of each matching substring
te The extents of each matched token surrounded by (...) in pat
$m \quad$ A cell array of the text of each match
$t \quad$ A cell array of the text of each token matched
$n m \quad$ A structure containing the text of each matched named token, with the name being used as the fieldname. A named token is denoted by (?<name>...).
$s p \quad$ A cell array of the text not returned by match, i.e., what remains if you split the string based on pat.

Particular output arguments, or the order of the output arguments, can be selected by additional opt arguments. These are strings and the correspondence between the output arguments and the optional argument are

| 'start' | $s$ |
| :--- | :--- |
| 'end' | $e$ |
| 'tokenExtents' | $t e$ |
| 'match' | $m$ |
| 'tokens' | $t$ |
| 'names' | $n m$ |
| 'split' | $s p$ |

Additional arguments are summarized below.
'once' Return only the first occurrence of the pattern.
'matchcase'
Make the matching case sensitive. (default)
Alternatively, use (?-i) in the pattern.
'ignorecase'
Ignore case when matching the pattern to the string.
Alternatively, use (?i) in the pattern.
'stringanchors'
Match the anchor characters at the beginning and end of the string. (default)
Alternatively, use (?-m) in the pattern.
'lineanchors'
Match the anchor characters at the beginning and end of the line.
Alternatively, use (?m) in the pattern.
'dotall' The pattern . matches all characters including the newline character. (default)
Alternatively, use (?s) in the pattern.
'dotexceptnewline'
The pattern . matches all characters except the newline character.
Alternatively, use (?-s) in the pattern.
'literalspacing'
All characters in the pattern, including whitespace, are significant and are used in pattern matching. (default)
Alternatively, use (?-x) in the pattern.
'freespacing'
The pattern may include arbitrary whitespace and also comments beginning with the character ' $\#$ '.
Alternatively, use (?x) in the pattern.
'noemptymatch'
Zero-length matches are not returned. (default)
'emptymatch'
Return zero-length matches.
regexp ('a', 'b*', 'emptymatch') returns [1 2] because there are zero or more ' b ' characters at positions 1 and end-of-string.

Stack Limitation Note: Pattern searches are done with a recursive function which can overflow the program stack when there are a high number of matches. For example,

```
regexp (repmat ('a', 1, 1e5), '(a)+')
```

may lead to a segfault. As an alternative, consider constructing pattern searches that reduce the number of matches (e.g., by creatively using set complement), and then further processing the return variables (now reduced in size) with successive regexp searches.
See also: [regexpi], page 93, [strfind], page 88, [regexprep], page 93.
[s, e, te, m, t, nm, sp] = regexpi (str, pat) [...] = regexpi (str, pat, "opt1", ...)

Case insensitive regular expression string matching.
Search for pat in UTF-8 encoded str and return the positions and substrings of any matches, or empty values if there are none. See [regexp], page 91, for details on the syntax of the search pattern.
See also: [regexp], page 91.

```
outstr = regexprep (string, pat, repstr)
```

outstr = regexprep (string, pat, repstr, "opt1", ...)

Replace occurrences of pattern pat in string with repstr.
The pattern is a regular expression as documented for regexp. See [regexp], page 91. All strings must be UTF-8 encoded.

The replacement string may contain $\$$ i, which substitutes for the ith set of parentheses in the match string. For example,
regexprep ("Bill Dunn", '(\w+) (\w+)', '\$2, \$1')
returns "Dunn, Bill"
Options in addition to those of regexp are
'once' Replace only the first occurrence of pat in the result.
'warnings'
This option is present for compatibility but is ignored.
Implementation Note: For compatibility with matlab, escape sequences in pat (e.g., " n " $=>$ newline) are expanded even when pat has been defined with single quotes. To disable expansion use a second backslash before the escape sequence (e.g., "<br>n") or use the regexptranslate function.
See also: [regexp], page 91, [regexpi], page 93, [strrep], page 90.
regexptranslate (op, s)
Translate a string for use in a regular expression.
This may include either wildcard replacement or special character escaping.
The behavior is controlled by op which can take the following values
"wildcard"
The wildcard characters ., *, and ? are replaced with wildcards that are appropriate for a regular expression. For example:

```
regexptranslate ("wildcard", "*.m")
    \(\Rightarrow\) '.*\.m'
```

"escape" The characters \$.? [], that have special meaning for regular expressions are escaped so that they are treated literally. For example:

```
regexptranslate ("escape", "12.5")
    \(\Rightarrow\) '12\.5'
```

See also: [regexp], page 91, [regexpi], page 93, [regexprep], page 93.

### 5.4 Converting Strings

Octave offers several kinds of conversion functions for Strings.

### 5.4.1 String encoding

```
native_bytes = unicode2native (utf8_str, codepage)
native_bytes = unicode2native (utf8_str)
Convert UTF-8 string utf8_str to byte stream using codepage.
The character vector utf8_str is converted to a byte stream native_bytes using the code page given by codepage. The string codepage must be an identifier of a valid code page. Examples for valid code pages are "ISO-8859-1", "Shift-JIS", or "UTF-16". For a list of supported code pages, see https://www.gnu.org/software/libiconv. If codepage is omitted or empty, the system default codepage is used.
```

If any of the characters cannot be mapped into the codepage codepage, they are replaced with the appropriate substitution sequence for that codepage.

See also: [native2unicode], page 95.
utf8_str = native2unicode (native_bytes, codepage)
utf8_str = native2unicode (native_bytes)
Convert byte stream native_bytes to UTF-8 using codepage.
The numbers in the vector native_bytes are rounded and clipped to integers between 0 and 255. This byte stream is then mapped into the code page given by the string codepage and returned in the string utf8_str. Octave uses UTF-8 as its internal encoding. The string codepage must be an identifier of a valid code page. Examples for valid code pages are "ISO-8859-1", "Shift-JIS", or "UTF-16". For a list of supported code pages, see https://www.gnu.org/software/libiconv. If codepage is omitted or empty, the system default codepage is used.

If native_bytes is a string vector, it is returned as is.
See also: [unicode2native], page 94.

### 5.4.2 Numerical Data and Strings

Apart from the string concatenation functions (see Section 5.3.2 [Concatenating Strings], page 77) which cast numerical data to the corresponding UTF-8 encoded characters, there are several functions that format numerical data as strings. mat2str and num2str convert real or complex matrices, while int2str converts integer matrices. int2str takes the real part of complex values and round fractional values to integer. A more flexible way to format numerical data as strings is the sprintf function (see Section 14.2.4 [Formatted Output], page 298, [sprintf], page 299).

```
s = mat2str (x, n)
s = mat2str (x, n, "class")
Format real, complex, and logical matrices as strings.
```

The returned string may be used to reconstruct the original matrix by using the eval function.

The precision of the values is given by $n$. If $n$ is a scalar then both real and imaginary parts of the matrix are printed to the same precision. Otherwise $n(1)$ defines the precision of the real part and $n(2)$ defines the precision of the imaginary part. The default for $n$ is 15 .

If the argument "class" is given then the class of $x$ is included in the string in such a way that eval will result in the construction of a matrix of the same class.

```
mat2str ([ -1/3 + i/7; 1/3 - i/7 ], [4 2])
    => "[-0.3333+0.14i;0.3333-0.14i]"
mat2str ([ -1/3 +i/7; 1/3 -i/7 ], [4 2])
        => "[-0.3333+0i 0+0.14i;0.3333+0i -0-0.14i]"
mat2str (int16 ([1 -1]), "class")
    => "int16([1 -1])"
mat2str (logical (eye (2)))
    => "[true false;false true]"
isequal (x, eval (mat2str (x)))
    # 1
```

See also: [sprintf], page 299, [num2str], page 96, [int2str], page 97.
num2str ( $x$ )
num2str (x, precision)
num2str (x, format)
Convert a number (or array) to a string (or a character array).
The optional second argument may either give the number of significant digits (precision) to be used in the output or a format template string (format) as in sprintf (see Section 14.2.4 [Formatted Output], page 298). num2str can also process complex numbers.
Examples:

```
num2str (123.456)
    \(\Rightarrow 123.456\)
num2str (123.456, 4)
    \(\Rightarrow 123.5\)
s = num2str ([1, 1.34; 3, 3.56], "\%5.1f")
    \(\Rightarrow\) s \(=\)
        \(1.0 \quad 1.3\)
        \(3.0 \quad 3.6\)
whos s
    \(\Rightarrow\) Variables in the current scope:
        Attr Name Size Bytes Class
            ==== ==== ==== ===== =====
            s \(\quad 2 \mathrm{x} 8\)
                            2x8
                            16 char
        Total is 16 elements using 16 bytes
num2str (1.234 + 27.3i)
    \(\Rightarrow 1.234+27.3 \mathrm{i}\)
```

The num2str function is not very flexible. For better control over the results, use sprintf (see Section 14.2.4 [Formatted Output], page 298).

## Programming Notes:

For matlab compatibility, leading spaces are stripped before returning the string.
Integers larger than flintmax may not be displayed correctly.
For complex $x$, the format string may only contain one output conversion specification and nothing else. Otherwise, results will be unpredictable.

Any optional format specified by the programmer is used without modification. This is in contrast to matlab which tampers with the format based on internal heuristics.

See also: [sprintf], page 299, [int2str], page 97, [mat2str], page 95.
int2str (n)
Convert an integer (or array of integers) to a string (or a character array).

```
int2str (123)
            => 123
s = int2str ([1, 2, 3; 4, 5, 6])
        # s =
            1 2 3
            4 6
        whos s
            => Variables in the current scope:
                Attr Name Size Bytes Class
                ==== ==== ==== ===== =====
                    s 2x7
                                    1 4 ~ c h a r
```

            Total is 14 elements using 14 bytes
    This function is not very flexible. For better control over the results, use sprintf (see Section 14.2.4 [Formatted Output], page 298).
Programming Notes:
Non-integers are rounded to integers before display. Only the real part of complex numbers is displayed.

See also: [sprintf], page 299, [num2str], page 96, [mat2str], page 95.
str2double (s)
Convert a string to a real or complex number.
The string must be in one of the following formats where a and b are real numbers and the complex unit is ' i ' or ' j ':

- $\mathrm{a}+\mathrm{bi}$
- $a+b^{*}$
- $a+i^{*} b$
- $\mathrm{bi}+\mathrm{a}$
- $\mathrm{b}^{*} \mathrm{i}+\mathrm{a}$
- i*b + a

If present, a and/or b are of the form [+-]d[..]d[[eE][+-]d] where the brackets indicate optional arguments and 'd' indicates zero or more digits. The special input values Inf, NaN, and NA are also accepted.
$s$ may be a character string, character matrix, or cell array. For character arrays the conversion is repeated for every row, and a double or complex array is returned. Empty rows in $s$ are deleted and not returned in the numeric array. For cell arrays each character string element is processed and a double or complex array of the same dimensions as $s$ is returned.
For unconvertible scalar or character string input str2double returns a NaN. Similarly, for character array input str2double returns a NaN for any row of $s$ that could not be converted. For a cell array, str2double returns a NaN for any element of $s$ for which conversion fails. Note that numeric elements in a mixed string/numeric cell array are not strings and the conversion will fail for these elements and return NaN . str2double can replace str2num, and it avoids the security risk of using eval on unknown data.
See also: [str2num], page 98.
$x=s t r 2 n u m(s)$
[x, state] = str2num (s)
Convert the string (or character array) $s$ to a number (or an array).
Examples:

```
str2num ("3.141596")
            => 3.141596
str2num (["1, 2, 3"; "4, 5, 6"])
        # 1}\begin{array}{lll}{1}&{2}&{3}\\{4}&{5}&{6}
```

The optional second output, state, is logically true when the conversion is successful. If the conversion fails the numeric output, $x$, is empty and state is false.
Caution: As str2num uses the eval function to do the conversion, str2num will execute any code contained in the string $s$. Use str2double for a safer and faster conversion.
For cell array of strings use str2double.
See also: [str2double], page 97, [eval], page 175.
bin2dec (s)
Return the decimal number corresponding to the binary number represented by the string $s$.
For example:

```
bin2dec ("1110")
    => 14
```

Spaces are ignored during conversion and may be used to make the binary number more readable.

```
bin2dec ("1000 0001")
    => 129
```

If $s$ is a string matrix, return a column vector with one converted number per row of $s$; Invalid rows evaluate to NaN .

If $s$ is a cell array of strings, return a column vector with one converted number per cell element in $s$.

See also: [dec2bin], page 99, [base2dec], page 100, [hex2dec], page 100.
bstr $=$ dec2bin (d)
bstr $=\operatorname{dec} 2 b i n(d, l e n)$
Return a string of ones and zeros representing the conversion of the integer $d$ to a binary number.
If $d$ is a matrix or cell array, return a string matrix with one row for each element in $d$, padded with leading zeros to the width of the largest value.

The optional second argument, len, specifies the minimum number of digits in the result.

For negative elements of $d$, return the binary value of the two's complement. The result is padded with leading ones to $8,16,32$, or 64 bits as appropriate for the magnitude of the input. Positive input elements are padded with leading zeros to the same width. If the second argument len exceeds that calculated width, the result is further padded with leading zeros, for compatibility with matlab.

Examples:

```
dec2bin (14)
    => "1110"
dec2bin (-14)
    => "11110010"
```

Known matlab Incompatibility: matlab's dec2bin allows non-integer values for $d$ as of Release 2022b, but is inconsistent with truncation versus rounding and is also inconsistent with its own dec2hex function. For self-consistency, Octave gives an error for non-integer inputs. Users requiring compatible code for non-integer inputs should make use of fix or round as appropriate.

See also: [bin2dec], page 98, [dec2base], page 100, [dec2hex], page 99.

```
hstr = dec2hex (d)
hstr = dec2hex (d, len)
```

Return a string representing the conversion of the integer $d$ to a hexadecimal (base16) number.
If $d$ is negative, return the hexadecimal equivalent of the two's complement binary value of $d$. If $d$ is a matrix or cell array, return a string matrix with one row for each element in $d$, padded with leading zeros to the width of the largest value.
The optional second argument, len, specifies the minimum number of digits in the result.

Examples:

```
dec2hex (2748)
    => "ABC"
dec2hex (-2)
    => "FE"
```

See also: [hex2dec], page 100, [dec2base], page 100, [dec2bin], page 99.

Return the integer corresponding to the hexadecimal number represented by the string s.

For example:

```
hex2dec ("12B")
            => 299
hex2dec ("12b")
            => 299
```

If $s$ is a string matrix, return a column vector with one converted number per row of $s$; Invalid rows evaluate to NaN .
If $s$ is a cell array of strings, return a column vector with one converted number per cell element in $s$.

See also: [dec2hex], page 99, [base2dec], page 100, [bin2dec], page 98.
dec2base (d, base)
dec2base (d, base, len)
Return a string of symbols in base base corresponding to the non-negative integer $d$.

```
dec2base (123, 3)
    => "11120"
```

If $d$ is a matrix or cell array, return a string matrix with one row per element in $d$, padded with leading zeros to the width of the largest value.
If base is a string then the characters of base are used as the symbols for the digits of $d$. Whitespace (spaces, tabs, newlines, , etc.) may not be used as a symbol.

```
dec2base (123, "aei")
        => "eeeia"
```

The optional third argument, len, specifies the minimum number of digits in the result.

See also: [base2dec], page 100, [dec2bin], page 99, [dec2hex], page 99.
base2dec (s, base)
Convert $s$ from a string of digits in base base to a decimal integer (base 10).

```
base2dec ("11120", 3)
        # 123
```

If $s$ is a string matrix, return a column vector with one value per row of $s$. If a row contains invalid symbols then the corresponding value will be NaN.
If $s$ is a cell array of strings, return a column vector with one value per cell element in $S$.

If base is a string, the characters of base are used as the symbols for the digits of $s$. Space (' ') may not be used as a symbol.

```
base2dec ("yyyzx", "xyz")
```

    \(\Rightarrow 123\)
    See also: [dec2base], page 100, [bin2dec], page 98, [hex2dec], page 100.
$s=$ num2hex ( $n$ )
$s=$ num2hex (n, "cell")
Convert a numeric array to an array of hexadecimal strings.
For example:

```
num2hex ([-1, 1, e, Inf])
=> "bff0000000000000
    3ff0000000000000
    4005bf0a8b145769
    7ff0000000000000"
```

If the argument $n$ is a single precision number or vector, the returned string has a length of 8 . For example:

```
num2hex (single ([-1, 1, e, Inf]))
=> "bf800000
    3f800000
    402df854
    7f800000"
```

With the optional second argument "cell", return a cell array of strings instead of a character array.
See also: [hex2num], page 101, [hex2dec], page 100, [dec2hex], page 99.
$n=$ hex2num (s)
$n=$ hex2num (s, class)
Typecast a hexadecimal character array or cell array of strings to an array of numbers.
By default, the input array is interpreted as a hexadecimal number representing a double precision value. If fewer than 16 characters are given the strings are right padded with ' 0 ' characters.
Given a string matrix, hex2num treats each row as a separate number.

```
hex2num (["4005bf0a8b145769"; "4024000000000000"])
    # [2.7183; 10.000]
```

The optional second argument class may be used to cause the input array to be interpreted as a different value type. Possible values are

## Option Characters

"int8" 2
"uint8" 2
"int16" 4
"uint16" 4
"int32" 8
"uint32" 8
"int64" 16
"uint64" 16
"char" 2
"single" 8
"double" 16
For example:
hex2num (["402df854"; "41200000"], "single")
$\Rightarrow \quad[2.7183 ; 10.000]$
See also: [num2hex], page 101, [hex2dec], page 100, [dec2hex], page 99.
[a, ...] = strread (str)
[a, ...] = strread (str, format)
[a, ...] = strread (str, format, format_repeat)
[a, ...] = strread (str, format, prop1, value1,...)
[a, ...] = strread (str, format, format_repeat, prop1, value1, ...)
This function is obsolete. Use textscan instead.
Read data from a string.
The string str is split into words that are repeatedly matched to the specifiers in format. The first word is matched to the first specifier, the second to the second specifier and so forth. If there are more words than specifiers, the process is repeated until all words have been processed.
The string format describes how the words in str should be parsed. It may contain any combination of the following specifiers:
$\%$ The word is parsed as a string.
\%f
$\%$ n The word is parsed as a number and converted to double.
\%d
$\%$ The word is parsed as a number and converted to int32.
\%*
\%*f
$\%$ *s The word is skipped.
For $\%$ s and $\% \mathrm{~d}, \% \mathrm{f}, \% \mathrm{n}, \% \mathrm{u}$ and the associated $\% *$ s ... specifiers an optional width can be specified as $\% \mathrm{Ns}$, etc. where N is an integer $>1$. For \%f, format specifiers like $\% \mathrm{~N} . \mathrm{Mf}$ are allowed.
literals In addition the format may contain literal character strings; these will be skipped during reading.

Parsed word corresponding to the first specifier are returned in the first output argument and likewise for the rest of the specifiers.
By default, format is "\% $\ddagger$ ", meaning that numbers are read from str. This will do if str contains only numeric fields.

For example, the string

```
str = "\
Bunny Bugs 5.5\n\
Duck Daffy -7.5e-5\n\
Penguin Tux 6"
```

can be read using

```
[a, b, c] = strread (str, "%s %s %f");
```

Optional numeric argument format_repeat can be used for limiting the number of items read:
-1 (default) read all of the string until the end.
$\mathrm{N} \quad$ Read N times nargout items. 0 (zero) is an acceptable value for format_repeat.

The behavior of strread can be changed via property-value pairs. The following properties are recognized:
"commentstyle"
Parts of str are considered comments and will be skipped. value is the comment style and can be any of the following.

- "shell" Everything from \# characters to the nearest end-of-line is skipped.
- "c" Everything between /* and */ is skipped.
- "c++" Everything from // characters to the nearest end-of-line is skipped.
- "matlab" Everything from \% characters to the nearest end-of-line is skipped.
- user-supplied. Two options: (1) One string, or $1 \times 1$ cell string: Skip everything to the right of it; (2) 2x1 cell string array: Everything between the left and right strings is skipped.


## "delimiter"

Any character in value will be used to split str into words (default value $=$ any whitespace). Note that whitespace is implicitly added to the set of delimiter characters unless a "\%s" format conversion specifier is supplied; see "whitespace" parameter below. The set of delimiter characters cannot be empty; if needed Octave substitutes a space as delimiter.
"emptyvalue"
Value to return for empty numeric values in non-whitespace delimited data. The default is NaN. When the data type does not support NaN (int32 for example), then default is zero.
"multipledelimsasone"
Treat a series of consecutive delimiters, without whitespace in between, as a single delimiter. Consecutive delimiter series need not be vertically "aligned".
"treatasempty"
Treat single occurrences (surrounded by delimiters or whitespace) of the string(s) in value as missing values.
"returnonerror"
If value true ( 1 , default), ignore read errors and return normally. If false (0), return an error.
"whitespace"
Any character in value will be interpreted as whitespace and trimmed; the string defining whitespace must be enclosed in double quotes for proper processing of special characters like " $\backslash \mathrm{t}$ ". In each data field, multiple consecutive whitespace characters are collapsed into one space and leading and trailing whitespace is removed. The default value for whitespace is " \b\r\n\t" (note the space). Whitespace is always added to the set of delimiter characters unless at least one "\%s" format conversion specifier is supplied; in that case only whitespace explicitly specified in "delimiter" is retained as delimiter and removed from the set of whitespace characters. If whitespace characters are to be kept as-is (in e.g., strings), specify an empty value (i.e., "") for "whitespace"; obviously, whitespace cannot be a delimiter then.

When the number of words in str doesn't match an exact multiple of the number of format conversion specifiers, strread's behavior depends on the last character of str:
last character $=$ " $\backslash n$ "
Data columns are padded with empty fields or NaN so that all columns have equal length
last character is not " $\backslash \mathrm{n}$ "
Data columns are not padded; strread returns columns of unequal length
See also: [textscan], page 287, [sscanf], page 305.

### 5.4.3 JSON data encoding/decoding

JavaScript Object Notation, in short JSON, is a very common human readable and structured data format. GNU Octave supports encoding and decoding this format with the following two functions.

```
JSON_txt = jsonencode (object)
JSON_txt = jsonencode (..., "ConvertInfAndNaN", TF)
JSON_txt = jsonencode (..., "PrettyPrint",TF)
Encode Octave data types into JSON text.
The input object is an Octave variable to encode.
```

The output JSON_txt is the JSON text that contains the result of encoding object.
If the value of the option "ConvertInfAndNaN" is true then NaN, NA, -Inf, and Inf values will be converted to "null" in the output. If it is false then they will remain as their original values. The default value for this option is true.

If the value of the option "PrettyPrint" is true, the output text will have indentations and line feeds. If it is false, the output will be condensed and written without whitespace. The default value for this option is false.
Programming Notes:

- Complex numbers are not supported.
- classdef objects are first converted to structs and then encoded.
- To preserve escape characters (e.g., " $\backslash \mathrm{n} "$ ), use single-quoted strings.
- Every character after the null character ("\0") in a double-quoted string will be dropped during encoding.
- Encoding and decoding an array is not guaranteed to preserve the dimensions of the array. In particular, row vectors will be reshaped to column vectors.
- Encoding and decoding is not guaranteed to preserve the Octave data type because JSON supports fewer data types than Octave. For example, if you encode an int8 and then decode it, you will get a double.
This table shows the conversions from Octave data types to JSON data types:


## Octave data type

logical scalar
logical vector
logical array
numeric scalar
numeric vector
numeric array
$\mathrm{NaN}, \mathrm{NA}, \operatorname{Inf},-\operatorname{Inf}$
when "ConvertInfAndNaN" = true
NaN, NA, Inf, - Inf
when "ConvertInfAndNaN" = false
empty array
character vector
character array
empty character array
cell scalar
cell vector
cell array
struct scalar
struct vector
struct array
classdef object
Examples:

```
jsonencode ([1, NaN; 3, 4])
=> [[1,null],[3,4]]
```


## JSON data type

Boolean
Array of Boolean, reshaped to row vector
nested Array of Boolean
Number
Array of Number, reshaped to row vector
nested Array of Number
"null"
"NaN", "NaN", "Infinity",
"-Infinity"
" []"
String
Array of String
""
Array
Array, reshaped to row vector
Array, flattened to row vector
Object
Array of Object, reshaped to row vector nested Array of Object
Object

```
jsonencode ([1, NaN; 3, 4], "ConvertInfAndNaN", false)
=> [[1,NaN],[3,4]]
## Escape characters inside a single-quoted string
jsonencode ('\0\a\b\t\n\v\f\r')
=> "\\0\\a\\b\\t\\n\\v\\f\\r"
## Escape characters inside a double-quoted string
jsonencode ("\a\b\t\n\v\f\r")
=> "\u0007\b\t\n\u000B\f\r"
jsonencode ([true; false], "PrettyPrint", true)
=> ans = [
        true,
        false
    ]
jsonencode (['foo', 'bar'; 'foo', 'bar'])
=> ["foobar","foobar"]
jsonencode (struct ('a', Inf, 'b', [], 'c', struct ()))
=> {"a":null,"b":[],"c":{}}
jsonencode (struct ('structarray', struct ('a', {1; 3}, 'b', {2; 4})))
=> {"structarray":[{"a":1,"b":2},{"a":3,"b":4}]}
jsonencode ({'foo'; 'bar'; {'foo'; 'bar'}})
=> ["foo","bar",["foo","bar"]]
jsonencode (containers.Map({'foo'; 'bar'; 'baz'}, [1, 2, 3]))
=> {"bar":2,"baz":3,"foo":1}
```

See also: [jsondecode], page 106.

```
object = jsondecode (JSON_txt)
object = jsondecode (..., "ReplacementStyle", rs)
object = jsondecode (..., "Prefix", pfx)
object = jsondecode (..., "makeValidName", TF)
```

Decode text that is formatted in JSON.
The input JSON_txt is a string that contains JSON text.
The output object is an Octave object that contains the result of decoding JSON_txt.
For more information about the options "ReplacementStyle" and "Prefix", see [matlab.lang.makeValidName], page 137.
If the value of the option "makeValidName" is false then names will not be changed by matlab.lang.makeValidName and the "ReplacementStyle" and "Prefix" options will be ignored.

NOTE: Decoding and encoding JSON text is not guaranteed to reproduce the original text as some names may be changed by matlab.lang.makeValidName.

This table shows the conversions from JSON data types to Octave data types:

## JSON data type

Boolean
Number
String
Object
null, inside a numeric array
null, inside a non-numeric array
Array, of different data types
Array, of Booleans
Array, of Numbers
Array, of Strings
Array of Objects, same field names
Array of Objects, different field names

Octave data type
scalar logical
scalar double
vector of characters
scalar struct (field names of the struct may be different from the keys of the JSON object due to matlab_lang_ makeValidName
NaN
empty double array []
cell array
logical array
double array
cell array of character vectors (cellstr)
struct array
cell array of scalar structs

Examples:

```
jsondecode ('[1, 2, null, 3]')
        # ans =
            1
            2
        NaN
            3
jsondecode ('["foo", "bar", ["foo", "bar"]]')
    => ans =
        {
            [1,1] = foo
            [2,1] = bar
            [3,1] =
            {
            [1,1] = foo
            [2,1] = bar
        }
            }
```

```
jsondecode ('{"nu#m#ber": 7, "s#tr#ing": "hi"}', ...
            'ReplacementStyle', 'delete')
        => scalar structure containing the fields:
        number = 7
        string = hi
jsondecode ('{"nu#m#ber": 7, "s#tr#ing": "hi"}', ...
            'makeValidName', false)
        => scalar structure containing the fields:
        nu#m#ber = 7
        s#tr#ing = hi
jsondecode ('{"1": "one", "2": "two"}', 'Prefix', 'm_')
    => scalar structure containing the fields:
        m_1 = one
        m_2 = two
```

See also: [jsonencode], page 104, [matlab.lang.makeValidName], page 137.

### 5.5 Character Class Functions

Octave also provides the following character class test functions patterned after the functions in the standard C library. They all operate on string arrays and return matrices of zeros and ones. Elements that are nonzero indicate that the condition was true for the corresponding character in the string array. For example:

```
isalpha ("!Q@WERT^Y\&")
    \(\Rightarrow[0,1,0,1,1,1,1,0,1,0]\)
```

isalnum (s)
Return a logical array which is true where the elements of $s$ are letters or digits and false where they are not.
This is equivalent to (isalpha (s) | isdigit (s)).
See also: [isalpha], page 108, [isdigit], page 109, [ispunct], page 109, [isspace], page 109, [iscntrl], page 109.
isalpha (s)
Return a logical array which is true where the elements of $s$ are letters and false where they are not.
This is equivalent to (islower (s) | isupper (s)).
See also: [isdigit], page 109, [ispunct], page 109, [isspace], page 109, [iscntrl], page 109, [isalnum], page 108, [islower], page 109, [isupper], page 109.
isletter (s)
Return a logical array which is true where the elements of $s$ are letters and false where they are not.

This is an alias for the isalpha function.
See also: [isalpha], page 108, [isdigit], page 109, [ispunct], page 109, [isspace], page 109, [iscntrl], page 109, [isalnum], page 108.
islower (s)
Return a logical array which is true where the elements of $s$ are lowercase letters and false where they are not.

See also: [isupper], page 109, [isalpha], page 108, [isletter], page 108, [isalnum], page 108.
isupper (s)
Return a logical array which is true where the elements of $s$ are uppercase letters and false where they are not.
See also: [islower], page 109, [isalpha], page 108, [isletter], page 108, [isalnum], page 108.
isdigit (s)
Return a logical array which is true where the elements of $s$ are decimal digits (0-9) and false where they are not.
See also: [isxdigit], page 109, [isalpha], page 108, [isletter], page 108, [ispunct], page 109, [isspace], page 109, [iscntrl], page 109.
isxdigit (s)
Return a logical array which is true where the elements of $s$ are hexadecimal digits (0-9 and a-fA-F).
See also: [isdigit], page 109.
ispunct (s)
Return a logical array which is true where the elements of $s$ are punctuation characters and false where they are not.

See also: [isalpha], page 108, [isdigit], page 109, [isspace], page 109, [iscntrl], page 109.
isspace (s)
Return a logical array which is true where the elements of $s$ are whitespace characters (space, formfeed, newline, carriage return, tab, and vertical tab) and false where they are not.
See also: [iscntrl], page 109, [ispunct], page 109, [isalpha], page 108, [isdigit], page 109.
iscntrl (s)
Return a logical array which is true where the elements of $s$ are control characters and false where they are not.
See also: [ispunct], page 109, [isspace], page 109, [isalpha], page 108, [isdigit], page 109.
isgraph (s)
Return a logical array which is true where the elements of $s$ are printable characters (but not the space character) and false where they are not.
See also: [isprint], page 110.
isprint (s)
Return a logical array which is true where the elements of $s$ are printable characters (including the space character) and false where they are not.
See also: [isgraph], page 109.
isascii (s)
Return a logical array which is true where the elements of $s$ are ASCII characters (in the range 0 to 127 decimal) and false where they are not.
isstrprop (str, prop)
Test character string properties.
For example:

```
isstrprop ("abc123", "alpha")
# [1, 1, 1, 0, 0, 0]
```

If str is a cell array, isstrpop is applied recursively to each element of the cell array. Numeric arrays are converted to character strings.
The second argument prop must be one of
"alpha" True for characters that are alphabetic (letters).
"alnum"
"alphanum"
True for characters that are alphabetic or digits.
"lower" True for lowercase letters.
"upper" True for uppercase letters.
"digit" True for decimal digits (0-9).
"xdigit" True for hexadecimal digits (a-fA-F0-9).
"space"
"wspace" True for whitespace characters (space, formfeed, newline, carriage return, tab, vertical tab).
"punct" True for punctuation characters (printing characters except space or letter or digit).
"cntrl" True for control characters.
"graph"
"graphic"
True for printing characters except space.
"print" True for printing characters including space.
"ascii" True for characters that are in the range of ASCII encoding.
See also: [isalpha], page 108, [isalnum], page 108, [islower], page 109, [isupper], page 109, [isdigit], page 109, [isxdigit], page 109, [isspace], page 109, [ispunct], page 109, [iscntrl], page 109, [isgraph], page 109, [isprint], page 110, [isascii], page 110.

## 6 Data Containers

Octave includes support for three different mechanisms to contain arbitrary data types in the same variable: Structures, which are C-like, and are indexed with named fields; containers.Map objects, which store data in key/value pairs; and cell arrays, where each element of the array can have a different data type and or shape. Multiple input arguments and return values of functions are organized as another data container, the comma separated list.

### 6.1 Structures

Octave includes support for organizing data in structures. The current implementation uses an associative array with indices limited to strings, but the syntax is more like C-style structures.

### 6.1.1 Basic Usage and Examples

Here are some examples of using data structures in Octave.
Elements of structures can be of any value type. For example, the three expressions

```
x.a = 1;
x.b = [1, 2; 3, 4];
x.c = "string";
```

create a structure with three elements. The '.' character separates the structure name (in the example above $\mathbf{x}$ ) from the field name and indicates to Octave that this variable is a structure. To print the value of the structure you can type its name, just as for any other variable:
x

Note that Octave may print the elements in any order.
Structures may be copied just like any other variable:

```
y = x
    # y =
        scalar structure containing the fields:
            a = 1
            b =
                1 2
                3 4
            c = string
```

Since structures are themselves values, structure elements may reference other structures, as well. The following statement adds the field $d$ to the structure x . The value of field d is itself a data structure containing the single field a, which has a value of 3 .

```
x.d.a = 3;
x.d
    # ans =
        scalar structure containing the fields:
            a=3
x
    # x =
    scalar structure containing the fields:
        a = 1
        b =
            1 2
            34
        c = string
        d =
            scalar structure containing the fields:
            a=3
```

Note that when Octave prints the value of a structure that contains other structures, only a few levels are displayed. For example:

```
a.b.c.d.e = 1;
a
    # a =
            scalar structure containing the fields:
            b =
                scalar structure containing the fields:
                c =
                    scalar structure containing the fields:
                    d: 1x1 scalar struct
```

This prevents long and confusing output from large deeply nested structures. The number of levels to print for nested structures may be set with the function struct_levels_to_ print, and the function print_struct_array_contents may be used to enable printing of the contents of structure arrays.

```
val = struct_levels_to_print ()
```

old_val = struct_levels_to_print (new_val)
struct_levels_to_print (new_val, "local")

Query or set the internal variable that specifies the number of structure levels to display.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [print_struct_array_contents], page 113.

```
val = print_struct_array_contents ()
old_val = print_struct_array_contents (new_val)
print_struct_array_contents (new_val, "local")
```

Query or set the internal variable that specifies whether to print struct array contents.
If true, values of struct array elements are printed. This variable does not affect scalar structures whose elements are always printed. In both cases, however, printing will be limited to the number of levels specified by struct_levels_to_print.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [struct_levels_to_print], page 113.
Functions can return structures. For example, the following function separates the real and complex parts of a matrix and stores them in two elements of the same structure variable y.

```
function y = f (x)
    y.re = real (x);
    y.im = imag (x);
endfunction
```

When called with a complex-valued argument, the function $f$ returns the data structure containing the real and imaginary parts of the original function argument.

```
f (rand (2) + rand (2) * I)
    => ans =
        scalar structure containing the fields:
            re =
            0.040239 0.242160
            0.238081 0.402523
        im =
            0.26475 0.14828
            0.18436 0.83669
```

Function return lists can include structure elements, and they may be indexed like any other variable. For example:

```
[ x.u, x.s(2:3,2:3), x.v ] = svd ([1, 2; 3, 4]);
x
# x =
    scalar structure containing the fields:
        u =
            -0.40455 -0.91451
            -0.91451 0.40455
        s =
            0.00000 0.00000 0.00000
            0.00000 5.46499 0.00000
            0.00000 0.00000 0.36597
        v =
            -0.57605 0.81742
            -0.81742 -0.57605
```

It is also possible to cycle through all the elements of a structure in a loop, using a special form of the for statement (see Section 10.5.1 [Looping Over Structure Elements], page 185).

### 6.1.2 Structure Arrays

A structure array is a particular instance of a structure, where each of the fields of the structure is represented by a cell array. Each of these cell arrays has the same dimensions. Conceptually, a structure array can also be seen as an array of structures with identical fields. An example of the creation of a structure array is

```
x(1).a = "string1";
x(2).a = "string2";
x(1).b = 1;
x(2).b = 2;
```

which creates a 1-by-2 structure array with two fields. Another way to create a structure array is with the struct function (see Section 6.1.3 [Creating Structures], page 116). As previously, to print the value of the structure array, you can type its name:
x

```
# x =
    {
        1x2 struct array containing the fields:
```

            a
            b
    \}
    Individual elements of the structure array can be returned by indexing the variable like $x(1)$, which returns a structure with two fields:

```
x(1)
# ans =
    {
        a = string1
        b}=
    }
```

Furthermore, the structure array can return a comma separated list of field values (see Section 6.4 [Comma Separated Lists], page 134), if indexed by one of its own field names. For example:

$$
\begin{aligned}
\mathrm{x} \cdot \mathrm{a} & \\
& \\
& \text { ans }=\text { string1 } \\
& \text { ans }=\text { string2 }
\end{aligned}
$$

Here is another example, using this comma separated list on the left-hand side of an assignment:

```
[x.a] = deal ("new string1", "new string2");
x(1).a
        a ans = new string1
x(2).a
        a ans = new string2
```

Just as for numerical arrays, it is possible to use vectors as indices (see Section 8.1 [Index Expressions], page 151):

```
x(3:4) = x(1:2);
[x([1,3]).a] = deal ("other string1", "other string2");
x.a
    #
        ans = other string1
        ans = new string2
        ans = other string2
        ans = new string2
```

The function size will return the size of the structure. For the example above

```
size (x)
    a ans =
    14
```

Elements can be deleted from a structure array in a similar manner to a numerical array, by assigning the elements to an empty matrix. For example

```
in = struct ("call1", {x, Inf, "last"},
    "call2", {x, Inf, "first"})
    # in =
        {
            1x3 struct array containing the fields:
                call1
                call2
    }
in(1) = [];
in.call1
    #
    ans = Inf
    ans = last
```


### 6.1.3 Creating Structures

Besides the index operator ".", Octave can use dynamic naming "(var)" or the struct function to create structures. Dynamic naming uses the string value of a variable as the field name. For example:

```
a = "field2";
x.a = 1;
x.(a) = 2;
x
    # x =
        {
            a = 1
            field2 = 2
        }
```

Dynamic indexing also allows you to use arbitrary strings, not merely valid Octave identifiers (note that this does not work on matlab):

```
a = "long field with spaces (and funny char\$)";
x.a = 1;
\(x .(a)=2 ;\)
x
        \(\Rightarrow \mathrm{x}=\)
        \{
            \(\mathrm{a}=1\)
            long field with spaces (and funny char\$) = 2
        \}
```

The warning id Octave:language-extension can be enabled to warn about this usage. See [warning_ids], page 248.

More realistically, all of the functions that operate on strings can be used to build the correct field name before it is entered into the data structure.

```
names = ["Bill"; "Mary"; "John"];
ages = [37; 26; 31];
for i = 1:rows (names)
    database.(names(i,:)) = ages(i);
endfor
database
        | database =
        {
            Bill = 37
                        Mary = 26
                        John = 31
            }
```

The third way to create structures is the struct command. struct takes pairs of arguments, where the first argument in the pair is the fieldname to include in the structure and the second is a scalar or cell array, representing the values to include in the structure or structure array. For example:

```
struct ("field1", 1, "field2", 2)
m ans =
    {
        field1 = 1
        field2 = 2
    }
```

If the values passed to struct are a mix of scalar and cell arrays, then the scalar arguments are expanded to create a structure array with a consistent dimension. For example:

```
s = struct ("field1", {1, "one"}, "field2", {2, "two"},
    "field3", 3);
s.field1
    #
        ans = 1
            ans = one
s.field2
        #
        ans = 2
        ans = two
s.field3
    #
        ans = 3
        ans = 3
```

If you want to create a struct which contains a cell array as an individual field, you must wrap it in another cell array as shown in the following example:

```
    struct ("field1", \{\{1, "one"\}\}, "field2", 2)
        \(\Rightarrow\) ans \(=\)
            \{
                field1 =
            \{
                \([1,1]=1\)
                \([1,2]=\) one
        \}
```

            field2 \(=2\)
        \}
    $s=$ struct ()
$s=$ struct (field1, value1, field2, value2, ...)
$s=$ struct (obj)

Create a scalar or array structure and initialize its values.
The field1, field2, ... variables are strings specifying the names of the fields and the value1, value $2, \ldots$ variables can be of any type.
If the values are cell arrays, create a structure array and initialize its values. The dimensions of each cell array of values must match. Singleton cells and non-cell values are repeated so that they fill the entire array. If the cells are empty, create an empty structure array with the specified field names.
If the argument is an object, return the underlying struct.
Observe that the syntax is optimized for struct arrays. Consider the following examples:

```
struct ("foo", 1)
    # scalar structure containing the fields:
        foo = 1
struct ("foo", {})
    | 0x0 struct array containing the fields:
        foo
struct ("foo", { {} })
    # scalar structure containing the fields:
        foo = {}(0x0)
struct ("foo", {1, 2, 3})
    | 1x3 struct array containing the fields:
        foo
```

The first case is an ordinary scalar struct - one field, one value. The second produces an empty struct array with one field and no values, since being passed an empty cell array of struct array values. When the value is a cell array containing a single entry, this becomes a scalar struct with that single entry as the value of the field. That single entry happens to be an empty cell array.
Finally, if the value is a non-scalar cell array, then struct produces a struct array.
See also: [cell2struct], page 134, [fieldnames], page 119, [getfield], page 121, [setfield], page 120, [rmfield], page 121, [isfield], page 120, [orderfields], page 121, [isstruct], page 119, [structfun], page 635.

The function isstruct can be used to test if an object is a structure or a structure array.
isstruct (x)
Return true if $x$ is a structure or a structure array.
See also: [ismatrix], page 68, [iscell], page 126, [isa], page 41.

### 6.1.4 Manipulating Structures

Other functions that can manipulate the fields of a structure are given below.
numfields (s)
Return the number of fields of the structure $s$.
See also: [fieldnames], page 119.

```
names = fieldnames (struct)
```

names $=$ fieldnames (obj)
names $=$ fieldnames (javaobj)
names = fieldnames ("javaclassname")

Return a cell array of strings with the names of the fields in the specified input.
When the input is a structure struct, the names are the elements of the structure.

When the input is an Octave object obj, the names are the public properties of the object.
When the input is a Java object javaobj or a string containing the name of a Java class javaclassname, the names are the public fields (data members) of the object or class.
See also: [numfields], page 119, [isfield], page 120, [orderfields], page 121, [struct], page 118, [properties], page 909.
isfield (x, "name")
isfield (x, name)
Return true if the $x$ is a structure and it includes an element named name.
If name is a cell array of strings then a logical array of equal dimension is returned.
See also: [fieldnames], page 119.
sout $=$ setfield (s, field, val)
sout $=$ setfield (s, sidx1, field1, fidx1, sidx2, field2, fidx2, ..., val)
Return a copy of the structure $s$ with the field member field set to the value val.
For example:

```
s = struct ();
s = setfield (s, "foo bar", 42);
```

This is equivalent to

```
s.("foo bar") = 42;
```

Note that ordinary structure syntax s.foo bar $=42$ cannot be used here, as the field name is not a valid Octave identifier because of the space character. Using arbitrary strings for field names is incompatible with matlab, and this usage will emit a warning if the warning ID Octave:language-extension is enabled. See [warning_ ids], page 248.
With the second calling form, set a field of a structure array. The input sidx selects an element of the structure array, field specifies the field name of the selected element, and fidx selects which element of the field (in the case of an array or cell array). The sidx, field, and fidx inputs can be repeated to address nested structure array elements. The structure array index and field element index must be cell arrays while the field name must be a string.
For example:

```
s = struct ("baz", 42);
setfield (s, {1}, "foo", {1}, "bar", 54)
#
    ans =
        scalar structure containing the fields:
            baz = 42
            foo =
                scalar structure containing the fields:
                bar = 54
```

The example begins with an ordinary scalar structure to which a nested scalar structure is added. In all cases, if the structure index sidx is not specified it defaults to

1 (scalar structure). Thus, the example above could be written more concisely as setfield (s, "foo", "bar", 54)
Finally, an example with nested structure arrays:

```
sa.foo = 1;
sa = setfield (sa, {2}, "bar", {3}, "baz", {1, 4}, 5);
sa(2).bar(3)
#
    ans =
        scalar structure containing the fields:
            baz = 0 0 0 5
```

Here sa is a structure array whose field at elements 1 and 2 is in turn another structure array whose third element is a simple scalar structure. The terminal scalar structure has a field which contains a matrix value.
Note that the same result as in the above example could be achieved by:

```
sa.foo = 1;
sa(2).bar(3).baz(1,4) = 5
```

See also: [getfield], page 121, [rmfield], page 121, [orderfields], page 121, [isfield], page 120, [fieldnames], page 119, [isstruct], page 119, [struct], page 118.
val = getfield (s, field)
val = getfield (s, sidx1, field1, fidx1,...)
Get the value of the field named field from a structure or nested structure $s$.
If $s$ is a structure array then sidx selects an element of the structure array, field specifies the field name of the selected element, and fidx selects which element of the field (in the case of an array or cell array). For a more complete description of the syntax, see [setfield], page 120 .
See also: [setfield], page 120, [rmfield], page 121, [orderfields], page 121, [isfield], page 120, [fieldnames], page 119, [isstruct], page 119, [struct], page 118.

```
sout = rmfield (s, "f")
sout = rmfield (s,f)
```

Return a copy of the structure (array) $s$ with the field $f$ removed.
If $f$ is a cell array of strings or a character array, remove each of the named fields.
See also: [orderfields], page 121, [fieldnames], page 119, [isfield], page 120.

```
sout = orderfields (s1)
sout = orderfields (s1, s2)
sout = orderfields (s1, {cellstr})
sout = orderfields (s1, p)
[sout, p] = orderfields (...)
```

Return a copy of s1 with fields arranged alphabetically, or as specified by the second input.
Given one input struct s1, arrange field names alphabetically.
If a second struct argument is given, arrange field names in s1 as they appear in s2. The second argument may also specify the order in a cell array of strings cellstr. The second argument may also be a permutation vector.

The optional second output argument $p$ is the permutation vector which converts the original name order to the new name order.

Examples:

```
s = struct ("d", 4, "b", 2, "a", 1, "c", 3);
t1 = orderfields (s)
    t1 =
        scalar structure containing the fields:
            a = 1
            b = 2
            c = 3
            d = 4
t = struct ("d", {}, "c", {}, "b", {}, "a", {});
t2 = orderfields (s, t)
        t2 =
            scalar structure containing the fields:
                d = 4
                    c = 3
            b = 2
            a}=
t3 = orderfields (s, [3, 2, 4, 1])
    t t3 =
            scalar structure containing the fields:
                    a = 1
            b = 2
            c = 3
            d = 4
[t4, p] = orderfields (s, {"d", "c", "b", "a"})
    t t4 =
            scalar structure containing the fields:
                    d = 4
                    c = 3
            b = 2
            a = 1
        p =
            1
            4
            2
            3
```

See also: [fieldnames], page 119, [getfield], page 121, [setfield], page 120, [rmfield], page 121, [isfield], page 120, [isstruct], page 119, [struct], page 118.
substruct (type, subs, ...)
Create a subscript structure for use with subsref or subsasgn.
For example:

```
idx = substruct ("()", {3, ":"})
    # idx =
        scalar structure containing the fields:
            type = ()
            subs =
            {
                [1,1] = 3
                    [1,2] = :
            }
x = [1, 2, 3;
            4, 5, 6;
            7, 8, 9];
subsref (x, idx)
    # 7 8 9
```

See also: [subsref], page 894, [subsasgn], page 896.

### 6.1.5 Processing Data in Structures

The simplest way to process data in a structure is within a for loop (see Section 10.5.1 [Looping Over Structure Elements], page 185). A similar effect can be achieved with the structfun function, where a user defined function is applied to each field of the structure. See [structfun], page 635.

Alternatively, to process the data in a structure, the structure might be converted to another type of container before being treated.

```
c = struct2cell (s)
```

Create a new cell array from the objects stored in the struct object.
If $f$ is the number of fields in the structure, the resulting cell array will have a dimension vector corresponding to $[f$ size $(s)]$. For example:

```
s = struct ("name", {"Peter", "Hannah", "Robert"},
    "age", {23, 16, 3});
c = struct2cell (s)
    => c = {2x1x3 Cell Array}
c(1,1,:)(:)
    #
        {
            [1,1] = Peter
            [2,1] = Hannah
            [3,1] = Robert
        }
c(2,1,:)(:)
        #
            {
            [1,1] = 23
            [2,1] = 16
            [3,1] = 3
            }
```

See also: [cell2struct], page 134, [namedargs2cell], page 124, [fieldnames], page 119.
$c=$ namedargs2cell (s)
Create a cell array of field name/value pairs from a scalar structure.
Example:

```
s.Name = "Peter";
s.Height = 185;
s.Age = 42;
c = namedargs2cell (s)
    => { "Name", "Peter", "Height", 185, "Age", 42 }
```

See also: [struct2cell], page 123.

## 6.2 containers.Map

```
m = containers.Map ()
m = containers.Map (keys, vals)
m = containers.Map (keys, vals, "UniformValues", is_uniform)
m = containers.Map ("KeyType", kt, "ValueType", vt)
```

Create an object of the containers.Map class that stores a list of key/value pairs.
keys is an array of unique keys for the map. The keys can be numeric scalars or strings. The type for numeric keys may be one of "double", "single", "int32", "uint32", "int64", or "uint64". Other numeric or logical keys will be converted to "double". A single string key may be entered as is. Multiple string keys are entered as a cell array of strings.
vals is an array of values for the map with the same number of elements as keys.
When called with no input arguments a default map is created with strings as the key type and "any" as the value type.
The "UniformValues" option specifies whether the values of the map must be strictly of the same type. If is_uniform is true, any values which would be added to the map are first validated to ensure they are of the correct type.
When called with "KeyType" and "ValueType" arguments, create an empty map with the specified types. The inputs $k t$ and vt are the types for the keys and values of the map respectively. Allowed values for kt are "char", "double", "single", "int32", "uint32", "int64", "uint64". Allowed values for vt are "any", "char", "double", "single", "int32", "uint32", "int64", "uint64", "logical".
The return value $m$ is an object of the containers.Map class.
See also: [struct], page 118.

### 6.3 Cell Arrays

It can be both necessary and convenient to store several variables of different size or type in one variable. A cell array is a container class able to do just that. In general cell arrays work just like $N$-dimensional arrays with the exception of the use of '\{' and ' $\}$ ' as allocation and indexing operators.

### 6.3.1 Basic Usage of Cell Arrays

As an example, the following code creates a cell array containing a string and a 2 -by- 2 random matrix

```
c = {"a string", rand(2, 2)};
```

To access the elements of a cell array, it can be indexed with the $\{$ and $\}$ operators. Thus, the variable created in the previous example can be indexed like this:

```
c{1}
    a ans = a string
```

As with numerical arrays several elements of a cell array can be extracted by indexing with a vector of indexes

```
c{1:2}
    ans = a string
    => ans =
\begin{tabular}{ll}
0.593993 & 0.627732 \\
0.377037 & 0.033643
\end{tabular}
```

The indexing operators can also be used to insert or overwrite elements of a cell array. The following code inserts the scalar 3 on the third place of the previously created cell array

```
c{3} = 3
    C c =
        {
            [1,1] = a string
            [1,2] =
            0.593993 0.627732
            0.377037 0.033643
            [1,3] = 3
        }
```

Details on indexing cell arrays are explained in Section 6.3.3 [Indexing Cell Arrays], page 130 .

In general nested cell arrays are displayed hierarchically as in the previous example. In some circumstances it makes sense to reference them by their index, and this can be performed by the celldisp function.

```
celldisp (c)
celldisp (c, name)
```

Recursively display the contents of a cell array.
By default the values are displayed with the name of the variable $c$. However, this name can be replaced with the variable name. For example:

```
c = {1, 2, {31, 32}};
celldisp (c, "b")
    #
        b{1} =
            1
            b{2} =
            2
            b{3}{1} =
                31
            b{3}{2} =
                32
```

See also: [disp], page 271.
To test if an object is a cell array, use the iscell function. For example:

```
iscell (c)
    ans = 1
iscell (3)
    ans = 0
```

iscell (x)

Return true if $x$ is a cell array object.
See also: [ismatrix], page 68, [isstruct], page 119, [iscellstr], page 133, [isa], page 41.

### 6.3.2 Creating Cell Arrays

The introductory example (see Section 6.3.1 [Basic Usage of Cell Arrays], page 125) showed how to create a cell array containing currently available variables. In many situations, however, it is useful to create a cell array and then fill it with data.

The cell function returns a cell array of a given size, containing empty matrices. This function is similar to the zeros function for creating new numerical arrays. The following example creates a 2 -by- 2 cell array containing empty matrices

```
c = cell (2,2)
    # c =
        {
            [1,1] = [] (0x0)
            [2,1] = [] (0x0)
            [1,2] = [](0x0)
            [2,2] = [] (0x0)
        }
```

Just like numerical arrays, cell arrays can be multi-dimensional. The cell function accepts any number of positive integers to describe the size of the returned cell array. It is also possible to set the size of the cell array through a vector of positive integers. In the following example two cell arrays of equal size are created, and the size of the first one is displayed

```
c1 = cell (3, 4, 5);
c2 = cell ( [3, 4, 5] );
size (c1)
    a ans =
    3 4 5
```

As can be seen, the [size], page 47, function also works for cell arrays. As do other functions describing the size of an object, such as [length], page 47, [numel], page 46, [rows], page 46, and [columns], page 46.

```
cell (n)
```

cell ( $m, n$ )
$\operatorname{cell}(m, n, k, \ldots)$
cell ([m n ...])

Create a new cell array object.
If invoked with a single scalar integer argument, return a square NxN cell array. If invoked with two or more scalar integer arguments, or a vector of integer values, return an array with the given dimensions.

See also: [cellstr], page 133, [mat2cell], page 128, [num2cell], page 127, [struct2cell], page 123.

As an alternative to creating empty cell arrays, and then filling them, it is possible to convert numerical arrays into cell arrays using the num2cell, mat2cell and cellslices functions.

```
C = num2cell (A)
C = num2cell (A, dim)
```

Convert the numeric matrix $A$ to a cell array.
When no $\operatorname{dim}$ is specified, each element of $A$ becomes a 1 x 1 element in the output $C$. If $\operatorname{dim}$ is defined then individual elements of $C$ contain all of the elements from $A$ along the specified dimension. dim may also be a vector of dimensions with the same rule applied.
For example:

```
x = [1,2;3,4]
#
    1 2
    3 4
## each element of A becomes a 1x1 element of C
num2cell (x)
    #
        {
            [1,1] = 1
            [2,1] = 3
            [1,2] = 2
            [2,2] = 4
            }
```

```
## all rows (dim 1) of A appear in each element of C
num2cell (x, 1)
    =>
        {
        [1,1] =
            1
            3
            [1,2] =
            2
            4
        }
## all columns (dim 2) of A appear in each element of C
num2cell (x, 2)
    #
        {
        [1,1] =
            1 2
        [2,1] =
            3 4
        }
## all rows and cols appear in each element of C
## (hence, only 1 output)
num2cell (x, [1, 2])
    =>
    {
        [1,1] =
            1 2
            34
    }
```

See also: [mat2cell], page 128.
$C=$ mat2cell (A, dim1, dim2, ..., dimi, ..., dimn)
$C=$ mat2cell (A, rowdim)
Convert the matrix $A$ to a cell array.
Each dimension argument (dim1, dim2, etc.) is a vector of integers which specifies how to divide that dimension's elements amongst the new elements in the output $C$. The number of elements in the $i$-th dimension is size ( $A, i$ ). Because all elements in $A$ must be partitioned, there is a requirement that sum (dimi) $==\operatorname{size}(A, i)$. The size of the output cell $C$ is numel (dim1) x numel (dim2) x . . x numel (dimn).
Given a single dimensional argument, rowdim, the output is divided into rows as specified. All other dimensions are not divided and thus all columns (dim 2), pages (dim 3), etc. appear in each output element.
Examples

```
x = reshape (1:12, [3, 4])'
#
    1 2 
```

```
        4 5
        7 8 9
        10}1
## The 4 rows (dim1) are divided in to two cell elements
## with 2 rows each.
## The 3 cols (dim2) are divided in to three cell elements
## with 1 col each.
mat2cell (x, [2,2], [1,1,1])
#
{
    [1,1] =
        1
        4
        [2,1] =
            7
            10
    [1,2] =
        2
        5
    [2,2] =
        8
        1 1
    [1,3] =
        3
        6
    [2,3] =
        9
        1 2
}
```

```
## The 4 rows (dim1) are divided in to two cell elements
## with a 3/1 split.
## All columns appear in each output element.
mat2cell (x, [3,1])
#
{
    [1,1] =
            1 2 3
            4 5 6
            7 8 9
            [2,1] =
            10}11\quad1
}
```

See also: [num2cell], page 127, [cell2mat], page 134.
sl = cellslices ( $x, l b, u b, d i m)$
Given an array $x$, this function produces a cell array of slices from the array determined by the index vectors $l b, u b$, for lower and upper bounds, respectively.
In other words, it is equivalent to the following code:

```
n = length (lb);
sl = cell (1, n);
for i = 1:length (lb)
    sl{i} = x(:,...,lb(i):ub(i),...,:);
endfor
```

The position of the index is determined by dim. If not specified, slicing is done along the first non-singleton dimension.
See also: [cell2mat], page 134, [cellindexmat], page 132, [cellfun], page 633.

### 6.3.3 Indexing Cell Arrays

As shown in see Section 6.3.1 [Basic Usage of Cell Arrays], page 125, elements can be extracted from cell arrays using the ' $\{$ ' and ' $\}$ ' operators. If you want to extract or access subarrays which are still cell arrays, you need to use the '(' and ')' operators. The following example illustrates the difference:

```
c = {"1", "2", "3"; "x", "y", "z"; "4", "5", "6"};
c{2,3}
    => ans = z
c (2,3)
    ans =
        {
            [1,1] = z
        }
```

So with ' $\}$ ' you access elements of a cell array, while with '()' you access a sub array of a cell array.

Using the '(' and ')' operators, indexing works for cell arrays like for multi-dimensional arrays. As an example, all the rows of the first and third column of a cell array can be set to 0 with the following command:

$$
\begin{aligned}
& c(:, \quad[1,3])=\{0\} \\
& \Rightarrow= \\
&\{ \\
& {[1,1] }=0 \\
& {[2,1] }=0 \\
& {[3,1] }=0 \\
& {[1,2] }=2 \\
& {[2,2] }=y \\
& {[3,2] }=5 \\
& {[1,3] }=0 \\
& {[2,3] }=0 \\
& {[3,3] }=0 \\
&\}
\end{aligned}
$$

Note, that the above can also be achieved like this:

$$
c(:,[1,3])=0 ;
$$

Here, the scalar ' 0 ' is automatically promoted to cell array ' $\{0\}$ ' and then assigned to the subarray of $c$.

To give another example for indexing cell arrays with '()', you can exchange the first and the second row of a cell array as in the following command:

```
c = {1, 2, 3; 4, 5, 6};
c([1, 2], :) = c([2, 1], :)
        =
            {
            [1,1] = 4
            [2,1] = 1
            [1,2] = 5
            [2,2] = 2
            [1,3] = 6
            [2,3] = 3
        }
```

Accessing multiple elements of a cell array with the ' $\{$ ' and ' $\}$ ' operators will result in a comma-separated list of all the requested elements (see Section 6.4 [Comma Separated Lists], page 134). Using the ' $\{$ ' and ' $\}$ ' operators the first two rows in the above example can be swapped back like this:

```
[c{[1,2], :}] = deal (c{[2, 1], :})
    =
        {
            [1,1] = 1
            [2,1] = 4
            [1,2] = 2
            [2,2] = 5
            [1,3] = 3
            [2,3] = 6
        }
```

As for struct arrays and numerical arrays, the empty matrix '[]' can be used to delete elements from a cell array:

```
x = \{"1", "2"; "3", "4"\};
\(x(1,:)=[]\)
    \(\Rightarrow \mathrm{x}=\)
        \{
        \([1,1]=3\)
    \([1,2]=4\)
    \}
```

The following example shows how to just remove the contents of cell array elements but not delete the space for them:

```
x = {"1", "2"; "3", "4"};
x(1, :) = {[]}
# x =
    {
        [1,1] = [] (0x0)
        [2,1] = 3
        [1,2] = [](0x0)
        [2,2] = 4
    }
```

The indexing operations operate on the cell array and not on the objects within the cell array. By contrast, cellindexmat applies matrix indexing to the objects within each cell array entry and returns the requested values.
$y=$ cellindexmat (x, varargin)
Perform indexing of matrices in a cell array.
Given a cell array of matrices $x$, this function computes

```
Y = cell (size (X));
for i = 1:numel (X)
    Y{i} = X{i}(varargin{1}, varargin{2}, ..., varargin{N});
endfor
```

The indexing arguments may be scalar (2), arrays ([1, 3]), ranges (1:3), or the colon operator (":"). However, the indexing keyword end is not available.

See also: [cellslices], page 130, [cellfun], page 633.

### 6.3.4 Cell Arrays of Strings

One common use of cell arrays is to store multiple strings in the same variable. It is also possible to store multiple strings in a character matrix by letting each row be a string. This, however, introduces the problem that all strings must be of equal length. Therefore, it is recommended to use cell arrays to store multiple strings. For cases, where the character matrix representation is required for an operation, there are several functions that convert a cell array of strings to a character array and back. char and strvcat convert cell arrays to a character array (see Section 5.3.2 [Concatenating Strings], page 77), while the function cellstr converts a character array to a cell array of strings:

```
    \(\mathrm{a}=\) ["hello"; "world"];
    \(c=c e l l s t r(a)\)
        \(\Rightarrow c=\)
            \{
                [1, 1] = hello
                    \([2,1]=\) world
            \}
```

cstr $=$ cellstr (strmat)

Create a new cell array object from the elements of the string array strmat.
Each row of strmat becomes an element of cstr. Any trailing spaces in a row are deleted before conversion.

To convert back from a cellstr to a character array use char.
See also: [cell], page 127, [char], page 78.
One further advantage of using cell arrays to store multiple strings is that most functions for string manipulations included with Octave support this representation. As an example, it is possible to compare one string with many others using the strcmp function. If one of the arguments to this function is a string and the other is a cell array of strings, each element of the cell array will be compared to the string argument:

```
c = {"hello", "world"};
strcmp ("hello", c)
    => ans =
        1 0
```

The following string functions support cell arrays of strings: char, strvcat, strcat (see Section 5.3.2 [Concatenating Strings], page 77), strcmp, strncmp, strcmpi, strncmpi (see Section 5.3.4 [Searching in Strings], page 84), str2double, deblank, strtrim, strtrunc, strfind, strmatch, , regexp, regexpi (see Section 5.3 [String Operations], page 74) and str2double (see Section 5.4 [Converting Strings], page 94).

The function iscellstr can be used to test if an object is a cell array of strings.
iscellstr (cell)
Return true if every element of the cell array cell is a character string.
See also: [ischar], page 73 , [isstring], page 73 .

### 6.3.5 Processing Data in Cell Arrays

Data that is stored in a cell array can be processed in several ways depending on the actual data. The simplest way to process that data is to iterate through it using one or more for loops. The same idea can be implemented more easily through the use of the cellfun function that calls a user-specified function on all elements of a cell array. See [cellfun], page 633 .

An alternative is to convert the data to a different container, such as a matrix or a data structure. Depending on the data this is possible using the cell2mat and cell2struct functions.
$m=$ cell2mat (c)
Convert the cell array $c$ into a matrix by concatenating all elements of $c$ into a hyperrectangle.
Elements of $c$ must be numeric, logical, or char matrices; or cell arrays; or structs; and cat must be able to concatenate them together.
See also: [mat2cell], page 128, [num2cell], page 127.
cell2struct (cell, fields)
cell2struct (cell, fields, dim)
Convert cell to a structure.
The number of fields in fields must match the number of elements in cell along dimension dim, that is numel (fields) == size (cell, dim). If dim is omitted, a value of 1 is assumed.

```
A = cell2struct ({"Peter", "Hannah", "Robert";
                            185, 170, 168},
                                {"Name","Height"}, 1);
A(1)
    #
        {
        Name = Peter
        Height = 185
        }
```

See also: [struct2cell], page 123, [cell2mat], page 134, [struct], page 118.

### 6.4 Comma Separated Lists

Comma separated lists ${ }^{1}$ are the basic argument type to all Octave functions - both for input and return arguments. In the example
$\max (a, b)$
' $a, b$ ' is a comma separated list. Comma separated lists can appear on both the right and left hand side of an assignment. For example

```
x = [1 0 1 0 0 1 1; 0 0 0 0 0 0 7];
[i, j] = find (x, 2, "last");
```

[^1]Here, ' $x, 2$, "last"' is a comma separated list constituting the input arguments of find. find returns a comma separated list of output arguments which is assigned element by element to the comma separated list ' $i, j$ '.

Another example of where comma separated lists are used is in the creation of a new array with [] (see Section 4.1 [Matrices], page 52) or the creation of a cell array with \{\} (see Section 6.3.1 [Basic Usage of Cell Arrays], page 125). In the expressions

$$
\begin{aligned}
& a=[1,2,3,4] ; \\
& c=\{4,5,6,7\} ;
\end{aligned}
$$

both ' $1,2,3,4$ ' and ' $4,5,6,7$ ' are comma separated lists.
Comma separated lists cannot be directly manipulated by the user. However, both structure arrays and cell arrays can be converted into comma separated lists, and thus used in place of explicitly written comma separated lists. This feature is useful in many ways, as will be shown in the following subsections.

### 6.4.1 Comma Separated Lists Generated from Cell Arrays

As has been mentioned above (see Section 6.3.3 [Indexing Cell Arrays], page 130), elements of a cell array can be extracted into a comma separated list with the \{ and \} operators. By surrounding this list with [ and ], it can be concatenated into an array. For example:

```
a = {1, [2, 3], 4, 5, 6};
b = [a{1:4}]
    => b =
\(\begin{array}{lllll}1 & 2 & 3 & 4 & 5\end{array}\)
```

Similarly, it is possible to create a new cell array containing cell elements selected with $\}$. By surrounding the list with ' $\{$ ' and ' $\}$ ' a new cell array will be created, as the following example illustrates:

```
a = {1, rand(2, 2), "three"};
b}={a{[1,3]}
    # b =
    {
        [1,1] = 1
        [1,2] = three
    }
```

Furthermore, cell elements (accessed by \{\}) can be passed directly to a function. The list of elements from the cell array will be passed as an argument list to a given function as if it is called with the elements as individual arguments. The two calls to printf in the following example are identical but the latter is simpler and can handle cell arrays of an arbitrary size:

```
c = {"GNU", "Octave", "is", "Free", "Software"};
printf ("%s ", c{1}, c{2}, c{3}, c{4}, c{5});
    GNU Octave is Free Software
printf ("%s ", c{:});
    GNU Octave is Free Software
```

If used on the left-hand side of an assignment, a comma separated list generated with \{\} can be assigned to. An example is
$\operatorname{in}\{1\}=[10,20,30] ;$

```
in\{2\} \(=\inf ;\)
in\{3\} = "last";
in\{4\} = "first";
out = cell (4, 1);
[out\{1:3\}] = in\{1 : 3\};
[out\{4:6\}] \(=\operatorname{in}\{[1,2,4]\})\)
    \(\Rightarrow\) out \(=\)
    \{
        \([1,1]=\)
            \(10 \quad 20 \quad 30\)
        \([2,1]=\operatorname{Inf}\)
        \([3,1]=\) last
        \([4,1]=\)
            \(10 \quad 20 \quad 30\)
        [5,1] = Inf
        [6,1] = first
    \}
```


### 6.4.2 Comma Separated Lists Generated from Structure Arrays

Structure arrays can equally be used to create comma separated lists. This is done by addressing one of the fields of a structure array. For example:

```
x = ceil (randn (10, 1));
in = struct ("call1", {x, 3, "last"},
    "call2", {x, inf, "first"});
out = struct ("call1", cell (2, 1), "call2", cell (2, 1));
[out.call1] = find (in.call1);
[out.call2] = find (in.call2);
```


## 7 Variables

Variables let you give names to values and refer to them later. You have already seen variables in many of the examples. The name of a variable must be a sequence of letters, digits and underscores, but it may not begin with a digit. Octave does not enforce a limit on the length of variable names, but it is seldom useful to have variables with names longer than about 30 characters. The following are all valid variable names

```
x
x15
__foo_bar_baz__
fucnrdthsucngtagdjb
```

However, names like __foo_bar_baz__ that begin and end with two underscores are understood to be reserved for internal use by Octave. You should not use them in code you write, except to access Octave's documented internal variables and built-in symbolic constants.

Case is significant in variable names. The symbols a and A are distinct variables.
A variable name is a valid expression by itself. It represents the variable's current value. Variables are given new values with assignment operators and increment operators. See Section 8.6 [Assignment Expressions], page 169.

There is one automatically created variable with a special meaning. The ans variable always contains the result of the last computation, where the output wasn't assigned to any variable. The code $\mathrm{a}=\cos$ (pi) will assign the value -1 to the variable a , but will not change the value of ans. However, the code cos (pi) will set the value of ans to -1 .

Variables in Octave do not have fixed types, so it is possible to first store a numeric value in a variable and then to later use the same name to hold a string value in the same program. Variables may not be used before they have been given a value. Doing so results in an error.
ans
[Automatic Variable]
The most recently computed result that was not explicitly assigned to a variable.
For example, after the expression

```
3^2 + 4^2
```

is evaluated, the value returned by ans is 25 .
isvarname (name)
Return true if name is a valid variable name.
A valid variable name is composed of letters, digits, and underscores (" - "), and the first character must not be a digit.
See also: [iskeyword], page 1099, [exist], page 144, [who], page 142.

```
varname = matlab.lang.makeValidName (str)
varname = matlab.lang.makeValidName (..., "ReplacementStyle", rs)
varname = matlab.lang.makeValidName (..., "Prefix", pfx)
[varname, ismodified] = matlab.lang.makeValidName (...)
Create valid variable name varname from str.
```

The input str must be a string or a cell array of strings. The output varname will be of the same type.
A valid variable name is a sequence of letters, digits, and underscores that does not begin with a digit.
The "ReplacementStyle" option specifies how invalid characters are handled. Acceptable values are

```
"underscore" (default)
```

Replace all invalid characters with an underscore ("_").
"delete" Remove any invalid character.
"hex" Replace all invalid characters with their hexadecimal representation.
Whitespace characters are always removed prior to the application of the "ReplacementStyle". Lowercase letters following a whitespace will be changed to uppercase.
The "Prefix" option specifies the string pfx to add as a prefix to the input if it begins with a digit. pfx must be a valid variable name itself. The default prefix is "x".

The optional output ismodified is a logical array indicating whether the respective element in str was a valid name or not.

See also: [iskeyword], page 1099, [isvarname], page 137, [matlab.lang.makeUniqueStrings], page 138.
uniqstr = matlab.lang.makeUniqueStrings (str)
uniqstr = matlab.lang.makeUniqueStrings (str, ex)
uniqstr = matlab.lang.makeUniqueStrings (str, ex, maxlength)
[uniqstr, ismodified] = matlab.lang.makeUniqueStrings (...)
Construct a list of unique strings from a list of strings.
The input str must be a string or a cell array of strings. The output uniqstr will be of the same type.
The algorithm makes two strings unique by appending an underscore ("_" and a numeric count to the second string.
If ex is a string or a cell array of strings, uniqstr will contain elements that are unique between themselves and with respect to ex.
If ex is an index array or a logical array for str then it selects the subset of str that are made unique. Unselected elements are not modified.
The optional input maxlength specifies the maximum length of any string in uniqstr. If an input string cannot be made unique without exceeding maxlength an error is emitted.
The optional output ismodified is a logical array indicating whether each element in str was modified to make it unique.

See also: [unique], page 791, [matlab.lang.makeValidName], page 137.
namelengthmax ()
Return the matlab compatible maximum variable name length.

Octave is capable of storing strings up to $2^{31}-1$ in length. However for matlab compatibility all variable, function, and structure field names should be shorter than the length returned by namelengthmax. In particular, variables stored to a matlab file format (*.mat) will have their names truncated to this length.

### 7.1 Global Variables

A global variable is one that may be accessed anywhere within Octave. This is in contrast to a local variable which can only be accessed outside of its current context if it is passed explicitly, such as by including it as a parameter when calling a function (fcn (local_var1, local_var2)).

A variable is declared global by using a global declaration statement. The following statements are all global declarations.

```
global a
global a b
global c = 2
global d = 3 e f = 5
```

Note that the global qualifier extends only to the next end-of-statement indicator which could be a comma (','), semicolon (';'), or newline (' $\backslash \mathrm{n}$ ''). For example, the following code declares one global variable, a, and one local variable $b$ to which the value 1 is assigned.

```
global a, b = 1
```

A global variable may only be initialized once in a global statement. For example, after executing the following code

```
global gvar = 1
global gvar = 2
```

the value of the global variable gvar is 1 , not 2 . Issuing a 'clear gvar' command does not change the above behavior, but 'clear all' does.

It is necessary declare a variable as global within a function body in order to access the one universal variable. For example,

```
global x
function \(f()\)
    \(\mathrm{x}=1\);
endfunction
f ()
```

does not set the value of the global variable x to 1 . Instead, a local variable, with name x , is created and assigned the value of 1 . In order to change the value of the global variable x , you must also declare it to be global within the function body, like this

```
function f ()
    global x;
    x = 1;
endfunction
```

Passing a global variable in a function parameter list will make a local copy and not modify the global value. For example, given the function

```
function f (x)
    \(\mathrm{x}=0\)
endfunction
```

and the definition of x as a global variable at the top level,
global $\mathrm{x}=13$
the expression
f ( x )
will display the value of $x$ from inside the function as 0 , but the value of $x$ at the top level remains unchanged, because the function works with a copy of its argument.

Programming Note: While global variables occasionally are the right solution to a coding problem, modern best practice discourages their use. Code which relies on global variables may behave unpredictably between different users and can be difficult to debug. This is because global variables can introduce systemic changes so that localizing a bug to a particular function, or to a particular loop within a function, becomes difficult.
isglobal (name)
Return true if name is a globally visible variable.
For example:

```
global x
isglobal ("x")
            # 1
```

See also: [isvarname], page 137, [exist], page 144.

### 7.2 Persistent Variables

A variable that has been declared persistent within a function will retain its contents in memory between subsequent calls to the same function. The difference between persistent variables and global variables is that persistent variables are local in scope to a particular function and are not visible elsewhere.

The following example uses a persistent variable to create a function that prints the number of times it has been called.

```
function count_calls ()
    persistent calls = 0;
    printf ("'count_calls' has been called %d times\n",
            ++calls);
endfunction
for i = 1:3
    count_calls ();
endfor
\dashv 'count_calls' has been called 1 times
\dashv'count_calls' has been called 2 times
\dashv'count_calls' has been called 3 times
```

As the example shows, a variable may be declared persistent using a persistent declaration statement. The following statements are all persistent declarations.

```
persistent a
persistent a b
persistent c = 2
persistent d = 3 e f = 5
```

The behavior of persistent variables is equivalent to the behavior of static variables in C.

One restriction for persistent variables is, that neither input nor output arguments of a function can be persistent:

```
function y = foo ()
    persistent y = 0; # Not allowed!
endfunction
foo ()
\dashv error: can't make function parameter y persistent
```

Like global variables, a persistent variable may only be initialized once. For example, after executing the following code

```
persistent pvar = 1
persistent pvar = 2
```

the value of the persistent variable pvar is 1 , not 2 .
If a persistent variable is declared but not initialized to a specific value, it will contain an empty matrix. So, it is also possible to initialize a persistent variable by checking whether it is empty, as the following example illustrates.

```
function count_calls ()
    persistent calls;
    if (isempty (calls))
        calls = 0;
    endif
    printf ("'count_calls' has been called %d times\n",
        ++calls);
endfunction
```

This implementation behaves in exactly the same way as the previous implementation of count_calls.

The value of a persistent variable is kept in memory until it is explicitly cleared. Assuming that the implementation of count_calls is saved on disk, we get the following behavior.

```
for i = 1:2
    count_calls ();
endfor
    \dashv'count_calls' has been called 1 times
    \dashv'count_calls' has been called 2 times
clear
```

```
for i = 1:2
    count_calls ();
endfor
    \dashv 'count_calls' has been called 3 times
    \dashv'count_calls' has been called 4 times
clear all
for i = 1:2
    count_calls ();
endfor
\dashv 'count_calls' has been called 1 times
\dashv'count_calls' has been called 2 times
clear count_calls
for i = 1:2
    count_calls ();
endfor
\dashv'count_calls' has been called 1 times
\dashv'count_calls' has been called 2 times
```

That is, the persistent variable is only removed from memory when the function containing the variable is removed. Note that if the function definition is typed directly into the Octave prompt, the persistent variable will be cleared by a simple clear command as the entire function definition will be removed from memory. If you do not want a persistent variable to be removed from memory even if the function is cleared, you should use the mlock function (see Section 11.10.6 [Function Locking], page 221).

### 7.3 Status of Variables

When creating simple one-shot programs it can be very convenient to see which variables are available at the prompt. The function who and its siblings whos and whos_line_format will show different information about what is in memory, as the following shows.

```
str = "A random string";
who
    \ Variables in the current scope:
    \dashv
    \dashv ans str
```

who
who pattern ...
who option pattern ...
C = who ("pattern", ...)
List currently defined variables matching the given patterns.
Valid pattern syntax is the same as described for the clear command. If no patterns are supplied, all variables are listed.
By default, only variables visible in the local scope are displayed.
The following are valid options, but may not be combined.
global List variables in the global scope rather than the current scope.
-regexp The patterns are considered to be regular expressions when matching the variables to display. The same pattern syntax accepted by the regexp function is used.
-file The next argument is treated as a filename. All variables found within the specified file are listed. No patterns are accepted when reading variables from a file.

If called as a function, return a cell array of defined variable names matching the given patterns.

See also: [whos], page 143, [isglobal], page 140, [isvarname], page 137, [exist], page 144, [regexp], page 91.
whos
whos pattern...
whos option pattern...
S = whos ("pattern", ...)
Provide detailed information on currently defined variables matching the given patterns.
Options and pattern syntax are the same as for the who command.
Extended information about each variable is summarized in a table with the following default entries.

Attr Attributes of the listed variable. Possible attributes are:
blank Variable in local scope
c Variable of complex type.
$\mathrm{f} \quad$ Formal parameter (function argument).
g Variable with global scope.
$\mathrm{p} \quad$ Persistent variable.
Name The name of the variable.
Size The logical size of the variable. A scalar is 1 x 1 , a vector is 1 xN or Nx 1 , a 2-D matrix is MxN.

Bytes The amount of memory currently used to store the variable.
Class The class of the variable. Examples include double, single, char, uint16, cell, and struct.

The table can be customized to display more or less information through the function whos_line_format.
If whos is called as a function, return a struct array of defined variable names matching the given patterns. Fields in the structure describing each variable are: name, size, bytes, class, global, sparse, complex, nesting, persistent.
See also: [who], page 142, [whos_line_format], page 144.
val = whos_line_format ()
old_val = whos_line_format (new_val)
whos_line_format (new_val, "local")
Query or set the format string used by the command whos.
A full format string is:

```
%[modifier]<command>[:width[:left-min[:balance]]];
```

The following command sequences are available:
$\% \mathrm{a} \quad$ Prints attributes of variables (c=complex, $\mathrm{s}=$ sparse, $\mathrm{f}=$ formal parameter, $\mathrm{g}=$ global, $\mathrm{p}=$ persistent).
$\% \mathrm{~b} \quad$ Prints number of bytes occupied by variables.
\%c Prints class names of variables.
\%e Prints elements held by variables.
$\% \mathrm{n} \quad$ Prints variable names.
\%s Prints dimensions of variables.
$\% \mathrm{t} \quad$ Prints type names of variables.
Every command may also have an alignment modifier:
1 Left alignment.
$r \quad$ Right alignment (default).
c Column-aligned (only applicable to command $\% \mathrm{~s}$ ).
The width parameter is a positive integer specifying the minimum number of columns used for printing. No maximum is needed as the field will auto-expand as required.
The parameters left-min and balance are only available when the column-aligned modifier is used with the command '\%s'. balance specifies the column number within the field width which will be aligned between entries. Numbering starts from 0 which indicates the leftmost column. left-min specifies the minimum field width to the left of the specified balance column.

The default format is:

```
" %la:5; %ln:6; %cs:16:6:1; %rb:12; %lc:-1;\n"
```

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [whos], page 143.
Instead of displaying which variables are in memory, it is possible to determine if a given variable is available. That way it is possible to alter the behavior of a program depending on the existence of a variable. The following example illustrates this.

```
if (! exist ("meaning", "var"))
    disp ("The program has no 'meaning'");
endif
```

$c=$ exist (name)
$c=$ exist (name, type)
Check for the existence of name as a variable, function, file, directory, or class.
The return code $c$ is one of
1 name is a variable.
2 name is an absolute filename, an ordinary file in Octave's path, or (after appending '. m ') a function file in Octave's path.
3
5 name is a built-in function.
7 name is a directory.
8 name is a classdef class.
103 name is a function not associated with a file (entered on the command line).
0 name does not exist.
If the optional argument type is supplied, check only for symbols of the specified type. Valid types are
"var" Check only for variables.
"builtin"
Check only for built-in functions.
"dir" Check only for directories.
"file" Check only for files and directories.
"class" Check only for classdef classes.
If no type is given, and there are multiple possible matches for name, exist will return a code according to the following priority list: variable, built-in function, octfile, directory, file, class.
exist returns 2 if a regular file called name is present in Octave's search path. For information about other types of files not on the search path use some combination of the functions file_in_path and stat instead.
Programming Note: If name is implemented by a buggy .oct/.mex file, calling exist may cause Octave to crash. To maintain high performance, Octave trusts .oct/.mex files instead of sandboxing them.

See also: [file_in_loadpath], page 215, [file_in_path], page 960, [dir_in_loadpath], page 216, [stat], page 956.

Usually Octave will manage the memory, but sometimes it can be practical to remove variables from memory manually. This is usually needed when working with large variables that fill a substantial part of the memory. On a computer that uses the IEEE floating point format, the following program allocates a matrix that requires around 128 MB memory.

```
large_matrix = zeros (4000, 4000);
```

Since having this variable in memory might slow down other computations, it can be necessary to remove it manually from memory. The clear or clearvars functions do this.
clear
clear pattern ...
clear options pattern ...
Delete the names matching the given patterns thereby freeing memory.
The pattern may contain the following special characters:
? Match any single character.

* Match zero or more characters.
[ list] Match the list of characters specified by list. If the first character is ! or ^, match all characters except those specified by list. For example, the pattern $[a-z A-Z]$ will match all lowercase and uppercase alphabetic characters.

For example, the command

```
clear foo b*r
```

clears the name foo and all names that begin with the letter ' $b$ ' and end with the letter 'r'.
If clear is called without any arguments, all user-defined variables are cleared from the current workspace (i.e., local variables). Any global variables present will no longer be visible in the current workspace, but they will continue to exist in the global workspace. Functions are unaffected by this form of clear.
The following options are available in both long and short form
all, -all, -a
Clear all local and global user-defined variables, and all functions from the symbol table.
-exclusive, -x
Clear variables that do not match the following pattern.
functions, -functions, -f
Clear function names from the function symbol table. Persistent variables will be re-initialized to their default value unless the function has been locked in memory with mlock.
global, -global, -g
Clear global variable names.
variables, -variables, -v
Clear local variable names.
classes, -classes, -c
Clear the class structure table and all objects.
-regexp, -r
The pattern arguments are treated as regular expressions and any matches will be cleared.
With the exception of -exclusive and -regexp, all long options can be used without the dash as well. Note that, aside from -exclusive, only one other option may appear. All options must appear before any patterns.

Programming Notes: The command clear name only clears the variable name when both a variable and a (shadowed) function named name are currently defined. For example, suppose you have defined a function foo, and then hidden it by performing the assignment foo $=2$. Executing the command clear foo once will clear the variable definition and restore the definition of $f \circ 0$ as a function. Executing clear foo a second time will clear the function definition.
When a local variable name, which is linked to a global variable, is cleared only the local copy of the variable is removed. The global copy is untouched and can be restored with global global_varname. Conversely, clear -g global_varname will remove both the local and global variables.

See also: [clearvars], page 147, [who], page 142, [whos], page 143, [exist], page 144, [mlock], page 222.
clearvars
clearvars pattern...
clearvars -regexp pattern...
clearvars ... -except pattern ...
clearvars . . - except -regexp pattern . . .
clearvars -global...
Delete the variables matching the given patterns from memory.
The pattern may contain the following special characters:
? Match any single character.

* Match zero or more characters.
[ list] Match the list of characters specified by list. If the first character is ! or ^, match all characters except those specified by list. For example, the pattern $[\mathrm{a}-\mathrm{zA}-\mathrm{Z}]$ will match all lowercase and uppercase alphabetic characters.

If the -regexp option is given then subsequent patterns are treated as regular expressions and any matches will be cleared.
If the -except option is given then subsequent patterns select variables that will not be cleared.
If the -global option is given then all patterns will be applied to global variables rather than local variables.
When called with no arguments, clearvars deletes all local variables.
Example Code:
Clear all variables starting with 'x' and the specific variable "foobar"
clearvars $x *$ foobar
Clear the specific variable "foobar" and use regular expressions to clear all variables starting with 'x' or 'y'.

```
clearvars foobar -regexp ^x "y
```

Clear all variables except for "foobar"
clearvars -except foobar

Clear all variables beginning with "foo", except for those ending in "bar"
clearvars foo* -except -regexp bar\$
See also: [clear], page 145, [who], page 142, [whos], page 143, [exist], page 144. pack ()

Consolidate workspace memory in matlab.
This function is provided for compatibility, but does nothing in Octave.
See also: [clear], page 145.
Information about a function or variable such as its location in the file system can also be acquired from within Octave. This is usually only useful during development of programs, and not within a program.
type name...
type -q name...
text = type ("name", ...)
Display the contents of name which may be a file, function (m-file), variable, operator, or keyword.
type normally prepends a header line describing the category of name such as function or variable; The -q option suppresses this behavior.
If no output variable is used the contents are displayed on screen. Otherwise, a cell array of strings is returned, where each element corresponds to the contents of each requested function.
which name...
Display the type of each name.
If name is defined from a function file, the full name of the file is also displayed.
See also: [help], page 21, [lookfor], page 21.
what
what dir
$\mathrm{w}=$ what (dir)
List the Octave specific files in directory dir.
If dir is not specified then the current directory is used.
If a return argument is requested, the files found are returned in the structure $w$. The structure contains the following fields:
path Full path to directory dir
$m \quad$ Cell array of m-files
mat Cell array of mat files
mex Cell array of mex files
oct Cell array of oct files
mdl Cell array of mdl files
slx Cell array of slx files
p Cell array of p-files
classes Cell array of class directories (@classname/)
packages Cell array of package directories (+pkgname/)
Compatibility Note: Octave does not support mdl, slx, and p files. what will always return an empty list for these categories.

See also: [which], page 148, [ls], page 982, [exist], page 144.

## 8 Expressions

Expressions are the basic building block of statements in Octave. An expression evaluates to a value, which you can print, test, store in a variable, pass to a function, or assign a new value to a variable with an assignment operator.

An expression can serve as a statement on its own. Most other kinds of statements contain one or more expressions which specify data to be operated on. As in other languages, expressions in Octave include variables, array references, constants, and function calls, as well as combinations of these with various operators.

### 8.1 Index Expressions

An index expression allows you to reference or extract selected elements of a vector, a matrix (2-D), or a higher-dimensional array.

Indices may be scalars, vectors, ranges, or the special operator ':', which selects entire rows, columns, or higher-dimensional slices.

An index expression consists of a set of parentheses enclosing $M$ expressions separated by commas. Each individual index value, or component, is used for the respective dimension of the object that it is applied to. In other words, the first index component is used for the first dimension (rows) of the object, the second index component is used for the second dimension (columns) of the object, and so on. The number of index components $M$ defines the dimensionality of the index expression. An index with two components would be referred to as a 2-D index because it has two dimensions.

In the simplest case, 1) all components are scalars, and 2) the dimensionality of the index expression $M$ is equal to the dimensionality of the object it is applied to. For example:

```
A = reshape (1:8, 2, 2, 2) # Create 3-D array
A =
ans(:,:,1) =
    1 3
    24
ans(:,:,2) =
    5 7
    6
A(2, 1, 2) # second row, first column of second slice
    # in third dimension: ans = 6
```

The size of the returned object in a specific dimension is equal to the number of elements in the corresponding component of the index expression. When all components are scalars, the result is a single output value. However, if any component is a vector or range then the returned values are the Cartesian product of the indices in the respective dimensions. For example:

```
A([1, 2], 1, 2) \equiv [A(1,1,2); A(2,1,2)]
#
ans =
    5
    6
```

The total number of returned values is the product of the number of elements returned for each index component. In the example above, the total is $2^{*} 1^{*} 1=2$ elements.

Notice that the size of the returned object in a given dimension is equal to the number of elements in the index expression for that dimension. In the code above, the first index component ([1, 2]) was specified as a row vector, but its shape is unimportant. The important fact is that the component specified two values, and hence the result must have a size of two in the first dimension; and because the first dimension corresponds to rows, the overall result is a column vector.

```
A(1, [2, 1, 1], 1) # result is a row vector: ans = [3, 1, 1]
A(ones (2, 2), 1, 1) # result is a column vector: ans = [1; 1; 1; 1]
```

The first line demonstrates again that the size of the output in a given dimension is equal to the number of elements in the respective indexing component. In this case, the output has three elements in the second dimension (which corresponds to columns), so the result is a row vector. The example also shows how repeating entries in the index expression can be used to replicate elements in the output. The last example further proves that the shape of the indexing component is irrelevant, it is only the number of elements $(2 x 2=4)$ which is important.

The above rules apply whenever the dimensionality of the index expression is greater than one $(M>1)$. However, for one-dimensional index expressions special rules apply and the shape of the output is determined by the shape of the indexing component. For example:

```
A([1, 2]) # result is a row vector: ans = [1, 2]
A([1; 2]) # result is a column vector: ans = [1; 2]
```

Note that it is permissible to use a $1-\mathrm{D}$ index with a multi-dimensional object (also called linear indexing). In this case, the elements of the multi-dimensional array are taken in column-first order like Fortran. That is, the columns of the array are imagined to be stacked on top of each other to form a column vector and then the single linear index is applied to this vector.

```
A(5) # linear indexing into three-dimensional array: ans = 5
A(3:5) # result has shape of index component: ans = [3, 4, 5]
```

A colon (':') may be used as an index component to select all of the elements in a specified dimension. Given the matrix,

$$
A=[1,2 ; 3,4]
$$

all of the following expressions are equivalent and select the first row of the matrix.

```
A(1, [1, 2]) # row 1, columns 1 and 2
A(1, 1:2) # row 1, columns in range 1-2
A(1, :) # row 1, all columns
```

When a colon is used in the special case of 1-D indexing the result is always a column vector. Creating column vectors with a colon index is a very frequently encountered code idiom and is faster and generally clearer than calling reshape for this case.

```
A(:) # result is column vector: ans = [1; 2; 3; 4]
A(:)' # result is row vector: ans = [1, 2, 3, 4]
```

In index expressions the keyword end automatically refers to the last entry for a particular dimension. This magic index can also be used in ranges and typically eliminates the needs to call size or length to gather array bounds before indexing. For example:

```
A(1:end/2) \# first half of \(A=>[1,2]\)
A (end +1 ) \(=5 ; \quad \#\) append element
A (end) = []; \# delete element
\(A(1: 2:\) end \() \quad \#\) odd elements of \(A=>[1,3]\)
A(2:2:end) \# even elements of \(A=>[2,4]\)
A (end:-1:1) \# reversal of \(A=>[4,3,2,1]\)
```


### 8.1.1 Advanced Indexing

When it is necessary to extract subsets of entries out of an array whose indices cannot be written as a Cartesian product of components, linear indexing together with the function sub2ind can be used. For example:

```
A = reshape (1:8, 2, 2, 2) # Create 3-D array
A =
ans(:,:,1) =
    1 3
    2 4
ans(:,:,2) =
    5 7
    6
A(sub2ind (size (A), [1, 2, 1], [1, 1, 2], [1, 2, 1]))
    ans = [A(1, 1, 1), A(2, 1, 2), A(1, 2, 1)]
```

An array with 'nd' dimensions can be indexed by an index expression which has from 1 to 'nd' components. For the ordinary and most common case, the number of components ' M ' matches the number of dimensions ' $n d$ '. In this case the ordinary indexing rules apply and each component corresponds to the respective dimension of the array.

However, if the number of indexing components exceeds the number of dimensions ( $M>n d$ ) then the excess components must all be singletons (1). Moreover, if $M<n d$, the behavior is equivalent to reshaping the input object so as to merge the trailing nd -M dimensions into the last index dimension M. Thus, the result will have the dimensionality of the index expression, and not the original object. This is the case whenever dimensionality of the index is greater than one ( $\mathrm{M}>1$ ), so that the special rules for linear indexing are not applied. This is easiest to understand with an example:

```
A = reshape (1:8, 2, 2, 2) # Create 3-D array
A =
```

```
ans(:,:,1) =
    1 3
    24
ans(:,:,2) =
    5 7
    6
## 2-D indexing causes third dimension to be merged into second dimension.
## Equivalent array for indexing, Atmp, is now 2x4.
Atmp = reshape (A, 2, 4)
Atmp =
    1 3 5 7
    4 6 8
```

```
A(2,1) # Reshape to 2x4 matrix, second entry of first column: ans = 2
A(2,4) # Reshape to 2x4 matrix, second entry of fourth column: ans = 8
A(:,:) # Reshape to 2x4 matrix, select all rows & columns, ans = Atmp
```

Note here the elegant use of the double colon to replace the call to the reshape function.
Another advanced use of linear indexing is to create arrays filled with a single value. This can be done by using an index of ones on a scalar value. The result is an object with the dimensions of the index expression and every element equal to the original scalar. For example, the following statements

```
a = 13;
a(ones (1, 4))
```

produce a row vector whose four elements are all equal to 13 .
Similarly, by indexing a scalar with two vectors of ones it is possible to create a matrix. The following statements

```
a = 13;
a(ones (1, 2), ones (1, 3))
```

create a $2 \times 3$ matrix with all elements equal to 13 . This could also have been written as
13(ones (2, 3))
It is more efficient to use indexing rather than the code construction scalar * ones ( M , N, ...) because it avoids the unnecessary multiplication operation. Moreover, multiplication may not be defined for the object to be replicated whereas indexing an array is always defined. The following code shows how to create a $2 \times 3$ cell array from a base unit which is not itself a scalar.

```
{"Hello"}(ones (2, 3))
```

It should be noted that ones (1, n) (a row vector of ones) results in a range object (with zero increment). A range is stored internally as a starting value, increment, end value, and
total number of values; hence, it is more efficient for storage than a vector or matrix of ones whenever the number of elements is greater than 4 . In particular, when ' $r$ ' is a row vector, the expressions

```
r(ones (1, n), :)
r(ones (n, 1), :)
```

will produce identical results, but the first one will be significantly faster, at least for ' $r$ ' and ' $n$ ' large enough. In the first case the index is held in compressed form as a range which allows Octave to choose a more efficient algorithm to handle the expression.

A general recommendation for users unfamiliar with these techniques is to use the function repmat for replicating smaller arrays into bigger ones, which uses such tricks.

A second use of indexing is to speed up code. Indexing is a fast operation and judicious use of it can reduce the requirement for looping over individual array elements, which is a slow operation.

Consider the following example which creates a 10 -element row vector $a$ containing the values $a_{i}=\sqrt{i}$.

```
for i = 1:10
    a(i) = sqrt (i);
endfor
```

It is quite inefficient to create a vector using a loop like this. In this case, it would have been much more efficient to use the expression

```
a = sqrt (1:10);
```

which avoids the loop entirely.
In cases where a loop cannot be avoided, or a number of values must be combined to form a larger matrix, it is generally faster to set the size of the matrix first (pre-allocate storage), and then insert elements using indexing commands. For example, given a matrix a,

```
[nr, nc] = size (a);
x = zeros (nr, n * nc);
for i = 1:n
    x(:,(i-1)*nc+1:i*nc) = a;
endfor
```

is considerably faster than

```
x = a;
for i = 1:n-1
    x = [x, a];
endfor
```

because Octave does not have to repeatedly resize the intermediate result.

```
ind = sub2ind (dims, i, j)
ind = sub2ind (dims, s1, s2, ..., sN)
```

Convert subscripts to linear indices.
The input dims is a dimension vector where each element is the size of the array in the respective dimension (see [size], page 47). The remaining inputs are scalars or vectors of subscripts to be converted.

The output vector ind contains the converted linear indices.
Background: Array elements can be specified either by a linear index which starts at 1 and runs through the number of elements in the array, or they may be specified with subscripts for the row, column, page, etc. The functions ind2sub and sub2ind interconvert between the two forms.
The linear index traverses dimension 1 (rows), then dimension 2 (columns), then dimension 3 (pages), etc. until it has numbered all of the elements. Consider the following 3 -by- 3 matrices:


The left matrix contains the subscript tuples for each matrix element. The right matrix shows the linear indices for the same matrix.
The following example shows how to convert the two-dimensional indices $(2,1)$ and $(2,3)$ of a 3 -by- 3 matrix to linear indices with a single call to sub2ind.

```
s1 = [2, 2];
s2 = [1, 3];
ind = sub2ind ([3, 3], s1, s2)
    # ind = 2 8
```

See also: [ind2sub], page 156, [size], page 47.
[s1, s2, ..., sN] = ind2sub (dims, ind)
Convert linear indices to subscripts.
The input dims is a dimension vector where each element is the size of the array in the respective dimension (see [size], page 47). The second input ind contains linear indies to be converted.
The outputs s1, ..., sN contain the converted subscripts.
Background: Array elements can be specified either by a linear index which starts at 1 and runs through the number of elements in the array, or they may be specified with subscripts for the row, column, page, etc. The functions ind2sub and sub2ind interconvert between the two forms.
The linear index traverses dimension 1 (rows), then dimension 2 (columns), then dimension 3 (pages), etc. until it has numbered all of the elements. Consider the following 3 -by- 3 matrices:

```
[1, 4, 7] [(1,1), (1,2), (1,3)]
[2, 5, 8] ==> [(2,1), (2,2), (2,3)]
[3, 6, 9] [(3,1), (3,2), (3,3)]
```

The left matrix contains the linear indices for each matrix element. The right matrix shows the subscript tuples for the same matrix.
The following example shows how to convert the two-dimensional indices $(2,1)$ and $(2,3)$ of a 3 -by- 3 matrix to linear indices with a single call to sub2ind.
The following example shows how to convert the linear indices 2 and 8 in a 3 -by- 3 matrix into subscripts.

```
ind = [2, 8];
[r, c] = ind2sub ([3, 3], ind)
    m = 2 2
    c}=1
```

If the number of output subscripts exceeds the number of dimensions, the exceeded dimensions are set to 1 . On the other hand, if fewer subscripts than dimensions are provided, the exceeding dimensions are merged into the final requested dimension. For clarity, consider the following examples:

```
ind = [2, 8];
dims = [3, 3];
## same as dims = [3, 3, 1]
[r, c, s] = ind2sub (dims, ind)
    r = 2 2
    "c=1 3
    m = 1 1
## same as dims = [9]
r = ind2sub (dims, ind)
    r = 2 8
```

See also: [sub2ind], page 155, [size], page 47.
isindex (ind)
isindex (ind, n)
Return true if ind is a valid index.
Valid indices are either positive integers (although possibly of real data type), or logical arrays.

If present, $n$ specifies the maximum extent of the dimension to be indexed. When possible the internal result is cached so that subsequent indexing using ind will not perform the check again.

Implementation Note: Strings are first converted to double values before the checks for valid indices are made. Unless a string contains the NULL character " $\backslash 0$ ", it will always be a valid index.

### 8.2 Calling Functions

A function is a name for a particular calculation. Because it has a name, you can ask for it by name at any point in the program. For example, the function sqrt computes the square root of a number.

A fixed set of functions are built-in, which means they are available in every Octave program. The sqrt function is one of these. In addition, you can define your own functions. See Chapter 11 [Functions and Scripts], page 191, for information about how to do this.

The way to use a function is with a function call expression, which consists of the function name followed by a list of arguments in parentheses. The arguments are expressions which give the raw materials for the calculation that the function will do. When there is more than one argument, they are separated by commas. If there are no arguments, you can
omit the parentheses, but it is a good idea to include them anyway, to clearly indicate that a function call was intended. Here are some examples:

```
sqrt (x^2 + y^2) # One argument
ones (n, m) # Two arguments
rand () # No arguments
```

Each function expects a particular number of arguments. For example, the sqrt function must be called with a single argument, the number to take the square root of:

```
sqrt (argument)
```

Some of the built-in functions take a variable number of arguments, depending on the particular usage, and their behavior is different depending on the number of arguments supplied.

Like every other expression, the function call has a value, which is computed by the function based on the arguments you give it. In this example, the value of sqrt (argument) is the square root of the argument. A function can also have side effects, such as assigning the values of certain variables or doing input or output operations.

Unlike most languages, functions in Octave may return multiple values. For example, the following statement

```
[u, s, v] = svd (a)
```

computes the singular value decomposition of the matrix a and assigns the three result matrices to $\mathrm{u}, \mathrm{s}$, and v .

The left side of a multiple assignment expression is itself a list of expressions, that is, a list of variable names potentially qualified by index expressions. See also Section 8.1 [Index Expressions], page 151, and Section 8.6 [Assignment Ops], page 169.

### 8.2.1 Call by Value

In Octave, unlike Fortran, function arguments are passed by value, which means that each argument in a function call is evaluated and assigned to a temporary location in memory before being passed to the function. There is currently no way to specify that a function parameter should be passed by reference instead of by value. This means that it is impossible to directly alter the value of a function parameter in the calling function. It can only change the local copy within the function body. For example, the function

```
function f (x, n)
    while (n-- > 0)
        disp (x);
    endwhile
endfunction
```

displays the value of the first argument $n$ times. In this function, the variable $n$ is used as a temporary variable without having to worry that its value might also change in the calling function. Call by value is also useful because it is always possible to pass constants for any function parameter without first having to determine that the function will not attempt to modify the parameter.

The caller may use a variable as the expression for the argument, but the called function does not know this: it only knows what value the argument had. For example, given a function called as

```
foo = "bar";
fcn (foo)
```

you should not think of the argument as being "the variable foo." Instead, think of the argument as the string value, "bar".

Even though Octave uses pass-by-value semantics for function arguments, values are not copied unnecessarily. For example,

```
x = rand (1000);
f (x);
```

does not actually force two 1000 by 1000 element matrices to exist unless the function $f$ modifies the value of its argument. Then Octave must create a copy to avoid changing the value outside the scope of the function $f$, or attempting (and probably failing!) to modify the value of a constant or the value of a temporary result.

### 8.2.2 Recursion

With some restrictions ${ }^{1}$, recursive function calls are allowed. A recursive function is one which calls itself, either directly or indirectly. For example, here is an inefficient ${ }^{2}$ way to compute the factorial of a given integer:

```
function retval \(=\) fact ( \(n\) )
    if ( \(\mathrm{n}>0\) )
        retval \(=\mathrm{n} *\) fact ( \(\mathrm{n}-1\) );
    else
        retval = 1;
    endif
endfunction
```

This function is recursive because it calls itself directly. It eventually terminates because each time it calls itself, it uses an argument that is one less than was used for the previous call. Once the argument is no longer greater than zero, it does not call itself, and the recursion ends.

The function max_recursion_depth may be used to specify a limit to the recursion depth and prevents Octave from recursing infinitely. Similarly, the function max_stack_ depth may be used to specify limit to the depth of function calls, whether recursive or not. These limits help prevent stack overflow on the computer Octave is running on, so that instead of exiting with a signal, the interpreter will throw an error and return to the command prompt.

```
val = max_recursion_depth ()
old_val = max_recursion_depth (new_val)
max_recursion_depth (new_val, "local")
```

Query or set the internal limit on the number of times a function may be called recursively.

[^2]If the limit is exceeded, an error message is printed and control returns to the top level.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [max_stack_depth], page 160.

```
val = max_stack_depth ()
old_val = max_stack_depth (new_val)
max_stack_depth (new_val, "local")
```

Query or set the internal limit on the number of times a function may be called recursively.

If the limit is exceeded, an error message is printed and control returns to the top level.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [max_recursion_depth], page 159.

### 8.2.3 Access via Handle

A function may be abstracted and referenced via a function handle acquired using the special operator ' $₫$ '. For example,

```
f = @plus;
f (2, 2)
# 4
```

is equivalent to calling plus $(2,2)$ directly. Beyond abstraction for general programming, function handles find use in callback methods for figures and graphics by adding listeners to properties or assigning pre-existing actions, such as in the following example:

```
function mydeletefcn (h, ~ , msg)
    printf (msg);
endfunction
sombrero;
set (gcf, "deletefcn", {@mydeletefcn, "Bye!\n"});
close;
```

The above will print "Bye!" to the terminal upon the closing (deleting) of the figure. There are many graphics property actions for which a callback function may be assigned, including, buttondownfcn, windowscrollwheelfcn, createfcn, deletefcn, keypressfcn, etc.

Note that the ' $@$ ' character also plays a role in defining class functions, i.e., methods, but not as a syntactical element. Rather it begins a directory name containing methods for a class that shares the directory name sans the '@' character. See Chapter 34 [Object Oriented Programming], page 889.

### 8.3 Arithmetic Operators

The following arithmetic operators are available, and work on scalars and matrices. The element-by-element operators and functions broadcast (see Section 19.2 [Broadcasting], page 627).
$x+y \quad$ Addition (always works element by element). If both operands are matrices, the number of rows and columns must both agree, or they must be broadcastable to the same shape.
$x-y \quad$ Subtraction (always works element by element). If both operands are matrices, the number of rows and columns of both must agree, or they must be broadcastable to the same shape.
$x * y \quad$ Matrix multiplication. The number of columns of $x$ must agree with the number of rows of $y$.
$x$.* $y$ Element-by-element multiplication. If both operands are matrices, the number of rows and columns must both agree, or they must be broadcastable to the same shape.
$x / y \quad$ Right division. This is conceptually equivalent to the expression

```
(inv (y') * x')'
```

but it is computed without forming the inverse of $y^{\prime}$ '.
If the system is not square, or if the coefficient matrix is singular, a minimum norm solution is computed.
$x$./ y Element-by-element right division.
$x \backslash y \quad$ Left division. This is conceptually equivalent to the expression
inv (x) * y
but it is computed without forming the inverse of $x$.
If the system is not square, or if the coefficient matrix is singular, a minimum norm solution is computed.
$x$. $\$ y Element-by-element left division. Each element of $y$ is divided by each corresponding element of $x$.
$x^{\wedge} y \quad$ Power operator. If $x$ and $y$ are both scalars, this operator returns $x$ raised to the power $y$. If $x$ is a scalar and $y$ is a square matrix, the result is computed using an eigenvalue expansion. If $x$ is a square matrix, the result is computed by repeated multiplication if $y$ is an integer, and by an eigenvalue expansion if $y$ is not an integer. An error results if both $x$ and $y$ are matrices.
The implementation of this operator needs to be improved.
$x$.^y Element-by-element power operator. If both operands are matrices, the number of rows and columns must both agree, or they must be broadcastable to the same shape. If several complex results are possible, the one with smallest nonnegative argument (angle) is taken. This rule may return a complex root even when a real root is also possible. Use realpow, realsqrt, cbrt, or nthroot if a real result is preferred.
$-x \quad$ Negation.
$+x \quad$ Unary plus. This operator has no effect on the operand.
$x^{\prime} \quad$ Complex conjugate transpose. For real arguments, this operator is the same as the transpose operator. For complex arguments, this operator is equivalent to the expression
conj (x.')
x.' Transpose.

Note that because Octave's element-by-element operators begin with a '.', there is a possible ambiguity for statements like
$1 . / m$
because the period could be interpreted either as part of the constant or as part of the operator. To resolve this conflict, Octave treats the expression as if you had typed
(1) . $/ \mathrm{m}$
and not
(1.) / m

Although this is inconsistent with the normal behavior of Octave's lexer, which usually prefers to break the input into tokens by preferring the longest possible match at any given point, it is more useful in this case.
ctranspose (x)
Return the complex conjugate transpose of $x$.
This function and $x^{\prime}$ are equivalent.
See also: [transpose], page 164 .
ldivide (x, y)
Return the element-by-element left division of $x$ and $y$.
This function and $x . \backslash y$ are equivalent.
See also: [rdivide], page 163, [mldivide], page 162, [times], page 164, [plus], page 163.
minus ( $x, y$ )
This function and $x-y$ are equivalent.
See also: [plus], page 163, [uminus], page 164.
mldivide ( $x, y$ )
Return the matrix left division of $x$ and $y$.
This function and $x \backslash y$ are equivalent.
If the system is not square, or if the coefficient matrix is singular, a minimum norm solution is computed.
See also: [mrdivide], page 163, [ldivide], page 162, [rdivide], page 163, [linsolve], page 590 .
mpower ( $x, y$ )
Return the matrix power operation of $x$ raised to the $y$ power.
This function and $x^{\wedge} y$ are equivalent.
See also: [power], page 163, [mtimes], page 163, [plus], page 163, [minus], page 162.
mrdivide ( $x, y$ )
Return the matrix right division of $x$ and $y$.
This function and $x / y$ are equivalent.
If the system is not square, or if the coefficient matrix is singular, a minimum norm solution is computed.

See also: [mldivide], page 162, [rdivide], page 163, [plus], page 163, [minus], page 162.
mtimes ( $x, y$ )
mtimes (x1, x2, ...)
Return the matrix multiplication product of inputs.
This function and $x * y$ are equivalent. If more arguments are given, the multiplication is applied cumulatively from left to right:

$$
(\ldots((x 1 * x 2) * x 3) * \ldots)
$$

See also: [times], page 164, [plus], page 163, [minus], page 162, [rdivide], page 163, [mrdivide], page 163, [mldivide], page 162, [mpower], page 163.
plus ( $x, y$ )
plus (x1, x2, ...)
This function and $x+y$ are equivalent.
If more arguments are given, the summation is applied cumulatively from left to right:

$$
(\ldots((x 1+x 2)+x 3)+\ldots)
$$

See also: [minus], page 162, [uplus], page 164.
power ( $x, y$ )
Return the element-by-element operation of $x$ raised to the $y$ power.
This function and $\mathrm{x} .{ }^{\text {^ } \mathrm{y} \text { are equivalent. }}$
If several complex results are possible, returns the one with smallest non-negative argument (angle). Use realpow, realsqrt, cbrt, or nthroot if a real result is preferred.

See also: [mpower], page 163, [realpow], page 548, [realsqrt], page 548, [cbrt], page 548, [nthroot], page 548.
rdivide (x, y)
Return the element-by-element right division of $x$ and $y$.
This function and $x$./ y are equivalent.
See also: [ldivide], page 162, [mrdivide], page 163, [times], page 164, [plus], page 163.
times ( $x, y$ )
times (x1, x2, ...)
Return the element-by-element multiplication product of inputs.
This function and $x . * y$ are equivalent. If more arguments are given, the multiplication is applied cumulatively from left to right:

$$
(\ldots((x 1 . * x 2) . * x 3) . * \ldots)
$$

See also: [mtimes], page 163, [rdivide], page 163.
transpose (x)
Return the transpose of $x$.
This function and $x .{ }^{\prime}$ are equivalent.
See also: [ctranspose], page 162.
uminus ( $x$ )
This function and $-x$ are equivalent.
See also: [uplus], page 164, [minus], page 162.
uplus (x)
This function and $+x$ are equivalent.
See also: [uminus], page 164, [plus], page 163.

### 8.4 Comparison Operators

Comparison operators compare numeric values for relationships such as equality. They are written using relational operators.

All of Octave's comparison operators return a value of 1 if the comparison is true, or 0 if it is false. For matrix values, they all work on an element-by-element basis. Broadcasting rules apply. See Section 19.2 [Broadcasting], page 627. For example:

```
[1, 2; 3, 4] == [1, 3; 2, 4]
        # 1 0
    0 1
```

According to broadcasting rules, if one operand is a scalar and the other is a matrix, the scalar is compared to each element of the matrix in turn, and the result is the same size as the matrix.

```
x<y True if x is less than y.
x<= y True if x is less than or equal to y.
x== y True if }x\mathrm{ is equal to y.
x >= y True if x is greater than or equal to y.
x>y True if x is greater than y.
x != y
x ~}=y\quad\mathrm{ True if }x\mathrm{ is not equal to }y\mathrm{ .
```

For complex numbers, the following ordering is defined: $z 1<z 2$ if and only if
abs (z1) < abs (z2)
|| (abs (z1) == abs (z2) \&\& $\arg (z 1)<\arg (z 2)$ )
This is consistent with the ordering used by max, min and sort, but is not consistent with MATLAB, which only compares the real parts.

String comparisons may also be performed with the strcmp function, not with the comparison operators listed above. See Chapter 5 [Strings], page 71.
eq ( $x, y$ )
Return true if the two inputs are equal.
This function is equivalent to $x==y$.
See also: [ne], page 165, [isequal], page 165, [le], page 165, [ge], page 165, [gt], page 165, [ne], page 165, [lt], page 165.
ge ( $x, y$ )
This function is equivalent to $x>=y$.
See also: [le], page 165, [eq], page 165, [gt], page 165, [ne], page 165, [lt], page 165.
gt ( $x, y$ )
This function is equivalent to $x>y$.
See also: [le], page 165, [eq], page 165, [ge], page 165, [ne], page 165, [lt], page 165.
isequal (x1, x2, ...)
Return true if all of $x 1, x 2, \ldots$ are equal.
See also: [isequaln], page 165.
isequaln ( $x 1, x 2, \ldots$ )
Return true if all of $x 1, x 2, \ldots$ are equal under the additional assumption that NaN $==\mathrm{NaN}$ (no comparison of NaN placeholders in dataset).
See also: [isequal], page 165 .
le ( $x, y$ )
This function is equivalent to $x<=y$.
See also: [eq], page 165, [ge], page 165, [gt], page 165, [ne], page 165, [lt], page 165.
lt ( $x, y$ )
This function is equivalent to $x<y$.
See also: [le], page 165, [eq], page 165, [ge], page 165, [gt], page 165, [ne], page 165.
ne ( $x, y$ )
Return true if the two inputs are not equal.
This function is equivalent to $x!=y$.
See also: [eq], page 165, [isequal], page 165, [le], page 165, [ge], page 165, [lt], page 165.

### 8.5 Boolean Expressions

### 8.5.1 Element-by-element Boolean Operators

An element-by-element boolean expression is a combination of comparison expressions using the boolean operators "or" (' $I$ '), "and" (' $\&$ '), and "not" ('!'), along with parentheses to control nesting. The truth of the boolean expression is computed by combining the truth values of the corresponding elements of the component expressions. A value is considered to be false if it is zero, and true otherwise.

Element-by-element boolean expressions can be used wherever comparison expressions can be used. They can be used in if and while statements. However, a matrix value used as the condition in an if or while statement is only true if all of its elements are nonzero.

Like comparison operations, each element of an element-by-element boolean expression also has a numeric value ( 1 if true, 0 if false) that comes into play if the result of the boolean expression is stored in a variable, or used in arithmetic.

Here are descriptions of the three element-by-element boolean operators.

## boolean1 \& boolean2

Elements of the result are true if both corresponding elements of boolean1 and boolean2 are true.
boolean1 | boolean2
Elements of the result are true if either of the corresponding elements of boolean1 or boolean2 is true.
! boolean
~ boolean Each element of the result is true if the corresponding element of boolean is false.

These operators work on an element-by-element basis. For example, the expression
$[1,0 ; 0,1] \&[1,0 ; 2,3]$
returns a two by two identity matrix.
For the binary operators, broadcasting rules apply. See Section 19.2 [Broadcasting], page 627. In particular, if one of the operands is a scalar and the other a matrix, the operator is applied to the scalar and each element of the matrix.

For the binary element-by-element boolean operators, both subexpressions boolean1 and boolean2 are evaluated before computing the result. This can make a difference when the expressions have side effects. For example, in the expression

## a \& b++

the value of the variable $b$ is incremented even if the variable $a$ is zero.
This behavior is necessary for the boolean operators to work as described for matrixvalued operands.

```
\(z=\) and ( \(x, y\) )
\(z=\) and (x1, x2, ..)
    Return the logical AND of \(x\) and \(y\).
```

This function is equivalent to the operator syntax $x \& y$. If more than two arguments are given, the logical AND is applied cumulatively from left to right:
(...((x1 \& x2) \& x3) \& ...)

See also: [or], page 167, [not], page 167, [xor], page 511.
$z=\operatorname{not}(x)$
Return the logical NOT of $x$.
This function is equivalent to the operator syntax ! x.
See also: [and], page 166, [or], page 167, [xor], page 511.
$z=$ or ( $x, y$ )
$z=$ or (x1, x2, ...)
Return the logical OR of $x$ and $y$.
This function is equivalent to the operator syntax $x \mid y$. If more than two arguments are given, the logical OR is applied cumulatively from left to right:

$$
(\ldots((x 1 \mid x 2) \mid x 3) \mid \ldots)
$$

See also: [and], page 166, [not], page 167, [xor], page 511.

### 8.5.2 Short-circuit Boolean Operators

Combined with the implicit conversion to scalar values in if and while conditions, Octave's element-by-element boolean operators are often sufficient for performing most logical operations. However, it is sometimes desirable to stop evaluating a boolean expression as soon as the overall truth value can be determined. Octave's short-circuit boolean operators work this way.

## boolean1 \&\& boolean2

The expression boolean1 is evaluated and converted to a scalar using the equivalent of the operation all (boolean1(:)). If it is false, the result of the overall expression is 0 . If it is true, the expression boolean2 is evaluated and converted to a scalar using the equivalent of the operation all (boolean2(:)). If it is true, the result of the overall expression is 1 . Otherwise, the result of the overall expression is 0 .
Warning: there is one exception to the rule of evaluating all (boolean1(:)), which is when boolean1 is the empty matrix. The truth value of an empty matrix is always false so [] \&\& true evaluates to false even though all ([]) is true.
boolean1 || boolean2
The expression boolean1 is evaluated and converted to a scalar using the equivalent of the operation all (boolean1(:)). If it is true, the result of the overall expression is 1 . If it is false, the expression boolean2 is evaluated and converted to a scalar using the equivalent of the operation all (boolean2(:)). If it is true, the result of the overall expression is 1 . Otherwise, the result of the overall expression is 0 .
Warning: the truth value of an empty matrix is always false, see the previous list item for details.

The fact that both operands may not be evaluated before determining the overall truth value of the expression can be important. For example, in the expression
a \&\& b++
the value of the variable $b$ is only incremented if the variable a is nonzero.
This can be used to write somewhat more concise code. For example, it is possible write

```
function f (a, b, c)
    if (nargin > 2 && ischar (c))
        ...
```

instead of having to use two if statements to avoid attempting to evaluate an argument that doesn't exist. For example, without the short-circuit feature, it would be necessary to write

```
function f (a, b, c)
    if (nargin > 2)
        if (ischar (c))
```

Writing
function $f(a, b, c)$
if (nargin > 2 \& ischar (c))
...
would result in an error if $f$ were called with one or two arguments because Octave would be forced to try to evaluate both of the operands for the operator ' $\varepsilon$ '.
matlab has special behavior that allows the operators ' $\varepsilon$ ' and ' $I$ ' to short-circuit when used in the truth expression for if and while statements. Octave behaves the same way for compatibility, however, the use of the ' $\&$ ' and ' $l$ ' operators in this way is strongly discouraged and a warning will be issued. Instead, you should use the ' $\& \&$ ' and ' $\|$ ' operators that always have short-circuit behavior.

Finally, the ternary operator (?:) is not supported in Octave. If short-circuiting is not important, it can be replaced by the ifelse function.

```
merge (mask, tval, fval)
```

ifelse (mask, tval, fval)
Merge elements of true_val and false_val, depending on the value of mask.
If mask is a logical scalar, the other two arguments can be arbitrary values. Otherwise, mask must be a logical array, and tval, fval should be arrays of matching class, or cell arrays. In the scalar mask case, tval is returned if mask is true, otherwise fval is returned.
In the array mask case, both tval and fval must be either scalars or arrays with dimensions equal to mask. The result is constructed as follows:

```
result(mask) = tval(mask);
result(! mask) = fval(! mask);
```

mask can also be arbitrary numeric type, in which case it is first converted to logical.
See also: [logical], page 64, [diff], page 512.

### 8.6 Assignment Expressions

An assignment is an expression that stores a new value into a variable. For example, the following expression assigns the value 1 to the variable $\mathbf{z}$ :

```
z = 1
```

After this expression is executed, the variable $\mathbf{z}$ has the value 1 . Whatever old value $\mathbf{z}$ had before the assignment is forgotten. The ' $=$ ' sign is called an assignment operator.

Assignments can store string values also. For example, the following expression would store the value "this food is good" in the variable message:

```
thing = "food"
predicate = "good"
message = [ "this " , thing , " is " , predicate ]
```

(This also illustrates concatenation of strings.)
Most operators (addition, concatenation, and so on) have no effect except to compute a value. If you ignore the value, you might as well not use the operator. An assignment operator is different. It does produce a value, but even if you ignore the value, the assignment still makes itself felt through the alteration of the variable. We call this a side effect.

The left-hand operand of an assignment need not be a variable (see Chapter 7 [Variables], page 137). It can also be an element of a matrix (see Section 8.1 [Index Expressions], page 151) or a list of return values (see Section 8.2 [Calling Functions], page 157). These are all called lvalues, which means they can appear on the left-hand side of an assignment operator. The right-hand operand may be any expression. It produces the new value which the assignment stores in the specified variable, matrix element, or list of return values.

It is important to note that variables do not have permanent types. The type of a variable is simply the type of whatever value it happens to hold at the moment. In the following program fragment, the variable foo has a numeric value at first, and a string value later on:

```
octave:13> foo = 1
foo = 1
octave:13> foo = "bar"
foo = bar
```

When the second assignment gives foo a string value, the fact that it previously had a numeric value is forgotten.

Assignment of a scalar to an indexed matrix sets all of the elements that are referenced by the indices to the scalar value. For example, if a is a matrix with at least two columns,

$$
a(:, 2)=5
$$

sets all the elements in the second column of a to 5 .
Assigning an empty matrix '[]' works in most cases to allow you to delete rows or columns of matrices and vectors. See Section 4.1.1 [Empty Matrices], page 55. For example, given a 4 by 5 matrix $A$, the assignment

A (3, :) = []
deletes the third row of $A$, and the assignment
A (:, 1:2:5) = []
deletes the first, third, and fifth columns.
An assignment is an expression, so it has a value. Thus, $\mathbf{z}=1$ as an expression has the value 1 . One consequence of this is that you can write multiple assignments together:

$$
x=y=z=0
$$

stores the value 0 in all three variables. It does this because the value of $\mathbf{z}=0$, which is 0 , is stored into $y$, and then the value of $y=z=0$, which is 0 , is stored into $x$.

This is also true of assignments to lists of values, so the following is a valid expression

$$
[\mathrm{a}, \mathrm{~b}, \mathrm{c}]=[\mathrm{u}, \mathrm{~s}, \mathrm{v}]=\operatorname{svd}(\mathrm{a})
$$

that is exactly equivalent to

$$
\begin{aligned}
& {[u, s, v]=s v d \text { (a) }} \\
& a=u \\
& b=s \\
& c=v
\end{aligned}
$$

In expressions like this, the number of values in each part of the expression need not match. For example, the expression

```
[a, b] = [u, s, v] = svd (a)
```

is equivalent to
[u, s, v] = svd (a)
$\mathrm{a}=\mathrm{u}$
$b=s$
The number of values on the left side of the expression can, however, not exceed the number of values on the right side. For example, the following will produce an error.

```
[a, b, c, d] = [u, s, v] = svd (a);
\dashv \text { error: element number 4 undefined in return list}
```

The symbol ~ may be used as a placeholder in the list of lvalues, indicating that the corresponding return value should be ignored and not stored anywhere:

```
[~, s, v] = svd (a);
```

This is cleaner and more memory efficient than using a dummy variable. The nargout value for the right-hand side expression is not affected. If the assignment is used as an expression, the return value is a comma-separated list with the ignored values dropped.

A very common programming pattern is to increment an existing variable with a given value, like this

$$
a=a+2 ;
$$

This can be written in a clearer and more condensed form using the += operator

$$
\text { a += } 2 ;
$$

Similar operators also exist for subtraction ( $-=$ ), multiplication ( $*=$ ), and division (/=). An expression of the form

```
expr1 op= expr2
```

is evaluated as

```
expr1 = (expr1) op (expr2)
```

where $o p$ can be either,,$+- *$, or $/$, as long as expr2 is a simple expression with no side effects. If expr2 also contains an assignment operator, then this expression is evaluated as

```
temp = expr2
expr1 = (expr1) op temp
```

where temp is a placeholder temporary value storing the computed result of evaluating expr2. So, the expression
a $*=\mathrm{b}+1$
is evaluated as
$\mathrm{a}=\mathrm{a} *(\mathrm{~b}+1)$
and not
$\mathrm{a}=\mathrm{a} * \mathrm{~b}+1$
You can use an assignment anywhere an expression is called for. For example, it is valid to write $\mathrm{x}!=(\mathrm{y}=1)$ to set y to 1 and then test whether x equals 1 . But this style tends to make programs hard to read. Except in a one-shot program, you should rewrite it to get rid of such nesting of assignments. This is never very hard.

### 8.7 Increment Operators

Increment operators increase or decrease the value of a variable by 1. The operator to increment a variable is written as ' ++ '. It may be used to increment a variable either before or after taking its value.

For example, to pre-increment the variable $x$, you would write $++x$. This would add one to $x$ and then return the new value of $x$ as the result of the expression. It is exactly the same as the expression $x=x+1$.

To post-increment a variable $x$, you would write $x^{++}$. This adds one to the variable $x$, but returns the value that $x$ had prior to incrementing it. For example, if $x$ is equal to 2 , the result of the expression $x^{++}$is 2 , and the new value of $x$ is 3 .

For matrix and vector arguments, the increment and decrement operators work on each element of the operand.

The increment and decrement operators must "hug" their corresponding variable. That means, no white spaces are allowed between these operators and the variable they affect.

Here is a list of all the increment and decrement expressions.
$++x \quad$ This expression increments the variable $x$. The value of the expression is the new value of $x$. It is equivalent to the expression $x=x+1$.
$--x \quad$ This expression decrements the variable $x$. The value of the expression is the new value of x . It is equivalent to the expression $\mathrm{x}=\mathrm{x}-1$.
$x^{++} \quad$ This expression causes the variable $x$ to be incremented. The value of the expression is the old value of $x$.
x-- This expression causes the variable $x$ to be decremented. The value of the expression is the old value of $x$.

### 8.8 Operator Precedence

Operator precedence determines how operators are grouped, when different operators appear close by in one expression. For example, '*' has higher precedence than ' + '. Thus, the expression $\mathrm{a}+\mathrm{b} * \mathrm{c}$ means to multiply b and c , and then add a to the product (i.e., $\mathrm{a}+\mathrm{b}$ * c) ).

You can overrule the precedence of the operators by using parentheses. You can think of the precedence rules as saying where the parentheses are assumed if you do not write parentheses yourself. In fact, it is wise to use parentheses whenever you have an unusual combination of operators, because other people who read the program may not remember what the precedence is in this case. You might forget as well, and then you too could make a mistake. Explicit parentheses will help prevent any such mistake.

When operators of equal precedence are used together, the leftmost operator groups first, except for the assignment operators, which group in the opposite order. Thus, the expression $\mathrm{a}-\mathrm{b}+\mathrm{c}$ groups as $(\mathrm{a}-\mathrm{b})+\mathrm{c}$, but the expression $\mathrm{a}=\mathrm{b}=\mathrm{c}$ groups as $\mathrm{a}=(\mathrm{b}=$ c).

The precedence of prefix unary operators is important when another operator follows the operand. For example, $-x^{\wedge} 2$ means $-\left(x^{\wedge} 2\right)$, because ' - ' has lower precedence than ' ${ }^{\prime \prime}$. .

Here is a table of the operators in Octave, in order of decreasing precedence. Unless noted, all operators group left to right.

```
function call and array indexing, cell array indexing, and structure element
indexing
    (()' '{}' '.'
postfix increment, and postfix decrement
    '++' '_-'
    These operators group right to left.
transpose and exponentiation
    `''.', (~)'.^'
unary plus, unary minus, prefix increment, prefix decrement, and logical "not"
    '+' '_' '++' '_-, (~) '!'
multiply and divide
    '*'/'`\'`.\'`.*'`./'
add, subtract
    '+'`_'
colon ':'
relational
    '<' '<=' '==' '>=' '>' '!=' '~='
element-wise "and"
    '&'
element-wise "or"
    'l'
```

```
logical "and"
    '&&'
logical "or"
            `|'
assignment
    '=' '+=' '-=' '*=' '/=' '\=' 'r=' '.*=' './=' '.\=' '. `=' '|=' '&='
```

    These operators group right to left.
    
## 9 Evaluation

Normally, you evaluate expressions simply by typing them at the Octave prompt, or by asking Octave to interpret commands that you have saved in a file.

Sometimes, you may find it necessary to evaluate an expression that has been computed and stored in a string, which is exactly what the eval function lets you do.
eval (try)
eval (try, catch)
Parse the string try and evaluate it as if it were an Octave program.
If execution fails, evaluate the optional string catch.
The string try is evaluated in the current context, so any results remain available after eval returns.
The following example creates the variable $A$ with the approximate value of 3.1416 in the current workspace.

```
eval ("A = acos(-1);");
```

If an error occurs during the evaluation of try then the catch string is evaluated, as the following example shows:

```
eval ('error ("This is a bad example");',
    'printf ("This error occurred:\n%s\n", lasterr ());');
    \dashvThis error occurred:
        This is a bad example
```

Programming Note: if you are only using eval as an error-capturing mechanism, rather than for the execution of arbitrary code strings, Consider using try/catch blocks or unwind_protect/unwind_protect_cleanup blocks instead. These techniques have higher performance and don't introduce the security considerations that the evaluation of arbitrary code does.
See also: [evalin], page 178, [evalc], page 175, [assignin], page 178, [feval], page 176.
The evalc function additionally captures any console output produced by the evaluated expression.

```
s = evalc (try)
s = evalc (try, catch)
```

Parse and evaluate the string try as if it were an Octave program, while capturing the output into the return variable $s$.
If execution fails, evaluate the optional string catch.
This function behaves like eval, but any output or warning messages which would normally be written to the console are captured and returned in the string $s$.
The diary is disabled during the execution of this function. When system is used, any output produced by external programs is not captured, unless their output is captured by the system function itself.

```
s = evalc ("t = 42"), t
    # s = t = 42
    # t = 42
```

See also: [eval], page 175, [diary], page 34.

### 9.1 Calling a Function by its Name

The feval function allows you to call a function from a string containing its name. This is useful when writing a function that needs to call user-supplied functions. The feval function takes the name of the function to call as its first argument, and the remaining arguments are given to the function.

The following example is a simple-minded function using feval that finds the root of a user-supplied function of one variable using Newton's method.

```
function result = newtroot (fname, x)
# usage: newtroot (fname, x)
#
# fname : a string naming a function f(x).
# x : initial guess
    delta = tol = sqrt (eps);
    maxit = 200;
    fx = feval (fname, x);
    for i = 1:maxit
        if (abs (fx) < tol)
            result = x;
            return;
        else
            fx_new = feval (fname, x + delta);
            deriv = (fx_new - fx) / delta;
            x = x - fx / deriv;
            fx = fx_new;
        endif
endfor
result = x;
endfunction
```

Note that this is only meant to be an example of calling user-supplied functions and should not be taken too seriously. In addition to using a more robust algorithm, any serious code would check the number and type of all the arguments, ensure that the supplied function really was a function, etc. See Section 4.8 [Predicates for Numeric Objects], page 67, for a list of predicates for numeric objects, and see Section 7.3 [Status of Variables], page 142, for a description of the exist function.
feval (name, ...)
Evaluate the function named name.
Any arguments after the first are passed as inputs to the named function. For example,

$$
\begin{gathered}
\text { feval ("acos", -1) } \\
\Rightarrow 3.1416
\end{gathered}
$$

calls the function acos with the argument ' -1 '.
The function feval can also be used with function handles of any sort (see Section 11.12.1 [Function Handles], page 233). Historically, feval was the only way to call user-supplied functions in strings, but function handles are now preferred due to the cleaner syntax they offer. For example,

```
f = @exp;
feval (f, 1)
    =>2.7183
f(1)
    =>2.7183
```

are equivalent ways to call the function referred to by $f$. If it cannot be predicted beforehand whether $f$ is a function handle, function name in a string, or inline function then feval can be used instead.

A similar function run exists for calling user script files, that are not necessarily on the user path
run script
run ("script")
Run script in the current workspace.
Scripts which reside in directories specified in Octave's load path, and which end with the extension .m, can be run simply by typing their name. For scripts not located on the load path, use run.

The filename script can be a bare, fully qualified, or relative filename and with or without a file extension. If no extension is specified, Octave will first search for a script with the .m extension before falling back to the script name without an extension.

Implementation Note: If script includes a path component, then run first changes the working directory to the directory where script is found. Next, the script is executed. Finally, run returns to the original working directory unless script has specifically changed directories.

See also: [path], page 214, [addpath], page 213, [source], page 224.

### 9.2 Evaluation in a Different Context

Before you evaluate an expression you need to substitute the values of the variables used in the expression. These are stored in the symbol table. Whenever the interpreter starts a new function it saves the current symbol table and creates a new one, initializing it with the list of function parameters and a couple of predefined variables such as nargin. Expressions inside the function use the new symbol table.

Sometimes you want to write a function so that when you call it, it modifies variables in your own context. This allows you to use a pass-by-name style of function, which is similar to using a pointer in programming languages such as C.

Consider how you might write save and load as m-files. For example:

```
function create_data
    x = linspace (0, 10, 10);
    y = sin (x);
    save mydata x y
endfunction
```

With evalin, you could write save as follows:

```
function save (file, name1, name2)
    f = open_save_file (file);
    save_var (f, name1, evalin ("caller", name1));
    save_var (f, name2, evalin ("caller", name2));
endfunction
```

Here, 'caller' is the create_data function and name1 is the string " x ", which evaluates simply as the value of $x$.

You later want to load the values back from mydata in a different context:

```
function process_data
    load mydata
    ... do work ...
endfunction
```

With assignin, you could write load as follows:

```
function load (file)
    f = open_load_file (file);
    [name, val] = load_var (f);
    assignin ("caller", name, val);
    [name, val] = load_var (f);
    assignin ("caller", name, val);
endfunction
```

Here, 'caller' is the process_data function.
You can set and use variables at the command prompt using the context 'base' rather than 'caller'.

These functions are rarely used in practice. One example is the fail ('code', 'pattern') function which evaluates 'code' in the caller's context and checks that the error message it produces matches the given pattern. Other examples such as save and load are written in C++ where all Octave variables are in the 'caller' context and evalin is not needed.

```
evalin (context, try)
evalin (context, try, catch)
```

Like eval, except that the expressions are evaluated in the context context, which may be either "caller" or "base".
See also: [eval], page 175, [assignin], page 178.
assignin (context, varname, value)
Assign value to varname in context context, which may be either "base" or "caller".
See also: [evalin], page 178 .

## 10 Statements

Statements may be a simple constant expression or a complicated list of nested loops and conditional statements.

Control statements such as if, while, and so on control the flow of execution in Octave programs. All the control statements start with special keywords such as if and while, to distinguish them from simple expressions. Many control statements contain other statements; for example, the if statement contains another statement which may or may not be executed.

Each control statement has a corresponding end statement that marks the end of the control statement. For example, the keyword endif marks the end of an if statement, and endwhile marks the end of a while statement. You can use the keyword end anywhere a more specific end keyword is expected, but using the more specific keywords is preferred because if you use them, Octave is able to provide better diagnostics for mismatched or missing end tokens.

The list of statements contained between keywords like if or while and the corresponding end statement is called the body of a control statement.

### 10.1 The if Statement

The if statement is Octave's decision-making statement. There are three basic forms of an if statement. In its simplest form, it looks like this:

```
if (condition)
    then-body
endif
```

condition is an expression that controls what the rest of the statement will do. The thenbody is executed only if condition is true.

The condition in an if statement is considered true if its value is nonzero, and false if its value is zero. If the value of the conditional expression in an if statement is a vector or a matrix, it is considered true only if it is non-empty and all of the elements are nonzero. The conceptually equivalent code when condition is a matrix is shown below.

```
if (matrix) \equiv if (all (matrix(:)))
```

The second form of an if statement looks like this:

```
if (condition)
    then-body
else
    else-body
endif
```

If condition is true, then-body is executed; otherwise, else-body is executed.
Here is an example:
if (rem (x, 2) $==0$ )
printf ("x is even\n");
else
printf ("x is odd\n");
endif

In this example, if the expression rem $(x, 2)=0$ is true (that is, the value of $x$ is divisible by 2 ), then the first printf statement is evaluated, otherwise the second printf statement is evaluated.

The third and most general form of the if statement allows multiple decisions to be combined in a single statement. It looks like this:

```
if (condition)
    then-body
elseif (condition)
    elseif-body
else
    else-body
endif
```

Any number of elseif clauses may appear. Each condition is tested in turn, and if one is found to be true, its corresponding body is executed. If none of the conditions are true and the else clause is present, its body is executed. Only one else clause may appear, and it must be the last part of the statement.

In the following example, if the first condition is true (that is, the value of x is divisible by 2 ), then the first printf statement is executed. If it is false, then the second condition is tested, and if it is true (that is, the value of x is divisible by 3 ), then the second printf statement is executed. Otherwise, the third printf statement is performed.

```
if (rem (x, 2) == 0)
    printf ("x is even\n");
elseif (rem (x, 3) == 0)
    printf ("x is odd and divisible by 3\n");
else
    printf ("x is odd\n");
endif
```

Note that the elseif keyword must not be spelled else if, as is allowed in Fortran. If it is, the space between the else and if will tell Octave to treat this as a new if statement within another if statement's else clause. For example, if you write

```
if (c1)
    body-1
else if (c2)
    body-2
endif
```

Octave will expect additional input to complete the first if statement. If you are using Octave interactively, it will continue to prompt you for additional input. If Octave is reading this input from a file, it may complain about missing or mismatched end statements, or, if you have not used the more specific end statements (endif, endfor, etc.), it may simply produce incorrect results, without producing any warning messages.

It is much easier to see the error if we rewrite the statements above like this,

```
if (c1)
    body-1
else
    if (c2)
            body-2
    endif
```

using the indentation to show how Octave groups the statements. See Chapter 11 [Functions and Scripts], page 191.

### 10.2 The switch Statement

It is very common to take different actions depending on the value of one variable. This is possible using the if statement in the following way

```
if (X == 1)
    do_something ();
elseif (X == 2)
    do_something_else ();
else
    do_something_completely_different ();
endif
```

This kind of code can however be very cumbersome to both write and maintain. To overcome this problem Octave supports the switch statement. Using this statement, the above example becomes

```
switch (X)
    case 1
        do_something ();
    case 2
        do_something_else ();
    otherwise
        do_something_completely_different ();
endswitch
```

This code makes the repetitive structure of the problem more explicit, making the code easier to read, and hence maintain. Also, if the variable $X$ should change its name, only one line would need changing compared to one line per case when if statements are used.

The general form of the switch statement is

```
switch (expression)
    case label
        command_list
    case label
        command_list
    ...
    otherwise
        command_list
endswitch
```

where label can be any expression. However, duplicate label values are not detected, and only the command_list corresponding to the first match will be executed. For the switch statement to be meaningful at least one case label command_list clause must be present, while the otherwise command_list clause is optional.

If label is a cell array the corresponding command_list is executed if any of the elements of the cell array match expression. As an example, the following program will print 'Variable is either 6 or 7 '.

```
A = 7;
switch (A)
    case { 6, 7 }
        printf ("variable is either 6 or 7\n");
        otherwise
        printf ("variable is neither 6 nor 7\n");
endswitch
```

As with all other specific end keywords, endswitch may be replaced by end, but you can get better diagnostics if you use the specific forms.

One advantage of using the switch statement compared to using if statements is that the labels can be strings. If an if statement is used it is not possible to write

```
if (X == "a string") # This is NOT valid
```

since a character-to-character comparison between X and the string will be made instead of evaluating if the strings are equal. This special-case is handled by the switch statement, and it is possible to write programs that look like this

```
switch (X)
    case "a string"
        do_something
endswitch
```


### 10.2.1 Notes for the C Programmer

The switch statement is also available in the widely used C programming language. There are, however, some differences between the statement in Octave and C

- Cases are exclusive, so they don't 'fall through' as do the cases in the switch statement of the C language.
- The command_list elements are not optional. Making the list optional would have meant requiring a separator between the label and the command list. Otherwise, things like

```
switch (foo)
    case (1) -2
```

would produce surprising results, as would

```
switch (foo)
    case (1)
    case (2)
        doit ();
```

particularly for C programmers. If doit() should be executed if foo is either 1 or 2, the above code should be written with a cell array like this

```
switch (foo)
    case { 1, 2 }
        doit ();
```


### 10.3 The while Statement

In programming, a loop means a part of a program that is (or at least can be) executed two or more times in succession.

The while statement is the simplest looping statement in Octave. It repeatedly executes a statement as long as a condition is true. As with the condition in an if statement, the condition in a while statement is considered true if its value is nonzero, and false if its value is zero. If the value of the conditional expression in a while statement is a vector or a matrix, it is considered true only if it is non-empty and all of the elements are nonzero.

Octave's while statement looks like this:

```
while (condition)
    body
endwhile
```

Here body is a statement or list of statements that we call the body of the loop, and condition is an expression that controls how long the loop keeps running.

The first thing the while statement does is test condition. If condition is true, it executes the statement body. After body has been executed, condition is tested again, and if it is still true, body is executed again. This process repeats until condition is no longer true. If condition is initially false, the body of the loop is never executed.

This example creates a variable fib that contains the first ten elements of the Fibonacci sequence.

```
fib = ones (1, 10);
i = 3;
while (i <= 10)
    fib (i) = fib (i-1) + fib (i-2);
    i++;
endwhile
```

Here the body of the loop contains two statements.
The loop works like this: first, the value of $i$ is set to 3 . Then, the while tests whether $i$ is less than or equal to 10 . This is the case when $i$ equals 3 , so the value of the $i$-th element of fib is set to the sum of the previous two values in the sequence. Then the i++ increments the value of $i$ and the loop repeats. The loop terminates when i reaches 11 .

A newline is not required between the condition and the body; but using one makes the program clearer unless the body is very simple.

### 10.4 The do-until Statement

The do-until statement is similar to the while statement, except that it repeatedly executes a statement until a condition becomes true, and the test of the condition is at the end of the loop, so the body of the loop is always executed at least once. As with the condition in an if statement, the condition in a do-until statement is considered true if its value is nonzero, and false if its value is zero. If the value of the conditional expression in a do-until statement is a vector or a matrix, it is considered true only if it is non-empty and all of the elements are nonzero.

Octave's do-until statement looks like this:

```
do
    body
until (condition)
```

Here body is a statement or list of statements that we call the body of the loop, and condition is an expression that controls how long the loop keeps running.

This example creates a variable fib that contains the first ten elements of the Fibonacci sequence.

```
fib = ones (1, 10);
i = 2;
do
    i++;
    fib (i) = fib (i-1) + fib (i-2);
until (i == 10)
```

A newline is not required between the do keyword and the body; but using one makes the program clearer unless the body is very simple.

### 10.5 The for Statement

The for statement makes it more convenient to count iterations of a loop. The general form of the for statement looks like this:

```
for var = expression
    body
endfor
```

where body stands for any statement or list of statements, expression is any valid expression, and var may take several forms. Usually it is a simple variable name or an indexed variable. If the value of expression is a structure, var may also be a vector with two elements. See Section 10.5.1 [Looping Over Structure Elements], page 185, below.

The assignment expression in the for statement works a bit differently than Octave's normal assignment statement. Instead of assigning the complete result of the expression, it assigns each column of the expression to var in turn. If expression is a range, a row vector, or a scalar, the value of var will be a scalar each time the loop body is executed. If var is a column vector or a matrix, var will be a column vector each time the loop body is executed.

The following example shows another way to create a vector containing the first ten elements of the Fibonacci sequence, this time using the for statement:

```
fib = ones (1, 10);
for i = 3:10
    fib(i) = fib(i-1) + fib(i-2);
endfor
```

This code works by first evaluating the expression $3: 10$, to produce a range of values from 3 to 10 inclusive. Then the variable $i$ is assigned the first element of the range and the body of the loop is executed once. When the end of the loop body is reached, the next value in the range is assigned to the variable $i$, and the loop body is executed again. This process continues until there are no more elements to assign.

Within Octave is it also possible to iterate over matrices or cell arrays using the for statement. For example consider

```
disp ("Loop over a matrix")
for i = [1,3;2,4]
    i
endfor
disp ("Loop over a cell array")
for i = {1,"two";"three",4}
    i
endfor
```

In this case the variable $i$ takes on the value of the columns of the matrix or cell matrix. So the first loop iterates twice, producing two column vectors [1;2], followed by [3;4], and likewise for the loop over the cell array. This can be extended to loops over multidimensional arrays. For example:

```
a = [1,3;2,4]; c = cat (3, a, 2*a);
for i = c
    i
endfor
```

In the above case, the multi-dimensional matrix $c$ is reshaped to a two-dimensional matrix as reshape (c, rows (c), prod (size (c)(2:end))) and then the same behavior as a loop over a two-dimensional matrix is produced.

Although it is possible to rewrite all for loops as while loops, the Octave language has both statements because often a for loop is both less work to type and more natural to think of. Counting the number of iterations is very common in loops and it can be easier to think of this counting as part of looping rather than as something to do inside the loop.

### 10.5.1 Looping Over Structure Elements

A special form of the for statement allows you to loop over all the elements of a structure:

```
for [ val, key ] = expression
    body
endfor
```

In this form of the for statement, the value of expression must be a structure. If it is, key and val are set to the name of the element and the corresponding value in turn, until there are no more elements. For example:

```
x.a = 1
x.b = [1, 2; 3, 4]
x.c = "string"
for [val, key] = x
    key
    val
endfor
\dashv key = a
\val = 1
\dashv key = b
\dagger val =
\dashv
\dashv1 2
\dashv 3 4
\dashv
\dashv key = c
\val = string
```

The elements are not accessed in any particular order. If you need to cycle through the list in a particular way, you will have to use the function fieldnames and sort the list yourself.

### 10.6 The break Statement

The break statement jumps out of the innermost while, do-until, or for loop that encloses it. The break statement may only be used within the body of a loop. The following example finds the smallest divisor of a given integer, and also identifies prime numbers:

```
num = 103;
div = 2;
while (div*div <= num)
    if (rem (num, div) == 0)
        break;
    endif
    div++;
endwhile
if (rem (num, div) == 0)
    printf ("Smallest divisor of %d is %d\n", num, div)
else
    printf ("%d is prime\n", num);
endif
```

When the remainder is zero in the first while statement, Octave immediately breaks out of the loop. This means that Octave proceeds immediately to the statement following the loop and continues processing. (This is very different from the exit statement which stops the entire Octave program.)

Here is another program equivalent to the previous one. It illustrates how the condition of a while statement could just as well be replaced with a break inside an if:

```
num = 103;
div = 2;
while (1)
    if (rem (num, div) == 0)
        printf ("Smallest divisor of %d is %d\n", num, div);
        break;
    endif
    div++;
    if (div*div > num)
        printf ("%d is prime\n", num);
        break;
    endif
endwhile
```


### 10.7 The continue Statement

The continue statement, like break, is used only inside while, do-until, or for loops. It skips over the rest of the loop body, causing the next cycle around the loop to begin immediately. Contrast this with break, which jumps out of the loop altogether. Here is an example:

```
# print elements of a vector of random
# integers that are even.
# first, create a row vector of 10 random
# integers with values between 0 and 100:
vec = round (rand (1, 10) * 100);
# print what we're interested in:
for x = vec
    if (rem (x, 2) != 0)
        continue;
    endif
    printf ("%d\n", x);
endfor
```

If one of the elements of vec is an odd number, this example skips the print statement for that element, and continues back to the first statement in the loop.

This is not a practical example of the continue statement, but it should give you a clear understanding of how it works. Normally, one would probably write the loop like this:

```
for x = vec
    if (rem (x, 2) == 0)
        printf ("%d\n", x);
    endif
endfor
```


### 10.8 The unwind_protect Statement

Octave supports a limited form of exception handling modeled after the unwind-protect form of Lisp.

The general form of an unwind_protect block looks like this:

```
unwind_protect
    body
unwind_protect_cleanup
    cleanup
end_unwind_protect
```

where body and cleanup are both optional and may contain any Octave expressions or commands. The statements in cleanup are guaranteed to be executed regardless of how control exits body.

This is useful to protect temporary changes to global variables from possible errors. For example, the following code will always restore the original value of the global variable frobnosticate even if an error occurs in the first part of the unwind_protect block.

```
save_frobnosticate = frobnosticate;
unwind_protect
    frobnosticate = true;
unwind_protect_cleanup
    frobnosticate = save_frobnosticate;
end_unwind_protect
```

Without unwind_protect, the value of frobnosticate would not be restored if an error occurs while evaluating the first part of the unwind_protect block because evaluation would stop at the point of the error and the statement to restore the value would not be executed.

In addition to unwind_protect, Octave supports another form of exception handling, the try block.

### 10.9 The try Statement

The original form of a try block looks like this:

```
try
    body
catch
    cleanup
end_try_catch
```

where body and cleanup are both optional and may contain any Octave expressions or commands. The statements in cleanup are only executed if an error occurs in body.

No warnings or error messages are printed while body is executing. If an error does occur during the execution of body, cleanup can use the functions lasterr or lasterror to access the text of the message that would have been printed, as well as its identifier. The alternative form,

```
try
    body
catch err
    cleanup
end_try_catch
```

will automatically store the output of lasterror in the structure err. See Chapter 12 [Errors and Warnings], page 239, for more information about the lasterr and lasterror functions.

### 10.10 Continuation Lines

In the Octave language, most statements end with a newline character and you must tell Octave to ignore the newline character in order to continue a statement from one line to the next. Lines that end with the characters ... are joined with the following line before they are divided into tokens by Octave's parser. For example, the lines

```
x = long_variable_name ...
    + longer_variable_name ...
    -42
```

form a single statement.
Any text between the continuation marker and the newline character is ignored. For example, the statement

```
x = long_variable_name ... # comment one
    + longer_variable_name ...comment two
    - 42 # last comment
```

is equivalent to the one shown above.
Inside double-quoted string constants, the character \has to be used as continuation marker. The $\backslash$ must appear at the end of the line just before the newline character:

```
s = "This text starts in the first line \
and is continued in the second line."
```

Input that occurs inside parentheses can be continued to the next line without having to use a continuation marker. For example, it is possible to write statements like

```
if (fine_dining_destination == on_a_boat
    || fine_dining_destination == on_a_train)
    seuss (i, will, not, eat, them, sam, i, am, i,
        will, not, eat, green, eggs, and, ham);
endif
```

without having to add to the clutter with continuation markers.

## 11 Functions and Scripts

Complicated Octave programs can often be simplified by defining functions. Functions can be defined directly on the command line during interactive Octave sessions, or in external files, and can be called just like built-in functions.

### 11.1 Introduction to Function and Script Files

There are seven different things covered in this section.

1. Typing in a function at the command prompt.
2. Storing a group of commands in a file - called a script file.
3. Storing a function in a file - called a function file.
4. Subfunctions in function files.
5. Multiple functions in one script file.
6. Private functions.
7. Nested functions.

Both function files and script files end with an extension of .m, for mATLAB compatibility. If you want more than one independent functions in a file, it must be a script file (see Section 11.11 [Script Files], page 223), and to use these functions you must execute the script file before you can use the functions that are in the script file.

### 11.2 Defining Functions

In its simplest form, the definition of a function named name looks like this:

```
function name
    body
endfunction
```

A valid function name is like a valid variable name: a sequence of letters, digits and underscores, not starting with a digit. Functions share the same pool of names as variables.

The function body consists of Octave statements. It is the most important part of the definition, because it says what the function should actually $d o$.

For example, here is a function that, when executed, will ring the bell on your terminal (assuming that it is possible to do so):

```
function wakeup
    printf ("\a");
endfunction
```

The printf statement (see Chapter 14 [Input and Output], page 271) simply tells Octave to print the string " $\backslash \mathrm{a}$ ". The special character ' a ' stands for the alert character (ASCII 7). See Chapter 5 [Strings], page 71.

Once this function is defined, you can ask Octave to evaluate it by typing the name of the function.

Normally, you will want to pass some information to the functions you define. The syntax for passing parameters to a function in Octave is

```
function name (arg-list)
    body
endfunction
```

where arg-list is a comma-separated list of the function's arguments. When the function is called, the argument names are used to hold the argument values given in the call. The list of arguments may be empty, in which case this form is equivalent to the one shown above.

To print a message along with ringing the bell, you might modify the wakeup to look like this:

```
function wakeup (message)
    printf ("\a%s\n", message);
endfunction
```

Calling this function using a statement like this

```
wakeup ("Rise and shine!");
```

will cause Octave to ring your terminal's bell and print the message 'Rise and shine!', followed by a newline character (the ' $\backslash n$ ' in the first argument to the printf statement).

In most cases, you will also want to get some information back from the functions you define. Here is the syntax for writing a function that returns a single value:

```
function ret-var = name (arg-list)
    body
endfunction
```

The symbol ret-var is the name of the variable that will hold the value to be returned by the function. This variable must be defined before the end of the function body in order for the function to return a value.

Variables used in the body of a function are local to the function. Variables named in arg-list and ret-var are also local to the function. See Section 7.1 [Global Variables], page 139, for information about how to access global variables inside a function.

For example, here is a function that computes the average of the elements of a vector:

```
function retval = avg (v)
    retval = sum (v) / length (v);
endfunction
```

If we had written avg like this instead,

```
function retval = avg (v)
    if (isvector (v))
        retval = sum (v) / length (v);
    endif
endfunction
```

and then called the function with a matrix instead of a vector as the argument, Octave would have printed an error message like this:
error: value on right hand side of assignment is undefined
because the body of the if statement was never executed, and retval was never defined. To prevent obscure errors like this, it is a good idea to always make sure that the return
variables will always have values, and to produce meaningful error messages when problems are encountered. For example, avg could have been written like this:

```
function retval \(=\) avg (v)
    retval = 0;
    if (isvector (v))
        retval \(=\) sum (v) / length (v);
    else
        error ("avg: expecting vector argument");
    endif
endfunction
```

There is still one remaining problem with this function. What if it is called without an argument? Without additional error checking, Octave will probably print an error message that won't really help you track down the source of the error. To allow you to catch errors like this, Octave provides each function with an automatic variable called nargin. Each time a function is called, nargin is automatically initialized to the number of arguments that have actually been passed to the function. For example, we might rewrite the avg function like this:

```
function retval = avg (v)
    retval = 0;
    if (nargin != 1)
        usage ("avg (vector)");
    endif
    if (isvector (v))
        retval = sum (v) / length (v);
    else
        error ("avg: expecting vector argument");
    endif
endfunction
```

Octave automatically reports an error for functions written in .m file code if they are called with more arguments than expected. Octave does not automatically report an error if a function is called with too few arguments, since functions in general may have default arguments, but any attempt to use a variable that has not been given a value will result in an error. Functions can check the arguments they are called with to avoid such problems and to provide more context-specific error messages.

```
nargin()
nargin (fcn)
```

Report the number of input arguments to a function.
Called from within a function, return the number of arguments passed to the function. At the top level, return the number of command line arguments passed to Octave.
If called with the optional argument fcn-a function name or handle - return the declared number of arguments that the function can accept.
If the last argument to $f \subset n$ is varargin the returned value is negative. For example, the function union for sets is declared as

```
function [y, ia, ib] = union (a, b, varargin)
and
nargin ("union")
# -3
```

Programming Note: nargin does not work on compiled functions (.oct files) such as built-in or dynamically loaded functions.
See also: [nargout], page 197, [narginchk], page 203, [varargin], page 199, [inputname], page 194.
inputname ( $n$ )
inputname ( $n$, ids_only)
Return the name of the $n$-th argument to the calling function.
If the argument is not a simple variable name, return an empty string. Examples which will return "" are numbers (5.1), expressions (y/2), and cell or structure indexing (c\{1\} or s.field).
inputname is only useful within a function. When used at the command line or within a script it always returns an empty string.
By default, return an empty string if the $n$-th argument is not a valid variable name. If the optional argument ids_only is false, return the text of the argument even if it is not a valid variable name. This is an Octave extension that allows the programmer to view exactly how the function was invoked even when the inputs are complex expressions.
See also: [nargin], page 193, [narginchk], page 203.
val = silent_functions ()
old_val = silent_functions (new_val)
silent_functions (new_val, "local")
Query or set the internal variable that controls whether internal output from a function is suppressed.
If this option is disabled, Octave will display the results produced by evaluating expressions within a function body that are not terminated with a semicolon.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

### 11.3 Returning from a Function

The body of a user-defined function can contain a return statement. This statement returns control to the rest of the Octave program. It looks like this:

## return

Unlike the return statement in C, Octave's return statement cannot be used to return a value from a function. Instead, you must assign values to the list of return variables that are part of the function statement. The return statement simply makes it easier to exit a function from a deeply nested loop or conditional statement.

Here is an example of a function that checks to see if any elements of a vector are nonzero.

```
function retval = any_nonzero (v)
    retval = 0;
    for i = 1:length (v)
            if (v (i) != 0)
            retval = 1;
            return;
        endif
    endfor
    printf ("no nonzero elements found\n");
endfunction
```

Note that this function could not have been written using the break statement to exit the loop once a nonzero value is found without adding extra logic to avoid printing the message if the vector does contain a nonzero element.
return
When Octave encounters the keyword return inside a function or script, it returns control to the caller immediately. At the top level, the return statement is ignored. A return statement is assumed at the end of every function definition.

### 11.4 Multiple Return Values

Unlike many other computer languages, Octave allows you to define functions that return more than one value. The syntax for defining functions that return multiple values is

```
function [ret-list] = name (arg-list)
    body
endfunction
```

where name, arg-list, and body have the same meaning as before, and ret-list is a commaseparated list of variable names that will hold the values returned from the function. The list of return values must have at least one element. If ret-list has only one element, this form of the function statement is equivalent to the form described in the previous section.

Here is an example of a function that returns two values, the maximum element of a vector and the index of its first occurrence in the vector.

```
function [max, idx] = vmax (v)
    idx = 1;
    max = v (idx);
    for i = 2:length (v)
        if (v (i) > max)
            max = v (i);
            idx = i;
        endif
    endfor
endfunction
```

In this particular case, the two values could have been returned as elements of a single array, but that is not always possible or convenient. The values to be returned may not
have compatible dimensions, and it is often desirable to give the individual return values distinct names.

It is possible to use the nthargout function to obtain only some of the return values or several at once in a cell array. See Section 3.1.5 [Cell Array Objects], page 46.
nthargout (n, func, ...)
nthargout ( $n$, ntot, func, ...)
Return the nth output argument of the function specified by the function handle or string func.
Any additional arguments are passed directly to func. The total number of arguments to call func with can be passed in ntot; by default ntot is $n$. The input $n$ can also be a vector of indices of the output, in which case the output will be a cell array of the requested output arguments.
The intended use nthargout is to avoid intermediate variables. For example, when finding the indices of the maximum entry of a matrix, the following two compositions of nthargout

```
m = magic (5);
cell2mat (nthargout ([1, 2], @ind2sub, size (m),
    nthargout (2, @max, m(:))))
# 5 3
```

are completely equivalent to the following lines:

```
m = magic (5);
[~}, idx] = max (M(:))
[i, j] = ind2sub (size (m), idx);
[i, j]
# 5
```

It can also be helpful to have all output arguments in a single cell in the following manner:

```
USV = nthargout ([1:3], @svd, hilb (5));
```

See also: [nargin], page 193, [nargout], page 197, [varargin], page 199, [varargout], page 198, [isargout], page 201.

In addition to setting nargin each time a function is called, Octave also automatically initializes nargout to the number of values that are expected to be returned. This allows you to write functions that behave differently depending on the number of values that the user of the function has requested. The implicit assignment to the built-in variable ans does not figure in the count of output arguments, so the value of nargout may be zero.

The svd and hist functions are examples of built-in functions that behave differently depending on the value of nargout. For example, hist will draw a histogram when called with no output variables, but if called with outputs it will return the frequency counts and/or bin centers without creating a plot.

It is possible to write functions that only set some return values. For example, calling the function

```
function \([\mathrm{x}, \mathrm{y}, \mathrm{z}]=\mathrm{f}\) ()
    \(\mathrm{x}=1\);
    z = 2;
endfunction
```

as
$[\mathrm{a}, \mathrm{b}, \mathrm{c}]=\mathrm{f}()$
produces:
$a=1$
b $=[](0 x 0)$
$c=2$
along with a warning.
nargout ()
nargout ( $f$ cn)
Report the number of output arguments from a function.
Called from within a function, return the number of values the caller expects to receive. At the top level, nargout with no argument is undefined and will produce an error.
If called with the optional argument fcn-a function name or handle - return the number of declared output values that the function can produce.

If the final output argument is varargout the returned value is negative.
For example,
f ()
will cause nargout to return 0 inside the function $f$ and

$$
[\mathrm{s}, \mathrm{t}]=\mathrm{f}()
$$

will cause nargout to return 2 inside the function $f$.
In the second usage,
nargout (@histc) \# or nargout ("histc") using a string input
will return 2, because histc has two outputs, whereas
nargout (@imread)
will return -2 , because imread has two outputs and the second is varargout.
Programming Note. nargout does not work for built-in functions and returns -1 for all anonymous functions.

See also: [nargin], page 193, [varargout], page 198, [isargout], page 201, [nthargout], page 196.

### 11.5 Variable-length Return Lists

It is possible to return a variable number of output arguments from a function using a syntax that's similar to the one used with the special varargin parameter name. To let a function return a variable number of output arguments the special output parameter name varargout is used. As with varargin, varargout is a cell array that will contain the requested output arguments.

As an example the following function sets the first output argument to 1 , the second to 2 , and so on.

```
function varargout = one_to_n ()
    for i = 1:nargout
        varargout{i} = i;
    endfor
endfunction
```

When called this function returns values like this

$$
\begin{aligned}
{[\mathrm{a}, \mathrm{~b}, \mathrm{c}] } & =\text { one_to_n () } \\
\Rightarrow \mathrm{a} & =1 \\
\Rightarrow \mathrm{~b} & =2 \\
\Rightarrow \mathrm{c} & =3
\end{aligned}
$$

If varargin (varargout) does not appear as the last element of the input (output) parameter list, then it is not special, and is handled the same as any other parameter name.
$[r 1, r 2, \ldots, r n]=\operatorname{deal}(a)$
$[r 1, r 2, \ldots, r n]=\operatorname{deal}(a 1, a 2, \ldots, a n)$
Copy the input parameters into the corresponding output parameters.
If only a single input parameter is supplied, its value is copied to each of the outputs.
For example,

$$
[\mathrm{a}, \mathrm{~b}, \mathrm{c}]=\operatorname{deal}(\mathrm{x}, \mathrm{y}, \mathrm{z}) ;
$$

is equivalent to

$$
\begin{aligned}
& \mathrm{a}=\mathrm{x} \\
& \mathrm{~b}=\mathrm{y} \\
& \mathrm{c}=\mathrm{z}
\end{aligned}
$$

and

$$
[\mathrm{a}, \mathrm{~b}, \mathrm{c}]=\operatorname{deal}(\mathrm{x}) ;
$$

is equivalent to

$$
\mathrm{a}=\mathrm{b}=\mathrm{c}=\mathrm{x} ;
$$

Programming Note: deal is often used with comma separated lists derived from cell arrays or structures. This is unnecessary as the interpreter can perform the same action without the overhead of a function call. For example:

```
c = {[11 2], "Three", 4};
[x, y, z] = c{:}
=>
    x =
            1 2
        y = Three
    z = 4
```

See also: [cell2struct], page 134, [struct2cell], page 123, [repmat], page 527.

### 11.6 Variable-length Argument Lists

Sometimes the number of input arguments is not known when the function is defined. As an example think of a function that returns the smallest of all its input arguments. For example:

```
a = smallest (1, 2, 3);
b = smallest (1, 2, 3, 4);
```

In this example both a and b would be 1 . One way to write the smallest function is

```
function val = smallest (arg1, arg2, arg3, arg4, arg5)
    body
endfunction
```

and then use the value of nargin to determine which of the input arguments should be considered. The problem with this approach is that it can only handle a limited number of input arguments.

If the special parameter name varargin appears at the end of a function parameter list it indicates that the function takes a variable number of input arguments. Using varargin the function looks like this

```
function val = smallest (varargin)
    body
endfunction
```

In the function body the input arguments can be accessed through the variable varargin. This variable is a cell array containing all the input arguments. See Section 6.3 [Cell Arrays], page 124, for details on working with cell arrays. The smallest function can now be defined like this

```
function val = smallest (varargin)
    val = min ([varargin{:}]);
endfunction
```

This implementation handles any number of input arguments, but it's also a very simple solution to the problem.

A slightly more complex example of varargin is a function print_arguments that prints all input arguments. Such a function can be defined like this

```
function print_arguments (varargin)
    for i = 1:length (varargin)
        printf ("Input argument %d: ", i);
        disp (varargin{i});
    endfor
endfunction
```

This function produces output like this

```
print_arguments (1, "two", 3);
    \dashv Input argument 1: 1
    - Input argument 2: two
    \dashvInput argument 3: 3
```

[reg, prop] = parseparams (params)
[reg, var1, ...] = parseparams (params, name1, default1, ...)

Return in reg the cell elements of param up to the first string element and in prop all remaining elements beginning with the first string element.
For example:

```
[reg, prop] = parseparams ({1, 2, "linewidth", 10})
reg =
{
    [1,1] = 1
    [1,2] = 2
}
prop =
{
    [1,1] = linewidth
    [1,2] = 10
}
```

The parseparams function may be used to separate regular numeric arguments from additional arguments given as property/value pairs of the varargin cell array.
In the second form of the call, available options are specified directly with their default values given as name-value pairs. If params do not form name-value pairs, or if an option occurs that does not match any of the available options, an error occurs.
When called from an m-file function, the error is prefixed with the name of the caller function.
The matching of options is case-insensitive.
See also: [varargin], page 199, [inputParser], page 208.

### 11.7 Ignoring Arguments

In the formal argument list, it is possible to use the dummy placeholder ~ instead of a name. This indicates that the corresponding argument value should be ignored and not stored to any variable.

```
function val = pick2nd (~, arg2)
    val = arg2;
endfunction
```

The value of nargin is not affected by using this declaration.
Return arguments can also be ignored using the same syntax. For example, the sort function returns both the sorted values, and an index vector for the original input which will result in a sorted output. Ignoring the second output is simple - don't request more than one output. But ignoring the first, and calculating just the second output, requires the use of the ~ placeholder.

```
x = [2, 3, 1];
[s, i] = sort (x)
=>
s =
    1 2 3
i =
    3 1 2
[~
#
i =
312
```

When using the ~ placeholder, commas-not whitespace - must be used to separate output arguments. Otherwise, the interpreter will view ~ as the logical not operator.
[~ i] = sort (x)
parse error:

```
invalid left hand side of assignment
```

Functions may take advantage of ignored outputs to reduce the number of calculations performed. To do so, use the isargout function to query whether the output argument is wanted. For example:

```
function [out1, out2] = long_function (x, y, z)
    if (isargout (1))
        ## Long calculation
        out1 = result;
    endif
endfunction
```

isargout (k)

Within a function, return a logical value indicating whether the argument $k$ will be assigned to a variable on output.
If the result is false, the argument has been ignored during the function call through the use of the tilde ( ${ }^{\sim}$ ) special output argument. Functions can use isargout to avoid performing unnecessary calculations for outputs which are unwanted.

If $k$ is outside the range $1: \max$ (nargout), the function returns false. $k$ can also be an array, in which case the function works element-by-element and a logical array is returned. At the top level, isargout returns an error.

See also: [nargout], page 197, [varargout], page 198, [nthargout], page 196.

### 11.8 Default Arguments

Since Octave supports variable number of input arguments, it is very useful to assign default values to some input arguments. When an input argument is declared in the argument list it is possible to assign a default value to the argument like this

```
function name (arg1 = val1, ...)
    body
endfunction
```

If no value is assigned to arg1 by the user, it will have the value val1.
As an example, the following function implements a variant of the classic "Hello, World" program.

```
function hello (who = "World")
    printf ("Hello, %s!\n", who);
endfunction
```

When called without an input argument the function prints the following

```
hello ();
    Hello, World!
```

and when it's called with an input argument it prints the following

```
hello ("Beautiful World of Free Software");
    Hello, Beautiful World of Free Software!
```

Sometimes it is useful to explicitly tell Octave to use the default value of an input argument. This can be done writing a ':' as the value of the input argument when calling the function.

```
hello (:);
    Hello, World!
```


### 11.9 Validating Arguments

Octave is a weakly typed programming language. Thus it is possible to call a function with arguments, that probably cause errors or might have undesirable side effects. For example calling a string processing function with a huge sparse matrix.

It is good practice at the head of a function to verify that it has been called correctly. Octave offers several functions for this purpose.

### 11.9.1 Validating the number of Arguments

In Octave the following idiom is seen frequently at the beginning of a function definition:

```
if (nargin < min_#_inputs || nargin > max_#_inputs)
    print_usage ();
endif
```

which stops the function execution and prints a message about the correct way to call the function whenever the number of inputs is wrong.

Similar error checking is provided by narginchk and nargoutchk.
narginchk (minargs, maxargs)
Check for correct number of input arguments.
Generate an error message if the number of arguments in the calling function is outside the range minargs and maxargs. Otherwise, do nothing.
Both minargs and maxargs must be scalar numeric values. Zero, Inf, and negative values are all allowed, and minargs and maxargs may be equal.
Note that this function evaluates nargin on the caller.
See also: [nargoutchk], page 203, [error], page 239, [nargout], page 197, [nargin], page 193.
nargoutchk (minargs, maxargs)
msgstr = nargoutchk (minargs, maxargs, nargs)
msgstr = nargoutchk (minargs, maxargs, nargs, "string")
msgstruct = nargoutchk (minargs, maxargs, nargs, "struct")
Check for correct number of output arguments.
In the first form, return an error if the number of arguments is not between minargs and maxargs. Otherwise, do nothing. Note that this function evaluates the value of nargout on the caller so its value must have not been tampered with.
Both minargs and maxargs must be numeric scalars. Zero, Inf, and negative are all valid, and they can have the same value.
For backwards compatibility, the other forms return an appropriate error message string (or structure) if the number of outputs requested is invalid.
This is useful for checking to that the number of output arguments supplied to a function is within an acceptable range.
See also: [narginchk], page 203, [error], page 239, [nargout], page 197, [nargin], page 193.

### 11.9.2 Validating the type of Arguments

Besides the number of arguments, inputs can be checked for various properties. validatestring is used for string arguments and validateattributes for numeric arguments.

```
validstr = validatestring (str, strarray)
validstr = validatestring (str, strarray, funcname)
validstr = validatestring (str, strarray, funcname, varname)
validstr = validatestring (..., position)
Verify that str is an element, or substring of an element, in strarray.
```

When str is a character string to be tested, and strarray is a cell array of strings of valid values, then validstr will be the validated form of $s t r$ where validation is defined as str being a member or substring of validstr. This is useful for both verifying and expanding short options, such as "r", to their longer forms, such as "red". If str is
a substring of validstr, and there are multiple matches, the shortest match will be returned if all matches are substrings of each other. Otherwise, an error will be raised because the expansion of str is ambiguous. All comparisons are case insensitive.
The additional inputs funcname, varname, and position are optional and will make any generated validation error message more specific.
Examples:

```
validatestring ("r", {"red", "green", "blue"})
=> "red"
validatestring ("b", {"red", "green", "blue", "black"})
=> error: validatestring: multiple unique matches were found for 'b':
    blue, black
```

See also: [strcmp], page 84, [strcmpi], page 85, [validateattributes], page 204, [inputParser], page 208.

```
validateattributes (A, classes, attributes)
validateattributes (A, classes, attributes, arg_idx)
validateattributes (A, classes, attributes, func_name)
validateattributes (A, classes, attributes, func_name, arg_name)
validateattributes (A, classes, attributes, func_name, arg_name,
arg_idx)
```

Check validity of input argument.
Confirms that the argument $A$ is valid by belonging to one of classes, and holding all of the attributes. If it does not, an error is thrown, with a message formatted accordingly. The error message can be made further complete by the function name fun_name, the argument name arg_name, and its position in the input arg_idx.
classes must be a cell array of strings (an empty cell array is allowed) with the name of classes (remember that a class name is case sensitive). In addition to the class name, the following categories names are also valid:
"float" Floating point value comprising classes "double" and "single".
"integer"
Integer value comprising classes (u)int8, (u)int16, (u)int32, (u)int64.
"numeric"
Numeric value comprising either a floating point or integer value.
attributes must be a cell array with names of checks for $A$. Some of them require an additional value to be supplied right after the name (see details for each below).
"<=" All values are less than or equal to the following value in attributes.
"く" All values are less than the following value in attributes.
" $>=$ =" All values are greater than or equal to the following value in attributes.
">" All values are greater than the following value in attributes.
"2d" A 2-dimensional matrix. Note that vectors and empty matrices have 2 dimensions, one of them being of length 1 , or both length 0 .
"3d" Has no more than 3 dimensions. A 2-dimensional matrix is a 3-D matrix whose 3 rd dimension is of length 1 .
"binary" All values are either 1 or 0 .
"column" Values are arranged in a single column.
"decreasing"
No value is NaN , and each is less than the preceding one.
"diag" Value is a diagonal matrix.
"even" All values are even numbers.
"finite" All values are finite.
"increasing"
No value is $N a N$, and each is greater than the preceding one.
"integer"
All values are integer. This is different than using isinteger which only checks its an integer type. This checks that each value in $A$ is an integer value, i.e., it has no decimal part.
"ncols" Has exactly as many columns as the next value in attributes.
"ndims" Has exactly as many dimensions as the next value in attributes.
"nondecreasing"
No value is $N a N$, and each is greater than or equal to the preceding one.
"nonempty"
It is not empty.
"nonincreasing"
No value is NaN , and each is less than or equal to the preceding one.
"nonnan" No value is a NaN.
"nonnegative"
All values are non negative.
"nonsparse"
It is not a sparse matrix.
"nonzero"
No value is zero.
"nrows" Has exactly as many rows as the next value in attributes.
"numel" Has exactly as many elements as the next value in attributes.
"odd" All values are odd numbers.
"positive"
All values are positive.
"real" It is a non-complex matrix.
"row" Values are arranged in a single row.
"scalar" It is a scalar.
"size" Its size has length equal to the values of the next in attributes. The next value must is an array with the length for each dimension. To ignore the check for a certain dimension, the value of NaN can be used.
"square" Is a square matrix.
"vector" Values are arranged in a single vector (column or vector).
See also: [isa], page 41, [validatestring], page 203, [inputParser], page 208.
As alternatives to validateattributes there are several shorter convenience functions to check for individual properties.
mustBeFinite ( $x$ )
Require that input $x$ is finite.
Raise an error if any element of the input $x$ is not finite, as determined by isfinite (x).

See also: [mustBeNonNan], page 207, [isfinite], page 513.
mustBeGreaterThan (x, c)
Require that input $x$ is greater than $c$.
Raise an error if any element of the input $x$ is not greater than $c$, as determined by $x>c$.
See also: [mustBeGreaterThanOrEqual], page 206, [mustBeLessThan], page 206, [gt], page 165.
mustBeGreaterThanOrEqual ( $x, c$ )
Require that input $x$ is greater than or equal to $c$.
Raise an error if any element of the input x is not greater than or equal to $c$, as determined by $\mathrm{x}>=c$.
See also: [mustBeGreaterThan], page 206, [mustBeLessThanOrEqual], page 207, [ge], page 165.
mustBeInteger ( $x$ )
Require that input $x$ is integer-valued (but not necessarily integer-typed).
Raise an error if any element of the input $x$ is not a finite, real, integer-valued numeric value, as determined by various checks.
See also: [mustBeNumeric], page 208.
mustBeLessThan ( $x, c$ )
Require that input $x$ is less than $c$.
Raise an error if any element of the input x is not less than $c$, as determined by $\mathrm{x}<$ c.

See also: [mustBeLessThanOrEqual], page 207, [mustBeGreaterThan], page 206, [lt], page 165.
mustBeLessThanOrEqual ( $x, c$ )
Require that input is less than or equal to a given value.
Raise an error if any element of the input $x$ is not less than or equal to $c$, as determined by $x<=c$.
See also: [mustBeLessThan], page 206, [mustBeGreaterThanOrEqual], page 206, [le], page 165.
mustBeMember (x, valid)
Require that input $x$ is a member of a set of given valid values.
Raise an error if any element of the input $x$ is not a member of the set valid, as determined by ismember ( x ).
Programming Note: char inputs may behave strangely because of the interaction between chars and cellstrings when calling ismember on them. But it will probably "do what you mean" if you just use it naturally. To guarantee operation, convert all char arrays to cellstrings with cellstr.

See also: [mustBeNonempty], page 207, [ismember], page 795.
mustBeNegative (x)
Require that input $x$ is negative.
Raise an error if any element of the input $x$ is not negative, as determined by $x<0$.
See also: [mustBeNonnegative], page 207.
mustBeNonempty (x)
Require that input $x$ is nonempty.
Raise an error if the input $x$ is empty, as determined by isempty ( $x$ ).
See also: [mustBeMember], page 207, [mustBeNonzero], page 208, [isempty], page 48.
mustBeNonNan (x)
Require that input $x$ is non-NaN.
Raise an error if any element of the input $x$ is NaN , as determined by isnan ( x ).
See also: [mustBeFinite], page 206, [mustBeNonempty], page 207, [isnan], page 512.
mustBeNonnegative (x)
Require that input $x$ is not negative.
Raise an error if any element of the input $x$ is negative, as determined by $x>=0$.
See also: [mustBeNonzero], page 208, [mustBePositive], page 208.
mustBeNonpositive ( $x$ )
Require that input $x$ is not positive.
Raise an error if any element of the input $x$ is positive, as determined by $x<=0$.
See also: [mustBeNegative], page 207, [mustBeNonzero], page 208.
mustBeNonsparse (x)
Require that input $x$ is not sparse.
Raise an error if the input $x$ is sparse, as determined by issparse ( $x$ ).
See also: [issparse], page 666.
mustBeNonzero (x)
Require that input $x$ is not zero.
Raise an error if any element of the input $x$ is zero, as determined by $x=0$.
See also: [mustBeNonnegative], page 207, [mustBePositive], page 208.
mustBeNumeric (x)
Require that input $x$ is numeric.
Raise an error if the input $x$ is not numeric, as determined by isnumeric ( $x$ ).
See also: [mustBeNumericOrLogical], page 208, [isnumeric], page 67.
mustBeNumericOrLogical (x)
Require that input $x$ is numeric or logical.
Raise an error if the input $x$ is not numeric or logical, as determined by isnumeric (x) || islogical (x).

See also: [mustBeNumeric], page 208, [isnumeric], page 67, [islogical], page 67.
mustBePositive (x)
Require that input $x$ is positive.
Raise an error if any element of the input $x$ is not positive, as determined by $x>0$.
See also: [mustBeNonnegative], page 207, [mustBeNonzero], page 208.
mustBeReal (x)
Require that input $x$ is real.
Raise an error if the input $x$ is not real, as determined by isreal ( $x$ ).
See also: [mustBeFinite], page 206, [mustBeNonNan], page 207, [isreal], page 67.

### 11.9.3 Parsing Arguments

If none of the preceding validation functions is sufficient there is also the class inputParser which can perform extremely complex input checking for functions.
$p=$ inputParser ()
Create object $p$ of the inputParser class.
This class is designed to allow easy parsing of function arguments. The class supports four types of arguments:

1. mandatory (see addRequired);
2. optional (see addOptional);
3. named (see addParameter);
4. switch (see addSwitch).

After defining the function API with these methods, the supplied arguments can be parsed with the parse method and the parsing results accessed with the Results accessor.
inputParser.Parameters
Return list of parameter names already defined.
inputParser.Results
Return structure with argument names as fieldnames and corresponding values.
inputParser.Unmatched
Return structure similar to Results, but for unmatched parameters. See the KeepUnmatched property.
inputParser.UsingDefaults
Return cell array with the names of arguments that are using default values.
inputParser.CaseSensitive $=$ boolean
Set whether matching of argument names should be case sensitive. Defaults to false.
inputParser.FunctionName $=$ name
Set function name to be used in error messages; Defaults to empty string.
inputParser.KeepUnmatched $=$ boolean
Set whether an error should be given for non-defined arguments. Defaults to false. If set to true, the extra arguments can be accessed through Unmatched after the parse method. Note that since Switch and Parameter arguments can be mixed, it is not possible to know the unmatched type. If argument is found unmatched it is assumed to be of the Parameter type and it is expected to be followed by a value.

## inputParser.StructExpand $=$ boolean

Set whether a structure can be passed to the function instead of parameter/value pairs. Defaults to true.

The following example shows how to use this class:

```
function check (varargin)
    p = inputParser (); # create object
    p.FunctionName = "check"; # set function name
    p.addRequired ("pack", @ischar); # mandatory argument
    p.addOptional ("path", pwd(), @ischar); # optional argument
    ## create a function handle to anonymous functions for validators
    val_mat = @(x) isvector (x) && all (x <= 1) && all (x >= 0);
    p.addOptional ("mat", [0 0], val_mat);
    ## create two arguments of type "Parameter"
    val_type = @(x) any (strcmp (x, {"linear", "quadratic"}));
    p.addParameter ("type", "linear", val_type);
    val_verb = @(x) any (strcmp (x, {"low", "medium", "high"}));
    p.addParameter ("tolerance", "low", val_verb);
    ## create a switch type of argument
    p.addSwitch ("verbose");
    p.parse (varargin{:}); # Run created parser on inputs
```

```
            ## the rest of the function can access inputs by using p.Results.
            ## for example, get the tolerance input with p.Results.tolerance
endfunction
check ("mech"); # valid, use defaults for other arguments
check (); # error, one argument is mandatory
check (1); # error, since ! ischar
check ("mech", "~/dev"); # valid, use defaults for other arguments
check ("mech", "~/dev", [0 1 0 0], "type", "linear"); # valid
## following is also valid. Note how the Switch argument type can
## be mixed into or before the Parameter argument type (but it
## must still appear after any Optional argument).
check ("mech", "~/dev", [0 1 0 0], "verbose", "tolerance", "high");
## following returns an error since not all optional arguments,
## 'path' and 'mat', were given before the named argument 'type'.
check ("mech", "~/dev", "type", "linear");
```

Note 1: A function can have any mixture of the four API types but they must appear in a specific order. Required arguments must be first and can be followed by any Optional arguments. Only the Parameter and Switch arguments may be mixed together and they must appear at the end.
Note 2: If both Optional and Parameter arguments are mixed in a function API then once a string Optional argument fails to validate it will be considered the end of the Optional arguments. The remaining arguments will be compared against any Parameter or Switch arguments.

See also: [nargin], page 193, [validateattributes], page 204, [validatestring], page 203, [varargin], page 199.

### 11.10 Function Files

Except for simple one-shot programs, it is not practical to have to define all the functions you need each time you need them. Instead, you will normally want to save them in a file so that you can easily edit them, and save them for use at a later time.

Octave does not require you to load function definitions from files before using them. You simply need to put the function definitions in a place where Octave can find them.

When Octave encounters an identifier that is undefined, it first looks for variables or functions that are already compiled and currently listed in its symbol table. If it fails to find a definition there, it searches a list of directories (the path) for files ending in .m that have the same base name as the undefined identifier. ${ }^{1}$ Once Octave finds a file with a name that matches, the contents of the file are read. If it defines a single function, it is compiled and executed. See Section 11.11 [Script Files], page 223, for more information about how you can define more than one function in a single file.
${ }^{1}$ The '.m' suffix was chosen for compatibility with matlab.

When Octave defines a function from a function file, it saves the full name of the file it read and the time stamp on the file. If the time stamp on the file changes, Octave may reload the file. When Octave is running interactively, time stamp checking normally happens at most once each time Octave prints the prompt. Searching for new function definitions also occurs if the current working directory changes.

Checking the time stamp allows you to edit the definition of a function while Octave is running, and automatically use the new function definition without having to restart your Octave session.

To avoid degrading performance unnecessarily by checking the time stamps on functions that are not likely to change, Octave assumes that function files in the directory tree octavehome/share/octave/version/m will not change, so it doesn't have to check their time stamps every time the functions defined in those files are used. This is normally a very good assumption and provides a significant improvement in performance for the function files that are distributed with Octave.

If you know that your own function files will not change while you are running Octave, you can improve performance by calling ignore_function_time_stamp ("all"), so that Octave will ignore the time stamps for all function files. Passing "system" to this function resets the default behavior.

```
edit name
edit field value
value = edit ("get", field)
value = edit ("get", "all")
```

Edit the named function, or change editor settings.
If edit is called with the name of a file or function as its argument it will be opened in the text editor defined by EDITOR.

- If the function name is available in a file on your path, then it will be opened in the editor. If no file is found, then the m-file variant, ending with ".m", will be considered. If still no file is found, then variants with a leading "@" and then with both a leading "@" and trailing ".m" will be considered.
- If name is the name of a command-line function, then an m-file will be created to contain that function along with its current definition.
- If name.cc is specified, then it will search for name.cc in the path and open it in the editor. If the file is not found, then a new .cc file will be created. If name happens to be an m-file or command-line function, then the text of that function will be inserted into the .cc file as a comment.
- If name.ext is on your path then it will be edited, otherwise the editor will be started with name.ext in the current directory as the filename.
Warning: You may need to clear name before the new definition is available. If you are editing a .cc file, you will need to execute mkoctfile name.cc before the definition will be available.

If edit is called with field and value variables, the value of the control field field will be set to value.
If an output argument is requested and the first input argument is get then edit will return the value of the control field field. If the control field does not exist, edit
will return a structure containing all fields and values. Thus, edit ("get", "all") returns a complete control structure.
The following control fields are used:
'author' This is the name to put after the "\#\# Author:" field of new functions. By default it guesses from the gecos field of the password database.
'email' This is the e-mail address to list after the name in the author field. By default it guesses <\$LOGNAME@\$HOSTNAME>, and if \$HOSTNAME is not defined it uses uname -n. You probably want to override this. Be sure to use the format user@host.
'license'
'gpl' GNU General Public License (default).
'bsd' BSD-style license without advertising clause.
'pd' Public domain.
'"text"' Your own default copyright and license.
Unless you specify 'pd', edit will prepend the copyright statement with "Copyright (C) YYYY Author".
'mode' This value determines whether the editor should be started in async mode (editor is started in the background and Octave continues) or sync mode (Octave waits until the editor exits). Set it to "sync" to start the editor in sync mode. The default is "async" (see [system], page 972).
'editinplace'
Determines whether files should be edited in place, without regard to whether they are modifiable or not. The default is true. Set it to false to have read-only function files automatically copied to 'home', if it exists, when editing them.
'home' This value indicates a directory that system m-files should be copied into before opening them in the editor. The intent is that this directory is also in the path, so that the edited copy of a system function file shadows the original. This setting only has an effect when 'editinplace' is set to false. The default is the empty matrix ([]), which means it is not used. The default in previous versions of Octave was $\sim /$ octave.

See also: [EDITOR], page 32, [path], page 214.
mfilename ()
mfilename ("fullpath")
mfilename ("fullpathext")
Return the name of the currently executing file.
The base name of the currently executing script or function is returned without any extension. If called from outside an m -file, such as the command line, return the empty string.
Given the argument "fullpath", include the directory part of the filename, but not the extension.

Given the argument "fullpathext", include the directory part of the filename and the extension.

See also: [inputname], page 194, [dbstack], page 263.
val = ignore_function_time_stamp ()
old_val = ignore_function_time_stamp (new_val)
Query or set the internal variable that controls whether Octave checks the time stamp on files each time it looks up functions defined in function files.
If the internal variable is set to "system", Octave will not automatically recompile function files in subdirectories of octave-home/share/version/m if they have changed since they were last compiled, but will recompile other function files in the search path if they change.
If set to "all", Octave will not recompile any function files unless their definitions are removed with clear.
If set to "none", Octave will always check time stamps on files to determine whether functions defined in function files need to recompiled.

### 11.10.1 Manipulating the Load Path

When a function is called, Octave searches a list of directories for a file that contains the function declaration. This list of directories is known as the load path. By default the load path contains a list of directories distributed with Octave plus the current working directory. To see your current load path call the path function without any input or output arguments.

It is possible to add or remove directories to or from the load path using addpath and rmpath. As an example, the following code adds ' $/$ /Octave' to the load path.

```
addpath ("~/Octave")
```

After this the directory ' $\sim$ /Octave' will be searched for functions.
addpath (dir1, ...)
addpath (dir1, ..., option)
Add named directories to the function search path.
If option is "-begin" or 0 (the default), prepend the directory name(s) to the current path. If option is "-end" or 1, append the directory name(s) to the current path. Directories added to the path must exist.
In addition to accepting individual directory arguments, lists of directory names separated by pathsep are also accepted. For example:

```
addpath ("dir1:/dir2:~/dir3")
```

The newly added paths appear in the load path in the same order that they appear in the arguments of addpath. When extending the load path to the front, the last path in the list of arguments is added first. When extending the load path to the end, the first path in the list of arguments is added first.
For each directory that is added, and that was not already in the path, addpath checks for the existence of a file named PKG_ADD (note lack of .m extension) and runs it if it exists.

See also: [path], page 214, [rmpath], page 214, [genpath], page 214, [pathdef], page 215, [savepath], page 214, [pathsep], page 215.
genpath (dir)
genpath (dir, skip,...)
Return a path constructed from dir and all its subdirectories.
The path does not include package directories (beginning with '+'), old-style class directories (beginning with ' $@$ '), private directories, or any subdirectories of these types.

If additional string parameters are given, the resulting path will exclude directories with those names.

See also: [path], page 214, [addpath], page 213.
rmpath (dir1, ...)
Remove dir1, ... from the current function search path.
In addition to accepting individual directory arguments, lists of directory names separated by pathsep are also accepted. For example:

```
rmpath ("dir1:/dir2:~/dir3")
```

For each directory that is removed, rmpath checks for the existence of a file named PKG_DEL (note lack of .m extension) and runs it if it exists.

See also: [path], page 214, [addpath], page 213, [genpath], page 214, [pathdef], page 215, [savepath], page 214, [pathsep], page 215.

```
savepath ()
savepath (file)
status = savepath (...)
```

Save the unique portion of the current function search path that is not set during Octave's initialization process to file.
If file is omitted, Octave looks in the current directory for a project-specific . octaverc file in which to save the path information. If no such file is present then the user's configuration file $\sim /$. octaverc is used.
If successful, savepath returns 0 .
The savepath function makes it simple to customize a user's configuration file to restore the working paths necessary for a particular instance of Octave. Assuming no filename is specified, Octave will automatically restore the saved directory paths from the appropriate .octaverc file when starting up. If a filename has been specified then the paths may be restored manually by calling source file.
See also: [path], page 214, [addpath], page 213, [rmpath], page 214, [genpath], page 214, [pathdef], page 215.
path ()
str = path ()
str $=$ path (path1, ...)
Modify or display Octave's load path.

If nargin and nargout are zero, display the elements of Octave's load path in an easy to read format.

If nargin is zero and nargout is greater than zero, return the current load path.
If nargin is greater than zero, concatenate the arguments, separating them with pathsep. Set the internal search path to the result and return it.
No checks are made for duplicate elements.
See also: [addpath], page 213, [rmpath], page 214, [genpath], page 214, [pathdef], page 215, [savepath], page 214, [pathsep], page 215.
val = pathdef ()
Return the default path for Octave.
The path information is extracted from one of four sources. The possible sources, in order of preference, are:

1. .octaverc
2. ~/.octaverc
3. <OCTAVE_HOME>/.../<version>/m/startup/octaverc
4. Octave's path prior to changes by any octaverc file.

See also: [path], page 214, [addpath], page 213, [rmpath], page 214, [genpath], page 214, [savepath], page 214.
val = pathsep ()
Query the character used to separate directories in a path.
See also: [filesep], page 960.
rehash ()
Reinitialize Octave's load path directory cache.
fname $=$ file_in_loadpath (file)
fname $=$ file_in_loadpath (file, "all")
Return the absolute name of file if it can be found in the list of directories specified by path.
If no file is found, return an empty character string.
When file is already an absolute name, the name is checked against the file system instead of Octave's loadpath. In this case, if file exists it will be returned in fname, otherwise an empty string is returned.
If the first argument is a cell array of strings, search each directory of the loadpath for element of the cell array and return the first that matches.
If the second optional argument "all" is supplied, return a cell array containing the list of all files that have the same name in the path. If no files are found, return an empty cell array.

See also: [file_in_path], page 960, [dir_in_loadpath], page 216, [path], page 214.
restoredefaultpath ()
Restore Octave's path to its initial state at startup.
See also: [path], page 214, [addpath], page 213, [rmpath], page 214, [genpath], page 214, [pathdef], page 215, [savepath], page 214, [pathsep], page 215.
command_line_path ()
Return the command line path variable.
See also: [path], page 214, [addpath], page 213, [rmpath], page 214, [genpath], page 214, [pathdef], page 215, [savepath], page 214, [pathsep], page 215.
dirname = dir_in_loadpath (dir)
dirname = dir_in_loadpath (dir, "all")
Return the absolute name of the loadpath element matching dir if it can be found in the list of directories specified by path.
If no match is found, return an empty character string.
The match is performed at the end of each path element. For example, if dir is "foo/bar", it matches the path element "/some/dir/foo/bar", but not "/some/dir/foo/bar/baz" "/some/dir/allfoo/bar". When dir is an absolute name, rather than just a path fragment, it is matched against the file system instead of Octave's loadpath. In this case, if dir exists it will be returned in dirname, otherwise an empty string is returned.
If the optional second argument is supplied, return a cell array containing all name matches rather than just the first.
See also: [file_in_path], page 960, [file_in_loadpath], page 215, [path], page 214.

```
current_encoding = dir_encoding (dir)
prev_encoding = dir_encoding (dir, encoding)
dir_encoding (...)
```

Set and query the encoding that is used for reading m-files in dir.
That encoding overrides the (globally set) m-file encoding.
The string $D I R$ must match the form how the directory would appear in the load path.
The encoding must be a valid encoding identifier or "delete". In the latter case, the (globally set) m-file encoding will be used for the given dir.

The currently or previously used encoding is returned in current_encoding or prev_encoding, respectively. The output argument must be explicitly requested.

The directory encoding is automatically read from the file .oct-config when a new path is added to the load path (for example with addpath). To set the encoding for all files in the same folder, that file must contain a line starting with "encoding=" followed by the encoding identifier.

For example to set the file encoding for all files in the same folder to ISO 8859-1 (Latin-1), create a file .oct-config with the following content:

```
encoding=iso8859-1
```

If the file encoding is changed after the files have already been parsed, the files have to be parsed again for that change to take effect. That can be done with the command clear all.
See also: [addpath], page 213, [path], page 214.

### 11.10.2 Subfunctions

A function file may contain secondary functions called subfunctions. These secondary functions are only visible to the other functions in the same function file. For example, a file f.m containing

```
function f ()
    printf ("in f, calling g\n");
    g ()
endfunction
function g ()
    printf ("in g, calling h\n");
    h ()
endfunction
function h ()
    printf ("in h\n")
endfunction
```

defines a main function $f$ and two subfunctions. The subfunctions $g$ and $h$ may only be called from the main function $f$ or from the other subfunctions, but not from outside the file $f . m$.
localfunctions ()
Return a list of all local functions, i.e., subfunctions, within the current file.
The return value is a column cell array of function handles to all local functions accessible from the function from which localfunctions is called. Nested functions are not included in the list.
If the call is from the command line, an anonymous function, or a script, the return value is an empty cell array.
See also: [functions], page 234.

### 11.10.3 Private Functions

In many cases one function needs to access one or more helper functions. If the helper function is limited to the scope of a single function, then subfunctions as discussed above might be used. However, if a single helper function is used by more than one function, then this is no longer possible. In this case the helper functions might be placed in a subdirectory, called "private", of the directory in which the functions needing access to this helper function are found.

As a simple example, consider a function func1, that calls a helper function func2 to do much of the work. For example:

```
function y = func1 (x)
    y = func2 (x);
endfunction
```

Then if the path to func1 is <directory>/func1.m, and if func2 is found in the directory <directory>/private/func2.m, then func2 is only available for use of the functions, like func1, that are found in <directory>.

### 11.10.4 Nested Functions

Nested functions are similar to subfunctions in that only the main function is visible outside the file. However, they also allow for child functions to access the local variables in their parent function. This shared access mimics using a global variable to share information but a global variable which is not visible to the rest of Octave. As a programming strategy, sharing data this way can create code which is difficult to maintain. It is recommended to use subfunctions in place of nested functions when possible.

As a simple example, consider a parent function foo, that calls a nested child function bar, with a shared variable $x$.

```
function y = foo ()
    x = 10;
    bar ();
    y = x;
    function bar ()
        x = 20;
    endfunction
endfunction
foo ()
    # 20
```

Notice that there is no special syntax for sharing $x$. This can lead to problems with accidental variable sharing between a parent function and its child. While normally variables are inherited, child function parameters and return values are local to the child function.

Now consider the function foobar that uses variables $x$ and $y$. foobar calls a nested function foo which takes $x$ as a parameter and returns $y$. foo then calls bat which does some computation.

```
function z = foobar ()
    x = 0;
    y = 0;
    z = foo (5);
    z += x + y;
    function y = foo (x)
        y = x + bat ();
        function z = bat ()
            z = x;
        endfunction
    endfunction
endfunction
foobar ()
    => 10
```

It is important to note that the $x$ and $y$ in foobar remain zero, as in foo they are a return value and parameter respectively. The $x$ in bat refers to the $x$ in foo.

Variable inheritance leads to a problem for eval and scripts. If a new variable is created in a parent function, it is not clear what should happen in nested child functions. For example, consider a parent function foo with a nested child function bar:

```
function y = foo (to_eval)
    bar ();
    eval (to_eval);
    function bar ()
        eval ("x = 100;");
        eval ("y = x;");
    endfunction
endfunction
foo ("x = 5;")
        => error: can not add variable "x" to a static workspace
foo ("y = 10;")
    => 10
foo("")
        # 100
```

The parent function foo is unable to create a new variable $x$, but the child function bar was successful. Furthermore, even in an eval statement $y$ in bar is the same $y$ as in its parent function foo. The use of eval in conjunction with nested functions is best avoided.

As with subfunctions, only the first nested function in a file may be called from the outside. Inside a function the rules are more complicated. In general a nested function may call:

0 . Globally visible functions

1. Any function that the nested function's parent can call
2. Sibling functions (functions that have the same parents)
3. Direct children

As a complex example consider a parent function ex_top with two child functions, ex_a and ex_b. In addition, ex_a has two more child functions, ex_aa and ex_ab. For example:

```
function ex_top ()
    ## Can call: ex_top, ex_a, and ex_b
    ## Can NOT call: ex_aa and ex_ab
    function ex_a ()
            ## Can call everything
            function ex_aa ()
                ## Can call everything
            endfunction
            function ex_ab ()
                ## Can call everything
            endfunction
    endfunction
    function ex_b ()
            ## Can call: ex_top, ex_a, and ex_b
            ## Can NOT call: ex_aa and ex_ab
    endfunction
endfunction
```


### 11.10.5 Overloading and Autoloading

Functions can be overloaded to work with different input arguments. For example, the operator ' + ' has been overloaded in Octave to work with single, double, uint8, int32, and many other arguments. The preferred way to overload functions is through classes and object oriented programming (see Section 34.4.1 [Function Overloading], page 899). Occasionally, however, one needs to undo user overloading and call the default function associated with a specific type. The builtin function exists for this purpose.

## [...] = builtin (f,...)

Call the base function $f$ even if $f$ is overloaded to another function for the given type signature.
This is normally useful when doing object-oriented programming and there is a requirement to call one of Octave's base functions rather than the overloaded one of a new class.
A trivial example which redefines the sin function to be the cos function shows how builtin works.

```
\(\sin (0)\)
    \(\Rightarrow 0\)
function \(y=\sin (x), y=\cos (x) ;\) endfunction
\(\sin (0)\)
    \(\Rightarrow 1\)
builtin ("sin", 0)
    \(\Rightarrow 0\)
```

A single dynamically linked file might define several functions. However, as Octave searches for functions based on the functions filename, Octave needs a manner in which to find each of the functions in the dynamically linked file. On operating systems that support symbolic links, it is possible to create a symbolic link to the original file for each of the functions which it contains.

However, there is at least one well known operating system that doesn't support symbolic links. Making copies of the original file for each of the functions is undesirable as it increases the amount of disk space used by Octave. Instead Octave supplies the autoload function, that permits the user to define in which file a certain function will be found.

```
autoload_map = autoload ()
autoload (function, file)
autoload (..., "remove")
```

Define function to autoload from file.
The second argument, file, should be an absolute filename or a file name in the same directory as the function or script from which the autoload command was run. file should not depend on the Octave load path.
Normally, calls to autoload appear in PKG_ADD script files that are evaluated when a directory is added to Octave's load path. To avoid having to hardcode directory names in file, if file is in the same directory as the PKG_ADD script then

```
autoload ("foo", "bar.oct");
```

will load the function foo from the file bar.oct. The above usage when bar.oct is not in the same directory, or usages such as

```
autoload ("foo", file_in_loadpath ("bar.oct"))
```

are strongly discouraged, as their behavior may be unpredictable.
With no arguments, return a structure containing the current autoload map.
If a third argument "remove" is given, the function is cleared and not loaded anymore during the current Octave session.

See also: $\left[P K G \_A D D\right]$, page 1004 .

### 11.10.6 Function Locking

It is sometime desirable to lock a function into memory with the mlock function. This is typically used for dynamically linked functions in oct-files or mex-files that contain some initialization, and it is desirable that calling clear does not remove this initialization.

As an example,

```
function my_function ()
    mlock ();
endfunction
```

prevents my_function from being removed from memory after it is called, even if clear is called. It is possible to determine if a function is locked into memory with the mislocked, and to unlock a function with munlock, which the following code illustrates.

```
my_function ();
mislocked ("my_function")
a ans = 1
munlock ("my_function");
mislocked ("my_function")
m ans = 0
```

A common use of mlock is to prevent persistent variables from being removed from memory, as the following example shows:

```
function count_calls ()
    mlock ();
    persistent calls = 0;
    printf ("count_calls() has been called %d times\n", ++calls);
endfunction
count_calls ();
\dashv count_calls() has been called 1 times
clear count_calls
count_calls ();
    \dashvcount_calls() has been called 2 times
```

mlock might also be used to prevent changes to an m-file, such as in an external editor, from having any effect in the current Octave session; A similar effect can be had with the ignore_function_time_stamp function.
mlock ()
Lock the current function into memory so that it can't be removed with clear.
See also: [munlock], page 222, [mislocked], page 222, [persistent], page 140, [clear], page 145 .
munlock ()
munlock ( $f \subset n$ )
Unlock the named function fcn so that it may be removed from memory with clear. If no function is named then unlock the current function.

See also: [mlock], page 222, [mislocked], page 222, [persistent], page 140, [clear], page 145.
mislocked ()
mislocked (fcn)
Return true if the named function $f c n$ is locked in memory.

If no function is named then return true if the current function is locked.
See also: [mlock], page 222, [munlock], page 222, [persistent], page 140, [clear], page 145 .

### 11.10.7 Function Precedence

Given the numerous different ways that Octave can define a function, it is possible and even likely that multiple versions of a function, might be defined within a particular scope. The precedence of which function will be used within a particular scope is given by

1. Subfunction A subfunction with the required function name in the given scope.
2. Private function A function defined within a private directory of the directory which contains the current function.
3. Class constructor A function that constructs a user class as defined in chapter Chapter 34 [Object Oriented Programming], page 889.
4. Class method An overloaded function of a class as in chapter Chapter 34 [Object Oriented Programming], page 889.
5. Command-line Function A function that has been defined on the command-line.
6. Autoload function A function that is marked as autoloaded with See [autoload], page 221.
7. A Function on the Path A function that can be found on the users load-path. There can also be Oct-file, mex-file or m-file versions of this function and the precedence between these versions are in that order.
8. Built-in function A function that is a part of core Octave such as numel, size, etc.

### 11.11 Script Files

A script file is a file containing (almost) any sequence of Octave commands. It is read and evaluated just as if you had typed each command at the Octave prompt, and provides a convenient way to perform a sequence of commands that do not logically belong inside a function.

Unlike a function file, a script file must not begin with the keyword function. If it does, Octave will assume that it is a function file, and that it defines a single function that should be evaluated as soon as it is defined.

A script file also differs from a function file in that the variables named in a script file are not local variables, but are in the same scope as the other variables that are visible on the command line.

Even though a script file may not begin with the function keyword, it is possible to define more than one function in a single script file and load (but not execute) all of them at once. To do this, the first token in the file (ignoring comments and other white space) must be something other than function. If you have no other statements to evaluate, you can use a statement that has no effect, like this:

```
# Prevent Octave from thinking that this
# is a function file:
```

1 ;
\# Define function one:

```
function one ()
```

    ...
    To have Octave read and compile these functions into an internal form, you need to make sure that the file is in Octave's load path (accessible through the path function), then simply type the base name of the file that contains the commands. (Octave uses the same rules to search for script files as it does to search for function files.)

If the first token in a file (ignoring comments) is function, Octave will compile the function and try to execute it, printing a message warning about any non-whitespace characters that appear after the function definition.

Note that Octave does not try to look up the definition of any identifier until it needs to evaluate it. This means that Octave will compile the following statements if they appear in a script file, or are typed at the command line,

```
# not a function file:
1;
function foo ()
    do_something ();
endfunction
function do_something ()
    do_something_else ();
endfunction
```

even though the function do_something is not defined before it is referenced in the function foo. This is not an error because Octave does not need to resolve all symbols that are referenced by a function until the function is actually evaluated.

Since Octave doesn't look for definitions until they are needed, the following code will always print 'bar = 3' whether it is typed directly on the command line, read from a script file, or is part of a function body, even if there is a function or script file called bar.m in Octave's path.

```
eval ("bar = 3");
bar
```

Code like this appearing within a function body could fool Octave if definitions were resolved as the function was being compiled. It would be virtually impossible to make Octave clever enough to evaluate this code in a consistent fashion. The parser would have to be able to perform the call to eval at compile time, and that would be impossible unless all the references in the string to be evaluated could also be resolved, and requiring that would be too restrictive (the string might come from user input, or depend on things that are not known until the function is evaluated).

Although Octave normally executes commands from script files that have the name file.m, you can use the function source to execute commands from any file.
source (file)
source (file, context)
Parse and execute the contents of file.
Without specifying context, this is equivalent to executing commands from a script file, but without requiring the file to be named file.m or to be on the execution path.
Instead of the current context, the script may be executed in either the context of the function that called the present function ("caller"), or the top-level context ("base").
See also: [run], page 177 .

### 11.11.1 Publish Octave Script Files

The function publish provides a dynamic possibility to document your script file. Unlike static documentation, publish runs the script file, saves any figures and output while running the script, and presents them alongside static documentation in a desired output format. The static documentation can make use of Section 11.11.2 [Publishing Markup], page 227 , to enhance and customize the output.

```
publish (file)
publish (file, output_format)
publish (file, option1, value1,...)
publish (file, options)
output_file = publish (file,...)
```

Generate a report from the Octave script file file in one of several output formats.
The generated reports interpret Publishing Markup in section comments, which is explained in detail in the GNU Octave manual. Section comments are comment blocks that start with a line with double comment character.
Assume the following example, using some Publishing Markup, to be the contents of the script file pub_example.m:

```
## Headline title
#
# Some *bold*, _italic_, or |monospaced| Text with
# a <https://www.octave.org link to *GNU Octave*>.
##
# "Real" Octave commands to be evaluated
sombrero ()
%% MATLAB comment style ('%') is supported as well
%
% * Bulleted list item 1
%* Bulleted list item 2
%
% # Numbered list item 1
% # Numbered list item 2
```

To publish this script file, type publish ("pub_example.m").

When called with one input argument, a HTML report is generated in a subdirectory html relative to the current working directory. Any Octave commands in pub_ example.m are evaluated in a separate context and any figures created while executing the script file are included in the report.
Using publish (file, output_format) is equivalent to the function call using a structure

```
options.format = output_format;
publish (file, options)
```

which is described below. The same holds for using option/value pairs

```
options.option1 = value1;
publish (file, options)
```

The structure options can have the following field names. If a field name is not specified, the default value is used:

- 'format' - Output format of the published script file, one of
'html' (default), 'doc', 'latex', 'ppt', 'pdf', or 'xml'.
The output formats 'doc', 'ppt', and 'xml' are not currently supported. To generate a 'doc' report, open a generated 'html' report with your office suite.
In Octave custom formats are supported by implementing all callback subfunctions in a function file named '__publish_<custom format>_output__.m'. To obtain a template for the HTML format type:

```
edit (fullfile (fileparts (which ("publish")), ...
    "private", "__publish_html_output__.m"))
```

- 'outputDir' - Full path of the directory where the generated report will be located. If no directory is given, the report is generated in a subdirectory html relative to the current working directory.
- 'stylesheet' - Not supported, only for matlab compatibility.
- 'createThumbnail' - Not supported, only for matlab compatibility.
- 'figureSnapMethod' - Not supported, only for matlab compatibility.
- 'imageFormat' - Desired format for any images produced while evaluating the code. The allowed image formats depend on the output format:
- 'html', 'xml' - 'png' (default), any image format supported by Octave
- 'latex' - 'epsc2' (default), any image format supported by Octave
- 'pdf' - 'jpg' (default) or 'bmp', note matlab uses 'bmp' as default
- 'doc' or 'ppt' - 'png' (default), 'jpg', 'bmp', or 'tiff'
- 'maxWidth' and 'maxHeight' - Maximum width (height) of the produced images in pixels. An empty value means no restriction. Both values must be set in order for the option to work properly.
'[]' (default), integer value $\geq 0$
- 'useNewFigure' - Use a new figure window for figures created by the evaluated code. This avoids side effects with already opened figure windows.
'true' (default) or 'false'
- 'evalCode' - Evaluate code of the Octave source file 'true' (default) or 'false'
- 'catchError' - Catch errors while evaluating code and continue 'true' (default) or 'false'
- 'codeToEvaluate' - Octave commands that should be evaluated prior to publishing the script file. These Octave commands do not appear in the generated report.
- 'maxOutputLines' - Maximum number of output lines from code evaluation which are included in output.
'Inf' (default) or integer value >0
- 'showCode' - Show the evaluated Octave commands in the generated report
'true' (default) or 'false'
The option output output_file is a string with path and file name of the generated report.
See also: [grabcode], page 227.
The counterpart to publish is grabcode:
grabcode (url)
grabcode (filename)
code_str $=$ grabcode (...)
Grab the code from a report created by the publish function.
The grabbed code inside the published report must be enclosed by the strings '\#\#\#\#\# SOURCE BEGIN \#\#\#\#\#' and '\#\#\#\#\# SOURCE END \#\#\#\#\#'. The publish function creates this format automatically.
If no return value is requested the code is saved to a temporary file and opened in the default editor. NOTE: The temporary file must be saved under a new or the code will be lost.
If an output is requested the grabbed code will be returned as string code_str.
Example:
publish ("my_script.m");
grabcode ("html/my_script.html");
The example above publishes my_script.m to the default location html/my_ script.html. Next, the published Octave script is grabbed to edit its content in a new temporary file.

See also: [publish], page 225.

### 11.11.2 Publishing Markup

### 11.11.2.1 Using Publishing Markup in Script Files

To use Publishing Markup, start by typing '\#\#' or '\%\%' at the beginning of a new line. For mATLAB compatibility $6 \% \% \%$ ' is treated the same way as $6 \% \%$ '.

The lines following '\#\#' or '\%\%' start with one of either '\#' or '\%' followed by at least one space. These lines are interpreted as section. A section ends at the first line not starting with '\#' or ' $\%$ ', or when the end of the document is reached.

A section starting in the first line of the document, followed by another start of a section that might be empty, is interpreted as a document title and introduction text.

See the example below for clarity:

```
%% Headline title
%
% Some *bold*, _italic_, or |monospaced| Text with
% a <https://www.octave.org link to GNU Octave>.
%%
# "Real" Octave commands to be evaluated
sombrero ()
## Octave comment style supported as well
#
# * Bulleted list item 1
# * Bulleted list item 2
#
# # Numbered list item 1
# # Numbered list item 2
```


### 11.11.2.2 Text Formatting

Basic text formatting is supported inside sections, see the example given below:

```
##
# *bold*, _italic_, or |monospaced| Text
```

Additionally two trademark symbols are supported, just embrace the letters ' TM ' or ' R '.

```
##
# (TM) or (R)
```


### 11.11.2.3 Sections

A section is started by typing '\#\#' or '\%\%' at the beginning of a new line. A section title can be provided by writing it, separated by a space, in the first line after '\#\#' or '\%\%'. Without a section title, the section is interpreted as a continuation of the previous section. For mATLAB compatibility ' $\% \% \%$ ' is treated the same way as ' $\% \%$ '.

```
some_code ();
## Section 1
#
## Section 2
some_code ();
##
# Still in section 2
some_code ();
%%% Section 3
%
%
```


### 11.11.2.4 Preformatted Code

To write preformatted code inside a section, indent the code by three spaces after '\#' at the beginning of each line and leave the lines above and below the code blank, except for ' $\#$ ' at the beginning of those lines.

```
##
# This is a syntax highlighted for-loop:
#
# for i = 1:5
# disp (i);
# endfor
#
# And more usual text.
```


### 11.11.2.5 Preformatted Text

To write preformatted text inside a section, indent the code by two spaces after ' $\#$ ' at the beginning of each line and leave the lines above and below the preformatted text blank, except for ' $\#$ ' at the beginning of those lines.

```
##
# This following text is preformatted:
#
# "To be, or not to be: that is the question:
# Whether 'tis nobler in the mind to suffer
# The slings and arrows of outrageous fortune,
# Or to take arms against a sea of troubles,
# And by opposing end them? To die: to sleep;"
#
# --"Hamlet" by W. Shakespeare
```


### 11.11.2.6 Bulleted Lists

To create a bulleted list, type

```
##
#
# * Bulleted list item 1
# * Bulleted list item 2
#
```

to get output like

- Bulleted list item 1
- Bulleted list item 2

Notice the blank lines, except for the '\#' or '\%' before and after the bulleted list!

### 11.11.2.7 Numbered Lists

To create a numbered list, type

```
##
#
# # Numbered list item 1
# # Numbered list item 2
#
```

to get output like

1. Numbered list item 1
2. Numbered list item 2

Notice the blank lines, except for the '\#' or ' $\%$ ' before and after the numbered list!

### 11.11.2.8 Including File Content

To include the content of an external file, e.g., a file called 'my_function.m' at the same location as the published Octave script, use the following syntax to include it with Octave syntax highlighting.

Alternatively, you can write the full or relative path to the file.

```
##
#
# <include>my_function.m</include>
#
# <include>/full/path/to/my_function.m</include>
#
# <include>../relative/path/to/my_function.m</include>
#
```


### 11.11.2.9 Including Graphics

To include external graphics, e.g., a graphic called 'my_graphic.png' at the same location as the published Octave script, use the following syntax.

Alternatively, you can write the full path to the graphic.

```
##
#
# <<my_graphic.png>>
#
# <</full/path/to/my_graphic.png>>
#
# <<../relative/path/to/my_graphic.png>>
#
```


### 11.11.2.10 Including URLs

Basically, a URL is written between an opening '<' and a closing ‘>' angle.

```
##
# <https://www.octave.org>
```

Text that is within these angles and separated by at least one space from the URL is a displayed text for the link.

```
##
# <https://www.octave.org GNU Octave>
```

A link starting with '<octave:' followed by the name of a GNU Octave function, optionally with a displayed text, results in a link to the online GNU Octave documentations function index.

```
##
# <octave:DISP The display function>
```


### 11.11.2.11 Mathematical Equations

One can insert $\mathrm{LT}_{\mathrm{E}} \mathrm{X}$ inline math, surrounded by single ' $\$$ ' signs, or displayed math, surrounded by double ' $\$ \$$ ' signs, directly inside sections.

```
##
# Some shorter inline equation $e^{ix} = \cos x + i\sin x$.
#
# Or more complicated formulas as displayed math:
# $$e^x = \lim_{n->\infty}\left(1+\dfrac{x}{n}\right)^{n}.$$
```


### 11.11.2.12 HTML Markup

If the published output is a HTML report, you can insert HTML markup, that is only visible in this kind of output.

```
##
# <html>
# <table style="border:1px solid black;">
# <tr><td>1</td><td>2</td></tr>
# <tr><td>3</td><td>3</td></tr>
# </html>
```


### 11.11.2.13 LaTeX Markup

If the published output is a ${ }_{A T} T_{E} \mathrm{X}$ or PDF report, you can insert $\mathrm{IAT}_{\mathrm{E}} \mathrm{X}$ markup, that is only visible in this kind of output.

```
##
# <latex>
# Some output only visible in LaTeX or PDF reports.
# \begin{equation}
# e^x = \lim\limits_{n->\infty}\left(1+\dfrac{x}{n}\right)^{n}
# \end{equation}
# </latex>
```


### 11.11.3 Jupyter Notebooks

Jupyter notebooks are one popular technique for displaying code, text, and graphical output together in a comprehensive manner. Octave can publish results to a Jupyter notebook with the function jupyter_notebook.
notebook = jupyter_notebook (notebook_filename)
notebook = jupyter_notebook (notebook_filename, options)
Run and fill the Jupyter Notebook in file notebook_filename from within GNU Octave.
Both text and graphical Octave outputs are supported.
This class has a public property notebook which is a structure representing the JSONdecoded Jupyter Notebook. This property is intentionally public to enable advanced notebook manipulations.

Note: Jupyter Notebook versions (nbformat) lower than 4.0 are not supported.
The optional second argument options is a struct with fields:

- tmpdir to set the temporary working directory.
\%plot magic is supported with the following settings:
- "\%plot -f <format>" or "\%plot -format <format>": specifies the image storage format. Supported formats are:
- PNG (default)
- SVG (Note: If SVG images do not appear in the notebook, it is most likely related to Jupyter Notebook security mechanisms and explicitly "trusting" them will be necessary).
- JPG
- "\%plot -r <number>" or "\%plot -resolution <number>": specifies the image resolution.
- "\%plot -w <number>" or "\%plot -width <number>": specifies the image width.
- "\%plot -h <number>" or "\%plot -height <number>": specifies the image height.

Examples:

```
## Run all cells and generate the filled notebook
## Instantiate an object from a notebook file
notebook = jupyter_notebook ("myNotebook.ipynb");
## Run the code and embed the results in the notebook property
notebook.run_all ();
## Generate a new notebook by overwriting the original notebook
notebook.generate_notebook ("myNotebook.ipynb");
## Run just the second cell and generate the filled notebook
## Instantiate an object from a notebook file
notebook = jupyter_notebook ("myNotebook.ipynb");
## Run the code and embed the results in the notebook property
notebook.run (2)
## Generate a new notebook in a new file
notebook.generate_notebook ("myNewNotebook.ipynb");
## Generate an Octave script from a notebook
## Instantiate an object from a notebook file
notebook = jupyter_notebook ("myNotebook.ipynb");
## Generate the Octave script
notebook.generate_octave_script ("jup_script.m");
```

See also: [jsondecode], page 106, [jsonencode], page 104.

### 11.12 Function Handles and Anonymous Functions

It can be very convenient store a function in a variable so that it can be passed to a different function. For example, a function that performs numerical minimization needs access to the function that should be minimized.

### 11.12.1 Function Handles

A function handle is a pointer to another function and is defined with the syntax @function-name
For example,

$$
f=@ \sin ;
$$

creates a function handle called $f$ that refers to the function sin.
Function handles are used to call other functions indirectly, or to pass a function as an argument to another function like quad or fsolve. For example:

```
f = @sin;
quad (f, 0, pi)
    => 2
```

You may use feval to call a function using function handle, or simply write the name of the function handle followed by an argument list. If there are no arguments, you must use an empty argument list '()'. For example:

```
    f = @sin;
    feval (f, pi/4)
        => 0.70711
    f (pi/4)
        => 0.70711
```

is_function_handle (x)

Return true if $x$ is a function handle.
See also: [isa], page 41, [typeinfo], page 41, [class], page 41, [functions], page 234.
$s=$ functions (fcn_handle)
Return a structure containing information about the function handle fcn_handle.
The structure $s$ always contains these three fields:
function The function name. For an anonymous function (no name) this will be the actual function definition.
type Type of the function. anonymous

The function is anonymous.
private The function is private.
overloaded
The function overloads an existing function.
simple The function is a built-in or m-file function.
subfunction
The function is a subfunction within an m-file.
nested The function is nested.
file The m-file that will be called to perform the function. This field is empty for anonymous and built-in functions.

In addition, some function types may return more information in additional fields.
Warning: functions is provided for debugging purposes only. Its behavior may change in the future and programs should not depend on any particular output format.

See also: [func2str], page 234, [str2func], page 234.
func2str (fcn_handle)
Return a string containing the name of the function referenced by the function handle fcn_handle.
See also: [str2func], page 234, [functions], page 234.
str2func (fcn_name)
Return a function handle constructed from the string fcn_name.
Previous versions of Octave accepted an optional second argument, "global", that caused str2func to ignore locally visible functions. This option is no longer supported.
See also: [func2str], page 234, [functions], page 234.
vars $=$ symvar (str)
Identify the symbolic variable names in the string str.
Common constant names such as i, j, pi, Inf and Octave functions such as sin or plot are ignored.
Any names identified are returned in a cell array of strings. The array is empty if no variables were found.

Example:

```
symvar ("x^2 + y^2 == 4")
=> {
    [1,1] = x
    [2,1] = y
    }
```


### 11.12.2 Anonymous Functions

Anonymous functions are defined using the syntax

```
@(argument-list) expression
```

Any variables that are not found in the argument list are inherited from the enclosing scope. Anonymous functions are useful for creating simple unnamed functions from expressions or for wrapping calls to other functions to adapt them for use by functions like quad. For example,

```
f = @(x) x.`2;
quad (f, 0, 10)
    # 333.33
```

creates a simple unnamed function from the expression $x . \wedge 2$ and passes it to quad,

```
quad (@(x) sin (x), 0, pi)
    =>2
```

wraps another function, and

```
a = 1;
b = 2;
quad (@(x) betainc (x, a, b), 0, 0.4)
    # 0.13867
```

adapts a function with several parameters to the form required by quad. In this example, the values of $a$ and $b$ that are passed to betainc are inherited from the current environment.

Note that for performance reasons it is better to use handles to existing Octave functions, rather than to define anonymous functions which wrap an existing function. The integration of $\sin (x)$ is 5 X faster if the code is written as

```
quad (@sin, 0, pi)
```

rather than using the anonymous function @ (x) sin (x). There are many operators which have functional equivalents that may be better choices than an anonymous function. Instead of writing
$f=@(x, y) x+y$
this should be coded as
f = @plus

See Section 34.4.2 [Operator Overloading], page 900, for a list of operators which also have a functional form.

### 11.13 Command Syntax and Function Syntax

Additionally to the function syntax described above (i.e., calling a function like fun (arg1, $\arg 2, \ldots$ ) , a function can be called using command syntax (for example, calling a function like fun $\arg 1 \arg 2 \ldots$. . . In that case, all arguments are passed to the function as strings. For example,
my_command hello world
is equivalent to
my_command ("hello", "world")
The general form of a command call is
cmdname arg1 arg2 ...
which translates directly to

```
cmdname ("arg1", "arg2", ...)
```

If an argument including spaces should be passed to a function in command syntax, (double-)quotes can be used. For example,

```
my_command "first argument" "second argument"
```

is equivalent to

```
my_command ("first argument", "second argument")
```

Any function can be used as a command if it accepts string input arguments. For example:

```
toupper lower_case_arg
    => ans = LOWER_CASE_ARG
```

Since the arguments are passed as strings to the corresponding function, it is not possible to pass input arguments that are stored in variables. In that case, a command must be called using the function syntax. For example:

```
strvar = "hello world";
toupper strvar
    a ans = STRVAR
toupper (strvar)
    a ans = HELLO WORLD
```

Additionally, the return values of functions cannot be assigned to variables using the command syntax. Only the first return argument is assigned to the built-in variable ans. If the output argument of a command should be assigned to a variable, or multiple output arguments of a function should be returned, the function syntax must be used.

### 11.14 Organization of Functions Distributed with Octave

Many of Octave's standard functions are distributed as function files. They are loosely organized by topic, in subdirectories of octave-home/share/octave/version/m, to make it easier to find them.

The following is a list of all the function file subdirectories, and the types of functions you will find there.
@ftp Class functions for the FTP object.
+containers
Package for the containers classes.
audio Functions for playing and recording sounds.
deprecated
Out-of-date functions which will eventually be removed from Octave.
elfun Elementary functions, principally trigonometric.
general Miscellaneous matrix manipulations, like flipud, rot90, and triu, as well as other basic functions, like ismatrix, narginchk, etc.
geometry Functions related to Delaunay triangulation.
gui Functions for GUI elements like dialog, message box, etc.
help Functions for Octave's built-in help system.
image Image processing tools. These functions require the X Window System.
io Input-output functions.
java Functions related to the Java integration.
linear-algebra
Functions for linear algebra.
miscellaneous
Functions that don't really belong anywhere else.
ode Functions to solve ordinary differential equations (ODEs).
optimization
Functions related to minimization, optimization, and root finding.
path Functions to manage the directory path Octave uses to find functions.
pkg Package manager for installing external packages of functions in Octave.
plot Functions for displaying and printing two- and three-dimensional graphs.
polynomial
Functions for manipulating polynomials.
prefs Functions implementing user-defined preferences.
set Functions for creating and manipulating sets of unique values.
signal Functions for signal processing applications.
sparse Functions for handling sparse matrices.
specfun Special functions such as bessel or factor.
special-matrix
Functions that create special matrix forms such as Hilbert or Vandermonde matrices.
startup Octave's system-wide startup file.
statistics
Statistical functions.
strings Miscellaneous string-handling functions.
testfun Functions for performing unit tests on other functions.
time Functions related to time and date processing.

## 12 Errors and Warnings

Octave includes several functions for printing error and warning messages. When you write functions that need to take special action when they encounter abnormal conditions, you should print the error messages using the functions described in this chapter.

Since many of Octave's functions use these functions, it is also useful to understand them, so that errors and warnings can be handled.

### 12.1 Handling Errors

An error is something that occurs when a program is in a state where it doesn't make sense to continue. An example is when a function is called with too few input arguments. In this situation the function should abort with an error message informing the user of the lacking input arguments.

Since an error can occur during the evaluation of a program, it is very convenient to be able to detect that an error occurred, so that the error can be fixed. This is possible with the try statement described in Section 10.9 [The try Statement], page 188.

### 12.1.1 Raising Errors

The most common use of errors is for checking input arguments to functions. The following example calls the error function if the function $f$ is called without any input arguments.

```
function f (arg1)
    if (nargin == 0)
        error ("not enough input arguments");
    endif
endfunction
```

When the error function is called, it prints the given message and returns to the Octave prompt. This means that no code following a call to error will be executed.

It is also possible to assign an identification string to an error. If an error has such an ID the user can catch this error as will be described in the next section. To assign an ID to an error, simply call error with two string arguments, where the first is the identification string, and the second is the actual error. Note that error IDs are in the format "NAMESPACE:ERROR-NAME". The namespace "Octave" is used for Octave's own errors. Any other string is available as a namespace for user's own errors.

```
error (template,...)
error (id, template,...)
```

Display an error message and stop m-file execution.
Format the optional arguments under the control of the template string template using the same rules as the printf family of functions (see Section 14.2.4 [Formatted Output], page 298) and print the resulting message on the stderr stream. The message is prefixed by the character string 'error: '.
Calling error also sets Octave's internal error state such that control will return to the top level without evaluating any further commands. This is useful for aborting from functions or scripts.

If the error message does not end with a newline character, Octave will print a traceback of all the function calls leading to the error. For example, given the following function definitions:

```
function f () g (); end
function g () h (); end
function h () nargin == 1 || error ("nargin != 1"); end
```

calling the function $f$ will result in a list of messages that can help you to quickly find the exact location of the error:

```
f ()
error: nargin != 1
error: called from:
error: h at line 1, column 27
error: g at line 1, column 15
error: f at line 1, column 15
```

If the error message ends in a newline character, Octave will print the message but will not display any traceback messages as it returns control to the top level. For example, modifying the error message in the previous example to end in a newline causes Octave to only print a single message:

```
function h () nargin == 1 || error ("nargin != 1\n"); end
f ()
error: nargin != 1
```

A null string (" ") input to error will be ignored and the code will continue running as if the statement were a NOP. This is for compatibility with matlab. It also makes it possible to write code such as

```
err_msg = "";
if (CONDITION 1)
    err_msg = "CONDITION 1 found";
elseif (CONDITION2)
    err_msg = "CONDITION 2 found";
endif
error (err_msg);
```

which will only stop execution if an error has been found.
Implementation Note: For compatibility with matlab, escape sequences in template (e.g., "\n" => newline) are processed regardless of whether template has been defined with single quotes, as long as there are two or more input arguments. To disable escape sequence expansion use a second backslash before the sequence (e.g., " $\backslash \backslash \mathrm{n} "$ ) or use the regexptranslate function.
See also: [warning], page 246, [lasterror], page 242.
Since it is common to use errors when there is something wrong with the input to a function, Octave supports functions to simplify such code. When the print_usage function is called, it reads the help text of the function calling print_usage, and presents a useful error. If the help text is written in Texinfo it is possible to present an error message that only contains the function prototypes as described by the @deftypefn parts of the help
text. When the help text isn't written in Texinfo, the error message contains the entire help message.

Consider the following function.

```
## -*- texinfo -*-
## @deftypefn {} f (@var{arg1})
## Function help text goes here...
## @end deftypefn
function f (arg1)
        if (nargin == 0)
            print_usage ();
        endif
endfunction
```

When it is called with no input arguments it produces the following error.
f ()
$\dashv$ error: Invalid call to f. Correct usage is:
$-1$
$\dashv \quad--\mathrm{f}$ (ARG1)
$\dashv$
$-1$
$\dashv$ Additional help for built-in functions and operators is
$\dashv$ available in the online version of the manual. Use the command
$\dashv$ 'doc <topic>' to search the manual index.
$\dashv$
$\dashv \quad$ Help and information about Octave is also available on the WWW
$\dashv$ at https://www.octave.org and via the help@octave.org
$\dashv$ mailing list.
print_usage ()
print_usage (name)
Print the usage message for the function name.
When called with no input arguments the print_usage function displays the usage message of the currently executing function.
See also: [help], page 21.
beep ()
Produce a beep from the speaker (or visual bell).
This function sends the alarm character " $\backslash \mathrm{a}$ " to the terminal. Depending on the user's configuration this may produce an audible beep, a visual bell, or nothing at all.
See also: [puts], page 297, [fputs], page 297, [printf], page 298, [fprintf], page 299.
val = beep_on_error ()
old_val = beep_on_error (new_val)
beep_on_error (new_val, "local")
Query or set the internal variable that controls whether Octave will try to ring the terminal bell before printing an error message.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

### 12.1.2 Catching Errors

When an error occurs, it can be detected and handled using the try statement as described in Section 10.9 [The try Statement], page 188. As an example, the following piece of code counts the number of errors that occurs during a for loop.

```
number_of_errors = 0;
for n = 1:100
    try
        ...
    catch
        number_of_errors++;
    end_try_catch
endfor
```

The above example treats all errors the same. In many situations it can however be necessary to discriminate between errors, and take different actions depending on the error. The lasterror function returns a structure containing information about the last error that occurred. As an example, the code above could be changed to count the number of errors related to the ' $*$ ' operator.

```
number_of_errors = 0;
for n = 1:100
    try
        ...
    catch
        msg = lasterror.message;
        if (strfind (msg, "operator *"))
            number_of_errors++;
        endif
    end_try_catch
endfor
```

Alternatively, the output of the lasterror function can be found in a variable indicated immediately after the catch keyword, as in the example below showing how to redirect an error as a warning:
try
...
catch err
warning(err.identifier, err.message);
...
end_try_catch
lasterr = lasterror ()
lasterror (err)
lasterror ("reset")
Query or set the last error message structure.

When called without arguments, return a structure containing the last error message and other information related to this error. The elements of the structure are:
message The text of the last error message
identifier
The message identifier of this error message
stack A structure containing information on where the message occurred. This may be an empty structure if the information cannot be obtained. The fields of the structure are:
file The name of the file where the error occurred
name The name of function in which the error occurred
line The line number at which the error occurred
column An optional field with the column number at which the error occurred

The last error structure may be set by passing a scalar structure, err, as input. Any fields of err that match those above are set while any unspecified fields are initialized with default values.
If lasterror is called with the argument "reset", all fields are set to their default values.

See also: [lasterr], page 243, [error], page 239, [lastwarn], page 248.

```
[msg, msgid] = lasterr ()
lasterr (msg)
lasterr (msg, msgid)
```

Query or set the last error message.
When called without input arguments, return the last error message and message identifier.

With one argument, set the last error message to msg.
With two arguments, also set the last message identifier.
See also: [lasterror], page 242, [error], page 239, [lastwarn], page 248.
The next example counts indexing errors. The errors are caught using the field identifier of the structure returned by the function lasterror.

```
number_of_errors = 0;
for n = 1:100
    try
        ...
    catch
        id = lasterror.identifier;
        if (strcmp (id, "Octave:invalid-indexing"))
            number_of_errors++;
        endif
    end_try_catch
endfor
```

The functions distributed with Octave can issue one of the following errors.
Octave:bad-alloc
Indicates that memory couldn't be allocated.
Octave:invalid-context
Indicates the error was generated by an operation that cannot be executed in the scope from which it was called. For example, the function print_usage () when called from the Octave prompt raises this error.

Octave:invalid-fun-call
Indicates that a function was called in an incorrect way, e.g., wrong number of input arguments.

Octave:invalid-indexing
Indicates that a data-type was indexed incorrectly, e.g., real-value index for arrays, nonexistent field of a structure.

Octave:invalid-input-arg
Indicates that a function was called with invalid input arguments.
Octave:undefined-function
Indicates a call to a function that is not defined. The function may exist but Octave is unable to find it in the search path.

When an error has been handled it is possible to raise it again. This can be useful when an error needs to be detected, but the program should still abort. This is possible using the rethrow function. The previous example can now be changed to count the number of errors related to the ' $*$ ' operator, but still abort if another kind of error occurs.

```
    number_of_errors = 0;
    for n = 1:100
    try
        ...
        catch
            msg = lasterror.message;
            if (strfind (msg, "operator *"))
                number_of_errors++;
            else
            rethrow (lasterror);
        endif
    end_try_catch
endfor
rethrow (err)
```

Reissue a previous error as defined by err.
err is a structure that must contain at least the "message" and "identifier" fields. err can also contain a field "stack" that gives information on the assumed location of the error. Typically err is returned from lasterror.

See also: [lasterror], page 242, [lasterr], page 243, [error], page 239.
err $=$ errno ()
err $=$ errno (val)
err $=$ errno (name)
Query or set the system-dependent variable errno.
When called with no inputs, return the current value of errno.
When called with a numeric input val, set the current value of errno to the specified value. The previous value of errno is returned as err.
When called with a character string name, return the numeric value of errno which corresponds to the specified error code. If name is not a recognized error code then -1 is returned.
See also: [errno_list], page 245.
errno_list ()
Return a structure containing the system-dependent errno values.
See also: [errno], page 245.

### 12.1.3 Recovering From Errors

Octave provides several ways of recovering from errors. There are try/catch blocks, unwind_protect/unwind_protect_cleanup blocks, and finally the onCleanup command.

The onCleanup command associates an ordinary Octave variable (the trigger) with an arbitrary function (the action). Whenever the Octave variable ceases to exist-whether due to a function return, an error, or simply because the variable has been removed with clear - then the assigned function is executed.

The function can do anything necessary for cleanup such as closing open file handles, printing an error message, or restoring global variables to their initial values. The last example is a very convenient idiom for Octave code. For example:

```
    function rand42
    old_state = rand ("state");
    restore_state = onCleanup (@() rand ("state", old_state));
    rand ("state", 42);
    endfunction # rand generator state restored by onCleanup
obj = onCleanup (function)
    Create a special object that executes a given function upon destruction.
If the object is copied to multiple variables (or cell or struct array elements) or returned from a function, function will be executed after clearing the last copy of the object. Note that if multiple local onCleanup variables are created, the order in which they are called is unspecified. For similar functionality See Section 10.8 [The unwind_protect Statement], page 188.
```


### 12.2 Handling Warnings

Like an error, a warning is issued when something unexpected happens. Unlike an error, a warning doesn't abort the currently running program. A simple example of a warning is
when a number is divided by zero. In this case Octave will issue a warning and assign the value $\operatorname{Inf}$ to the result.

```
a = 1/0
     warning: division by zero
    a a = Inf
```


### 12.2.1 Issuing Warnings

It is possible to issue warnings from any code using the warning function. In its most simple form, the warning function takes a string describing the warning as its input argument. As an example, the following code controls if the variable ' $a$ ' is non-negative, and if not issues a warning and sets ' $a$ ' to zero.

```
a = -1;
if (a < 0)
    warning ("'a' must be non-negative. Setting 'a' to zero.");
    a = 0;
endif
            \dashv'a' must be non-negative. Setting 'a' to zero.
```

Since warnings aren't fatal to a running program, it is not possible to catch a warning using the try statement or something similar. It is however possible to access the last warning as a string using the lastwarn function.

It is also possible to assign an identification string to a warning. If a warning has such an ID the user can enable and disable this warning as will be described in the next section. To assign an ID to a warning, simply call warning with two string arguments, where the first is the identification string, and the second is the actual warning. Note that warning IDs are in the format "NAMESPACE:WARNING-NAME". The namespace "Octave" is used for Octave's own warnings. Any other string is available as a namespace for user's own warnings.

```
warning (template,...)
warning (id, template, ...)
warning ("on", id)
warning ("off",id)
warning ("error", id)
warning ("query",id)
warning (state, id, "local")
warning (warning_struct)
warning_struct = warning (...)
warning (state, mode)
```

Display a warning message or control the behavior of Octave's warning system.
The first call form uses a template template and optional additional arguments to display a message on the stderr stream. The message is formatted using the same rules as the printf family of functions (see Section 14.2.4 [Formatted Output], page 298) and prefixed by the character string 'warning: '. You should use this function when you want to notify the user of an unusual condition, but only when it makes sense for your program to go on. For example:

```
warning ("foo: maybe something wrong here");
```

If the warning message does not end with a newline character, Octave will print a traceback of all the function calls leading to the warning. If the warning message does end in a newline character, Octave will suppress the traceback messages as it returns control to the top level. For more details and examples, see [error], page 239.
The optional warning identifier id allows users to enable or disable warnings tagged by this identifier. A message identifier is a string of the form "NAMESPACE:WARNING-NAME". Octave's own warnings use the "Octave" namespace (see [warning_ids], page 248). For example:

```
warning ("MyNameSpace:check-something",
    "foo: maybe something wrong here");
```

The second call form is meant to change and/or query the state of warnings. The first input argument must be a string state ("on", "off", "error", or "query") followed by an optional warning identifier id or "all" (default).
The optional output argument warning_struct is a structure or structure array with fields "state" and "identifier". The state argument may have the following values:
"on"|"off":
Enable or disable the display of warnings identified by id and optionally return their previous state stout.
"error": Turn warnings identified by id into errors and optionally return their previous state stout.
"query": Return the current state of warnings identified by id.
A structure or structure array warning_struct, with fields "state" and "identifier", may be given as an input to achieve equivalent results. The following example shows how to temporarily disable a warning and then restore its original state:

```
loglog (-1:10);
## Disable the previous warning and save its original state
[~, id] = lastwarn ();
warnstate = warning ("off", id);
loglog (-1:10);
## Restore its original state
warning (warnstate);
```

If a final argument "local" is provided then the warning state will be set temporarily until the end of the current function. Changes to warning states that are set locally affect the current function and all functions called from the current scope. The previous warning state is restored on return from the current function. The "local" option is ignored if used in the top-level workspace.
With no input argument warning () is equivalent to warning ("query", "all") except that in the absence of an output argument, the state of warnings is displayed on stderr.
The level of verbosity of the warning system may also be controlled by two modes mode:
"backtrace":
enable/disable the display of the stack trace after the warning message
"verbose":
enable/disable the display of additional information after the warning message

In this case the state argument may only be "on" or "off".
Implementation Note: For compatibility with matlab, escape sequences in template (e.g., " $\backslash \mathrm{n} "=>$ newline) are processed regardless of whether template has been defined with single quotes, as long as there are two or more input arguments. To disable escape sequence expansion use a second backslash before the sequence (e.g., "<br>n") or use the regexptranslate function.

See also: [warning_ids], page 248, [lastwarn], page 248, [error], page 239.
[msg, msgid] = lastwarn ()
lastwarn (msg)
lastwarn (msg, msgid)
Query or set the last warning message.
When called without input arguments, return the last warning message and message identifier.

With one argument, set the last warning message to msg.
With two arguments, also set the last message identifier.
See also: [warning], page 246, [lasterror], page 242, [lasterr], page 243.
The functions distributed with Octave can issue one of the following warnings.
Octave: abbreviated-property-match
If the Octave:abbreviated-property-match warning is enabled, a warning is printed if a non-exact or ambiguous match is being used for a operation specifying an object property. For example, for a graphics object, fig, with the property 'displayname', get (fig, 'dis') elicits a warning if the Octave:abbreviated-property-match warning is enabled. By default, the Octave:abbreviated-property-match warning is enabled.

Octave: addpath-pkg
If the Octave: addpath-pkg warning is enabled, Octave will warn when a package directory (i.e., +package_name) is added to the path. Typically, only the parent directory which contains the package directory should be added to the load path. By default, the Octave: addpath-pkg warning is enabled.

Octave:array-as-logical
If the Octave: array-as-logical warning is enabled, Octave will warn when an array of size greater than $1 \times 1$ is used as a truth value in an if, while, or until statement. By default, the Octave: array-as-logical warning is disabled.

Octave: array-to-scalar
If the Octave : array-to-scalar warning is enabled, Octave will warn when an implicit conversion from an array to a scalar value is attempted. By default, the Octave: array-to-scalar warning is disabled.

Octave: array-to-vector
If the Octave: array-to-vector warning is enabled, Octave will warn when an implicit conversion from an array to a vector value is attempted. By default, the Octave: array-to-vector warning is disabled.

Octave:assign-as-truth-value
If the Octave: assign-as-truth-value warning is enabled, a warning is issued for statements like

```
if (s = t)
```

    ...
    since such statements are not common, and it is likely that the intent was to write

```
if (s == t)
```

instead.
There are times when it is useful to write code that contains assignments within the condition of a while or if statement. For example, statements like

```
while (c = getc ())
```

are common in C programming.
It is possible to avoid all warnings about such statements by disabling the Octave:assign-as-truth-value warning, but that may also let real errors like

```
if (x = 1) # intended to test (x == 1)!
```

            ...
    slip by.
In such cases, it is possible suppress errors for specific statements by writing them with an extra set of parentheses. For example, writing the previous example as

```
while ((c = getc ()))
```

will prevent the warning from being printed for this statement, while allowing Octave to warn about other assignments used in conditional contexts.
By default, the Octave: assign-as-truth-value warning is enabled.
Octave: autoload-relative-file-name
If the Octave:autoload-relative-file-name is enabled, Octave will warn when parsing autoload() function calls with relative paths to function files. This usually happens when using autoload() calls in PKG_ADD files, when the PKG_ADD file is not in the same directory as the .oct file referred to by the autoload() command. By default, the Octave: autoload-relative-file-name warning is enabled.
Octave: charmat-truncated
If the Octave:charmat-truncated warning is enabled, a warning is printed when a character matrix with multiple rows is converted to a string. In this
case, the Octave interpreter keeps only the first row and discards the others. By default, the Octave:charmat-truncated warning is enabled.

Octave:classdef-to-struct
If the Octave:classdef-to-struct warning is enabled, a warning is issued when a classdef object is forcibly converted into a struct with struct (CLASSDEF_OBJ). Conversion removes the access restrictions from the object and makes private and protected properties visible. By default, the Octave:classdef-to-struct warning is enabled.

Octave:colon-complex-argument
If the Octave:colon-complex-argument warning is enabled, a warning is issued when one of the three arguments to the colon operator (base, increment, limit) is a complex value. For example, 1:3*i will cause a warning to be emitted. By default, the Octave:colon-complex-argument warning is enabled.

Octave:colon-nonscalar-argument
If the Octave:colon-nonscalar-argument warning is enabled, a warning is issued when one of the three arguments to the colon operator (base, increment, limit) is not a scalar. For example, $1:[3,5]$ will cause a warning to be emitted. By default, the Octave:colon-nonscalar-argument warning is enabled.

Octave:data-file-in-path
If the Octave: data-file-in-path warning is enabled, a warning is issued when Octave does not find the target of a file operation such as load or fopen directly, but is able to locate the file in Octave's search path for files. The warning could indicate that a different file target than the programmer intended is being used. By default, the Octave:data-file-in-path warning is enabled.

Octave:deprecated-function
If the Octave:deprecated-function warning is enabled, a warning is issued when Octave encounters a function that is obsolete and scheduled for removal from Octave. By default, the Octave:deprecated-function warning is enabled.

Octave: deprecated-keyword
If the Octave:deprecated-keyword warning is enabled, a warning is issued when Octave encounters a keyword that is obsolete and scheduled for removal from Octave. By default, the Octave:deprecated-keyword warning is enabled.

Octave:deprecated-option
If the Octave:deprecated-option warning is enabled, a warning is issued when an obsolete option or input to a function is used. By default, the Octave:deprecated-option warning is enabled.

Octave:deprecated-property
If the Octave:deprecated-property warning is enabled, a warning is issued when Octave encounters a graphics property that is obsolete and scheduled for removal from Octave. By default, the Octave: deprecated-property warning is enabled.

Octave:eigs:UnconvergedEigenvalues
If the Octave:eigs:UnconvergedEigenvalues warning is enabled then the eigs function will issue a warning if the number of calculated eigenvalues is less than the number of requested eigenvalues. By default, the Octave:eigs:UnconvergedEigenvalues warning is enabled.

Octave:empty-index
If the Octave: empty-index warning is enabled then Octave will emit a warning whenever indexing operators are used without an index, for example $x()$. By default, the Octave:empty-index warning is enabled.

Octave: erase: chararray
If the Octave: erase : chararray warning is enabled then the erase function will issue a warning if the input pattern is a character array rather than a string or cell array of strings. By default, the Octave:erase:chararray warning is enabled.

Octave: function-name-clash
If the Octave:function-name-clash warning is enabled, a warning is issued when Octave finds that the name of a function defined in a function file differs from the name of the file. (If the names disagree, the name declared inside the file is ignored.) By default, the Octave:function-name-clash warning is enabled.

Octave:future-time-stamp
If the Octave:future-time-stamp warning is enabled, Octave will print a warning if it finds a function file with a time stamp that is in the future. By default, the Octave:future-time-stamp warning is enabled.

Octave:glyph-render
If the Octave:glyph-render warning is enabled, Octave will print a warning if the glyph for a character couldn't be rendered with the current font. By default, the Octave:glyph-render warning is enabled.

Octave:imag-to-real
If the Octave:imag-to-real warning is enabled, a warning is printed for implicit conversions of complex numbers to real numbers. By default, the Octave:imag-to-real warning is disabled.

Octave:infinite-loop
If the Octave:infinite-loop warning is enabled, a warning is printed when an infinite loop is detected such as for $i=1$ :Inf or while (1). By default, the Octave:infinite-loop warning is enabled.

Octave: language-extension
Print warnings when using features that are unique to the Octave language and that may still be missing in matlab. By default, the Octave:language-extension warning is disabled. The --traditional or --braindead startup options for Octave may also be of use, see Section 2.1.1 [Command Line Options], page 15.

Octave:legacy-function
If the Octave:legacy-function warning is enabled, a warning is issued when Octave encounters a function that matlab has suggested should be avoided. The function may become obsolete at some point in the future and removed, in which case the warning will change to Octave:deprecated-function, and the function will continue to exist for two further versions of Octave before being removed. By default, the Octave:legacy-function warning is enabled.

## Octave:logical-conversion

If the Octave:logical-conversion warning is enabled, a warning is printed if an implicit conversion of an array from numerical to boolean occurs and any of the elements in the array are not equal to zero or one. By default, the Octave:logical-conversion warning is enabled.
Octave:lu:sparse_input
If the Octave:lu:sparse_input warning is enabled, Octave will warn when the lu function is called with a sparse input and less than four output arguments. In this case, sparsity-preserving column permutations are not performed and the result may be inaccurate. By default, the Octave:lu:sparse_input warning is enabled.

Octave:missing-glyph
If the Octave:glyph-render warning is enabled, Octave will print a warning if the current font doesn't provide a glyph for a used character. By default, the Octave:missing-glyph warning is enabled.

Octave:missing-semicolon
If the Octave:missing-semicolon warning is enabled, Octave will warn when statements in function definitions don't end in semicolons. By default the Octave:missing-semicolon warning is disabled.

Octave:mixed-string-concat
If the Octave:mixed-string-concat warning is enabled, print a warning when concatenating a mixture of double and single quoted strings. By default, the Octave:mixed-string-concat warning is disabled.
Octave:nearly-singular-matrix
Octave:singular-matrix
These warnings are emitted if a (nearly) singular matrix is inverted. By default, the Octave:nearly-singular-matrix and Octave:singular-matrix warnings are enabled.
Octave:neg-dim-as-zero
If the Octave:neg-dim-as-zero warning is enabled, print a warning for expressions like
eye (-1)
By default, the Octave:neg-dim-as-zero warning is disabled.
Octave:noninteger-range-as-index
If the Octave:noninteger-range-as-index warning is enabled, a warning is printed if an array is indexed with a range that contains non-integer values. For example,

```
a = [11 2 3 3 4 5];
b = 2.2:4.2
# 1.2 2.2 3.2
a(b)
# 2 3 4
```

elicits a warning if the Octave:noninteger-range-as-index warning is enabled. By default, the Octave:noninteger-range-as-index warning is enabled.

Octave: num-to-str
If the Octave:num-to-str warning is enabled, a warning is printed for implicit conversions of numbers to their UTF-8 encoded character equivalents when strings are constructed using a mixture of strings and numbers in matrix notation. For example,

```
[ "f", 111, 111 ]
=> "foo"
```

elicits a warning if the Octave:num-to-str warning is enabled. By default, the Octave:num-to-str warning is enabled.

Octave:possible-matlab-short-circuit-operator
If the Octave:possible-matlab-short-circuit-operator warning is enabled, Octave will warn about using the not short circuiting operators \& and | inside if or while conditions. They normally never short circuit, but they do short circuit when used in a condition. By default, the Octave:possible-matlab-short-circuit-operator warning is enabled.

Octave:recursive-path-search
If the Octave:recursive-path-search warning is enabled, Octave will issue a warning if addpath is used with double trailing slashes. By default, the Octave:recursive-path-search warning is enabled.

Octave:remove-init-dir
The path function changes the search path that Octave uses to find functions. It is possible to set the path to a value which excludes Octave's own built-in functions. If the Octave:remove-init-dir warning is enabled then Octave will warn when the path function has been used in a way that may render Octave unworkable. By default, the Octave:remove-init-dir warning is enabled.

Octave:reload-forces-clear
If several functions have been loaded from the same file, Octave must clear all the functions before any one of them can be reloaded. If the Octave:reload-forces-clear warning is enabled, Octave will warn you when this happens, and print a list of the additional functions that it is forced to clear. By default, the Octave:reload-forces-clear warning is enabled.

Octave: separator-insert
Print warning if commas or semicolons might be inserted automatically in literal matrices. By default, the Octave:separator-insert warning is disabled.

Octave:shadowed-function
If the Octave:shadowed-function warning is enabled, Octave will warn if a path is added to the search path that contains functions that shadow core functions. By default, the Octave:shadowed-function warning is enabled.

Octave:single-quote-string
Print warning if a single quote character is used to introduce a string constant. By default, the Octave:single-quote-string warning is disabled.

Octave:sqrtm:SingularMatrix
If the Octave:sqrtm:SingularMatrix warning is enabled, a warning is printed if the matrix square root function sqrtm is called with an input matrix that is singular. By default, the Octave:sqrtm:SingularMatrix warning is enabled.

Octave:str-to-num
If the Octave:str-to-num warning is enabled, a warning is printed for implicit conversions of strings to their numeric UTF-8 encoded byte sequences. For example,

```
"abc" + 0
# 97 98 99
```

elicits a warning if the Octave:str-to-num warning is enabled. By default, the Octave:str-to-num warning is disabled.

Octave:LaTeX:internal-error
If the Octave:LaTeX:internal-error warning is enabled, a warning is printed whenever the LaTeX renderer for text in plots encounters an issue. By default, the Octave:LaTeX:internal-error warning is enabled.

Octave: unimplemented-matlab-functionality
If the Octave:unimplemented-matlab-functionality warning is enabled, a warning is printed when a MATLAB code construct is used which the Octave interpreter parses as valid, but for which Octave does not yet implement the functionality. By default, the Octave: unimplemented-matlab-functionality warning is enabled.

Octave:variable-switch-label
If the Octave: variable-switch-label warning is enabled, Octave will print a warning if a switch label is not a constant or constant expression. By default, the Octave: variable-switch-label warning is disabled.

### 12.2.2 Enabling and Disabling Warnings

The warning function also allows you to control which warnings are actually printed to the screen. If the warning function is called with a string argument that is either "on" or "off" all warnings will be enabled or disabled.

It is also possible to enable and disable individual warnings through their string identifications. The following code will issue a warning

```
warning ("example:non-negative-variable",
    "'a' must be non-negative. Setting 'a' to zero.");
```

while the following won't issue a warning

```
warning ("off", "example:non-negative-variable");
warning ("example:non-negative-variable",
    "'a' must be non-negative. Setting 'a' to zero.");
```


## 13 Debugging

Octave includes a built-in debugger to aid in the development of scripts. This can be used to interrupt the execution of an Octave script at a certain point, or when certain conditions are met. Once execution has stopped, and debug mode is entered, the symbol table at the point where execution has stopped can be examined and modified to check for errors.

The normal command-line editing and history functions are available in debug mode.

### 13.1 Entering Debug Mode

There are two basic means of interrupting the execution of an Octave script. These are breakpoints (see Section 13.3 [Breakpoints], page 258), discussed in the next section, and interruption based on some condition.

Octave supports three means to stop execution based on the values set in the functions debug_on_interrupt, debug_on_warning, and debug_on_error.
val = debug_on_interrupt ()
old_val = debug_on_interrupt (new_val)
debug_on_interrupt (new_val, "local")
Query or set the internal variable that controls whether Octave will try to enter debugging mode when it receives an interrupt signal (typically generated with $C-c$ ).
If a second interrupt signal is received before reaching the debugging mode, a normal interrupt will occur.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [debug_on_error], page 257, [debug_on_warning], page 257.

```
val = debug_on_warning ()
```

old_val = debug_on_warning (new_val)
debug_on_warning (new_val, "local")
Query or set the internal variable that controls whether Octave will try to enter the debugger when a warning is encountered.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [debug_on_error], page 257, [debug_on_interrupt], page 257.
val = debug_on_error ()
old_val = debug_on_error (new_val)
debug_on_error (new_val, "local")
Query or set the internal variable that controls whether Octave will try to enter the debugger when an error is encountered.
This will also inhibit printing of the normal traceback message (you will only see the top-level error message).

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [debug_on_warning], page 257, [debug_on_interrupt], page 257.

### 13.2 Leaving Debug Mode

Use either dbcont or return to leave the debug mode and continue the normal execution of the script.
dbcont
Leave command-line debugging mode and continue code execution normally.
See also: [dbstep], page 263, [dbquit], page 258.
To quit debug mode and return directly to the prompt without executing any additional code use dbquit.
dbquit
dbquit all
Quit debugging mode immediately without further code execution. With no arguments, exit the current debugging level. With argument all, exit all debugging levels and return to the Octave prompt.
See also: [dbcont], page 258, [dbstep], page 263.
Finally, typing exit or quit at the debug prompt will result in Octave terminating normally.

### 13.3 Breakpoints

Breakpoints can be set in any m-file function by using the dbstop function.

```
dbstop func
dbstop func line
dbstop func line1 line2...
dbstop line1...
dbstop in func
dbstop in func at line
dbstop in func at line if "condition"
dbstop in class at method
dbstop if event
dbstop if event ID
dbstop (bp_struct)
rline = dbstop ...
```

Set breakpoints for the built-in debugger.
func is the name of a function on the current path. When already in debug mode the func argument can be omitted and the current function will be used. Breakpoints at subfunctions are set with the scope operator ' $>$ '. For example, If file.m has a subfunction func2, then a breakpoint in func2 can be specified by file>func2.
line is the line number at which to break. If line is not specified, it defaults to the first executable line in the file func.m. Multiple lines can be specified in a single command; when function syntax is used, the lines may also be passed as a single vector argument ([line1, line2, ...]).
condition is any Octave expression that can be evaluated in the code context that exists at the breakpoint. When the breakpoint is encountered, condition will be evaluated, and execution will stop if condition is true. If condition cannot be evaluated, for example because it refers to an undefined variable, an error will be thrown. Expressions with side effects (such as $\mathrm{y}++>1$ ) will alter variables, and should generally be avoided. Conditions containing quotes ('" ', ''’) or comment characters ('\#', '\%') must be enclosed in quotes. (This does not apply to conditions entered from the editor's context menu.) For example:

```
dbstop in axis at 246 if 'any (opt == "x")'
```

The form specifying event does not cause a specific breakpoint at a given function and line number. Instead it causes debug mode to be entered when certain unexpected events are encountered. Possible values are
error Stop when an error is reported. This is equivalent to specifying both debug_on_error (true) and debug_on_interrupt (true).
caught error
Stop when an error is caught by a try-catch block (not yet implemented).

```
interrupt
```

Stop when an interrupt (Ctrl-C) occurs.
naninf $\quad$ Stop when code returns a non-finite value (not yet implemented).
warning Stop when a warning is reported. This is equivalent to specifying debug_ on_warning (true).

The events error, caught error, and warning can all be followed by a string specifying an error ID or warning ID. If that is done, only errors with the specified ID will cause execution to stop. To stop on one of a set of IDs, multiple dbstop commands must be issued.

Breakpoints and events can be removed using the dbclear command with the same syntax.
It is possible to save all breakpoints and restore them at once by issuing the commands bp_state = dbstatus; ...; dbstop (bp_state).
The optional output rline is the real line number where the breakpoint was set. This can differ from the specified line if the line is not executable. For example, if a breakpoint attempted on a blank line then Octave will set the real breakpoint at the next executable line.
When a file is re-parsed, such as when it is modified outside the GUI, all breakpoints within the file are cleared.
See also: [dbclear], page 260, [dbstatus], page 260, [dbstep], page 263, [debug_on_error], page 257, [debug_on_warning], page 257, [debug_on_interrupt], page 257.

Breakpoints in class methods are also supported (e.g., dbstop ("@class/method")). However, breakpoints cannot be set in built-in functions (e.g., sin, etc.) or dynamically loaded functions (i.e., oct-files).

To set a breakpoint immediately upon entering a function use line number 1 , or omit the line number entirely and just give the function name. When setting the breakpoint Octave will ignore the leading comment block, and the breakpoint will be set on the first executable statement in the function. For example:

```
dbstop ("asind", 1)
# 29
```

Note that the return value of 29 means that the breakpoint was effectively set to line 29 . The status of breakpoints in a function can be queried with dbstatus.
dbstatus
dbstatus func
bp_list $=$ dbstatus.. .
Report the location of active breakpoints.
When called with no input or output arguments, print the list of all functions with breakpoints and the line numbers where those breakpoints are set.
If a function name func is specified then only report breakpoints for the named function and its subfunctions.
The optional return argument bp_list is a struct array with the following fields.
name The name of the function with a breakpoint. A subfunction, say func2 within an m-file, say file.m, is specified as file>func2.
file $\quad$ The name of the m -file where the function code is located.
line The line number with the breakpoint.
cond The condition that must be satisfied for the breakpoint to be active, or the empty string for unconditional breakpoints.

If dbstop if error is true but no explicit IDs are specified, the return value will have an empty field called "errs". If IDs are specified, the errs field will have one row per ID. If dbstop if error is false, there is no "errs" field. The "warn" field is set similarly by dbstop if warning.
See also: [dbstop], page 258, [dbclear], page 260, [dbwhere], page 262, [dblist], page 262, [dbstack], page 263.

Reusing the previous example, dbstatus ("asind") will return 29. The breakpoints listed can then be cleared with the dbclear function.

```
dbclear func
dbclear func line
dbclear func line1 line2 ...
dbclear line...
dbclear all
dbclear in func
dbclear in func at line
```

```
dbclear if event
dbclear ("func")
dbclear ("func", line)
dbclear ("func", line1, line2,...)
dbclear ("func", line1,...)
dbclear (line,...)
dbclear ("all")
```

Delete a breakpoint at line number line in the function func.
Arguments are
func Function name as a string variable. When already in debug mode this argument can be omitted and the current function will be used.
line Line number from which to remove a breakpoint. Multiple lines may be given as separate arguments or as a vector.
event An event such as error, interrupt, or warning (see [dbstop], page 258, for details).

When called without a line number specification all breakpoints in the named function are cleared.
If the requested line is not a breakpoint no action is performed.
The special keyword "all" will clear all breakpoints from all files.
See also: [dbstop], page 258, [dbstatus], page 260, [dbwhere], page 262.
A breakpoint may also be set in a subfunction. For example, if a file contains the functions

```
function y = func1 (x)
    y = func2 (x);
endfunction
function y = func2 (x)
    y = x + 1;
endfunction
```

then a breakpoint can be set at the start of the subfunction directly with

```
dbstop func1>func2
# 5
```

Note that ' $>$ ' is the character that distinguishes subfunctions from the m-file containing them.

Another simple way of setting a breakpoint in an Octave script is the use of the keyboard function.

```
keyboard ()
keyboard ("prompt")
```

Stop m-file execution and enter debug mode.
When the keyboard function is executed, Octave prints a prompt and waits for user input. The input strings are then evaluated and the results are printed. This makes it possible to examine the values of variables within a function, and to assign new values
if necessary. To leave the prompt and return to normal execution type 'return' or 'dbcont'. The keyboard function does not return an exit status.
If keyboard is invoked without arguments, a default prompt of 'debug> ' is used.
See also: [dbstop], page 258, [dbcont], page 258, [dbquit], page 258.
The keyboard function is placed in a script at the point where the user desires that the execution be stopped. It automatically sets the running script into the debug mode.

### 13.4 Debug Mode

There are three additional support functions that allow the user to find out where in the execution of a script Octave entered the debug mode, and to print the code in the script surrounding the point where Octave entered debug mode.

## dbwhere

In debugging mode, report the current file and line number where execution is stopped.
See also: [dbstack], page 263, [dblist], page 262, [dbstatus], page 260, [dbcont], page 258, [dbstep], page 263, [dbup], page 264, [dbdown], page 264.

```
dbtype
dbtype lineno
dbtype startl:endl
dbtype startl:end
dbtype func
dbtype func lineno
dbtype func startl:endl
dbtype func startl:end
```

Display a script file with line numbers.
When called with no arguments in debugging mode, display the script file currently being debugged.
An optional range specification can be used to list only a portion of the file. The special keyword "end" is a valid line number specification for the last line of the file.
When called with the name of a function, list that script file with line numbers.
See also: [dblist], page 262, [dbwhere], page 262, [dbstatus], page 260, [dbstop], page 258.
dblist
dblist $n$
In debugging mode, list $n$ lines of the function being debugged centered around the current line to be executed.
If unspecified $n$ defaults to 10 ( $+/-5$ lines)
See also: [dbwhere], page 262, [dbtype], page 262, [dbstack], page 263.
You may also use isdebugmode to determine whether the debugger is currently active.

## isdebugmode ()

Return true if in debugging mode, otherwise false.
See also: [dbwhere], page 262, [dbstack], page 263, [dbstatus], page 260.
Debug mode also allows single line stepping through a function using the command dbstep.

```
dbstep
```

dbstep $n$
dbstep in
dbstep out
dbnext ...

In debugging mode, execute the next $n$ lines of code.
If $n$ is omitted, execute the next single line of code. If the next line of code is itself defined in terms of an m-file remain in the existing function.
Using dbstep in will cause execution of the next line to step into any m-files defined on the next line.
Using dbstep out will cause execution to continue until the current function returns.
dbnext is an alias for dbstep.
See also: [dbcont], page 258, [dbquit], page 258.
When in debug mode the RETURN key will execute the last entered command. This is useful, for example, after hitting a breakpoint and entering dbstep once. After that, one can advance line by line through the code with only a single key stroke. This feature may be disabled using the auto_repeat_debug_command function.

```
val = auto_repeat_debug_command ()
old_val = auto_repeat_debug_command (new_val)
auto_repeat_debug_command (new_val, "local")
```

Query or set the internal variable that controls whether debugging commands are automatically repeated when the input line is empty (typing just RET).
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

### 13.5 Call Stack

The function being debugged may be the leaf node of a series of function calls. After examining values in the current subroutine it may turn out that the problem occurred in earlier pieces of code. Use dbup and dbdown to move up and down through the series of function calls to locate where variables first took on the wrong values. dbstack shows the entire series of function calls and at what level debugging is currently taking place.

```
dbstack
dbstack n
dbstack -completenames
```

[stack, idx] = dbstack (...)
Display or return current debugging function stack information.
With optional argument $n$, omit the $n$ innermost stack frames.
Although accepted, the argument -completenames is silently ignored. Octave always returns absolute filenames.

The arguments $n$ and -completenames can both be specified and may appear in any order.

The optional return argument stack is a struct array with the following fields:
file $\quad$ The name of the m -file where the function code is located.
name The name of the function with a breakpoint.
line $\quad$ The line number of an active breakpoint.
column The column number of the line where the breakpoint begins.
The return argument idx specifies which element of the stack struct array is currently active.

See also: [dbup], page 264, [dbdown], page 264, [dbwhere], page 262, [dblist], page 262, [dbstatus], page 260.

## dbup

dbup $n$
In debugging mode, move up the execution stack $n$ frames.
If $n$ is omitted, move up one frame.
See also: [dbstack], page 263, [dbdown], page 264.

## dbdown

dbdown $n$
In debugging mode, move down the execution stack $n$ frames.
If $n$ is omitted, move down one frame.
See also: [dbstack], page 263, [dbup], page 264.

### 13.6 Profiling

Octave supports profiling of code execution on a per-function level. If profiling is enabled, each call to a function (supporting built-ins, operators, functions in oct- and mex-files, userdefined functions in Octave code and anonymous functions) is recorded while running Octave code. After that, this data can aid in analyzing the code behavior, and is in particular helpful for finding "hot spots" in the code which use up a lot of computation time and are the best targets to spend optimization efforts on.

The main command for profiling is profile, which can be used to start or stop the profiler and also to query collected data afterwards. The data is returned in an Octave data structure which can then be examined or further processed by other routines or tools.

```
profile on
profile off
profile resume
profile clear
S = profile ("status")
T = profile ("info")
    Control the built-in profiler.
```

    profile on
    Start the profiler, clearing all previously collected data if there is any.
profile off
Stop profiling. The collected data can later be retrieved and examined
with $T=$ profile ("info").
profile clear
Clear all collected profiler data.
profile resume
Restart profiling without clearing the old data. All newly collected statis-
tics are added to the existing ones.
$S=$ profile ("status")
Return a structure with information about the current status of the pro-
filer. At the moment, the only field is ProfilerStatus which is either
"on" or "off".
T = profile ("info")
Return the collected profiling statistics in the structure $T$. The flat profile
is returned in the field FunctionTable which is an array of structures,
each entry corresponding to a function which was called and for which pro-
filing statistics are present. In addition, the field Hierarchical contains
the hierarchical call tree. Each node has an index into the FunctionTable
identifying the function it corresponds to as well as data fields for number
of calls and time spent at this level in the call tree.
See also: [profshow], page 265, [profexplore], page 266.

An easy way to get an overview over the collected data is profshow. This function takes the profiler data returned by profile as input and prints a flat profile, for instance:

| Function Attr |  | Time (s) | Calls |
| :---: | :---: | :---: | :---: |
| >myfib | R | 2.195 | 13529 |
| binary <= |  | 0.061 | 13529 |
| binary - |  | 0.050 | 13528 |
| binary + |  | 0.026 | 6764 |

This shows that most of the run time was spent executing the function 'myfib', and some minor proportion evaluating the listed binary operators. Furthermore, it is shown how often the function was called and the profiler also records that it is recursive.

```
profshow (data)
```

profshow (data, n)
profshow ()
profshow (n)

Display flat per-function profiler results.
Print out profiler data (execution time, number of calls) for the most critical $n$ functions. The results are sorted in descending order by the total time spent in each function. If $n$ is unspecified it defaults to 20 .
The input data is the structure returned by profile ("info"). If unspecified, profshow will use the current profile dataset.
The attribute column displays ' $R$ ' for recursive functions, and is blank for all other function types.

See also: [profexplore], page 266, [profile], page 264.
profexport (dir)
profexport (dir, data)
profexport (dir, name)
profexport (dir, name, data)
Export profiler data as HTML.
Export the profiling data in data into a series of HTML files in the folder dir. The initial file will be data/index.html.
If name is specified, it must be a string that contains a "name" for the profile being exported. This name is included in the HTML.
The input data is the structure returned by profile ("info"). If unspecified, profexport will use the current profile dataset.
See also: [profshow], page 265, [profexplore], page 266, [profile], page 264.
profexplore ()
profexplore (data)
Interactively explore hierarchical profiler output.
Assuming data is the structure with profile data returned by profile ("info"), this command opens an interactive prompt that can be used to explore the call-tree. Type help to get a list of possible commands. If data is omitted, profile ("info") is called and used in its place.
See also: [profile], page 264, [profshow], page 265.

### 13.7 Profiler Example

Below, we will give a short example of a profiler session. See Section 13.6 [Profiling], page 264, for the documentation of the profiler functions in detail. Consider the code:
global N A;
$N=300 ;$
$\mathrm{A}=\operatorname{rand}(\mathrm{N}, \mathrm{N})$;

```
function xt = timesteps (steps, x0, expM)
    global N;
    if (steps == 0)
            xt = NA (N, 0);
        else
            xt = NA (N, steps);
            x1 = expM * x0;
            xt(:, 1) = x1;
            xt(:, 2 : end) = timesteps (steps - 1, x1, expM);
    endif
endfunction
function foo ()
    global N A;
        initial = @(x) sin (x);
        x0 = (initial (linspace (0, 2 * pi, N)))';
        expA = expm (A);
        xt = timesteps (100, x0, expA);
endfunction
function fib = bar (N)
        if (N <= 2)
            fib = 1;
        else
                fib = bar (N - 1) + bar (N - 2);
        endif
endfunction
```

If we execute the two main functions, we get:

```
tic; foo; toc;
# Elapsed time is 2.37338 seconds.
tic; bar (20); toc;
# Elapsed time is 2.04952 seconds.
```

But this does not give much information about where this time is spent; for instance, whether the single call to expm is more expensive or the recursive time-stepping itself. To get a more detailed picture, we can use the profiler.

```
profile on;
foo;
profile off;
data = profile ("info");
profshow (data, 10);
```

This prints a table like:

| \# | Function Attr | Time (s) | Calls |
| :---: | :---: | :---: | :---: |
| 7 | expm | 1.034 | 1 |
| 3 | binary * | 0.823 | 117 |
| 41 | binary \} | 0.188 | 1 |
| 38 | binary | 0.126 | 2 |
| 43 | timesteps R | 0.111 | 101 |
| 44 | NA | 0.029 | 101 |
| 39 | binary + | 0.024 | 8 |
| 34 | norm | 0.011 | 1 |
| 40 | binary - | 0.004 | 101 |
| 33 | balance | 0.003 | 1 |

The entries are the individual functions which have been executed (only the 10 most important ones), together with some information for each of them. The entries like 'binary *' denote operators, while other entries are ordinary functions. They include both builtins like expm and our own routines (for instance timesteps). From this profile, we can immediately deduce that expm uses up the largest proportion of the processing time, even though it is only called once. The second expensive operation is the matrix-vector product in the routine timesteps. ${ }^{1}$

Timing, however, is not the only information available from the profile. The attribute column shows us that timesteps calls itself recursively. This may not be that remarkable in this example (since it's clear anyway), but could be helpful in a more complex setting. As to the question of why is there a 'binary $\backslash$ ' in the output, we can easily shed some light on that too. Note that data is a structure array (Section 6.1.2 [Structure Arrays], page 115) which contains the field FunctionTable. This stores the raw data for the profile shown. The number in the first column of the table gives the index under which the shown function can be found there. Looking up data.FunctionTable(41) gives:

```
scalar structure containing the fields:
    FunctionName = binary \
    TotalTime = 0.18765
    NumCalls = 1
    IsRecursive = 0
    Parents = 7
    Children = [](1x0)
```

Here we see the information from the table again, but have additional fields Parents and Children. Those are both arrays, which contain the indices of functions which have directly called the function in question (which is entry 7 , expm, in this case) or been called by it (no functions). Hence, the backslash operator has been used internally by expm.

Now let's take a look at bar. For this, we start a fresh profiling session (profile on does this; the old data is removed before the profiler is restarted):

[^3]```
profile on;
bar (20);
profile off;
profshow (profile ("info"));
This gives:
\begin{tabular}{|c|c|c|c|c|}
\hline \# & \multicolumn{2}{|l|}{Function Attr} & Time (s) & Calls \\
\hline 1 & bar & R & 2.091 & 13529 \\
\hline 2 & binary <= & & 0.062 & 13529 \\
\hline 3 & binary - & & 0.042 & 13528 \\
\hline 4 & binary + & & 0.023 & 6764 \\
\hline 5 & profile & & 0.000 & 1 \\
\hline 8 & false & & 0.000 & 1 \\
\hline 6 & nargin & & 0.000 & 1 \\
\hline 7 & binary != & & 0.000 & 1 \\
\hline 9 & __enable__ & & 0.000 & 1 \\
\hline
\end{tabular}
```

Unsurprisingly, bar is also recursive. It has been called 13,529 times in the course of recursively calculating the Fibonacci number in a suboptimal way, and most of the time was spent in bar itself.

Finally, let's say we want to profile the execution of both foo and bar together. Since we already have the run-time data collected for bar, we can restart the profiler without clearing the existing data and collect the missing statistics about foo. This is done by:

```
profile resume;
foo;
profile off;
profshow (profile ("info"), 10);
```

As you can see in the table below, now we have both profiles mixed together.

| \# | Function Attr |  | Time (s) | Calls |
| :---: | :---: | :---: | :---: | :---: |
| 1 | bar | R | 2.091 | 13529 |
| 16 | expm |  | 1.122 | 1 |
| 12 | binary * |  | 0.798 | 117 |
| 46 | binary \} |  | 0.185 | 1 |
| 45 | binary |  | 0.124 | 2 |
| 48 | timesteps | R | 0.115 | 101 |
| 2 | binary <= |  | 0.062 | 13529 |
| 3 | binary - |  | 0.045 | 13629 |
| 4 | binary + |  | 0.041 | 6772 |
| 49 | NA |  | 0.036 | 101 |

## 14 Input and Output

Octave supports several ways of reading and writing data to or from the prompt or a file. The simplest functions for data Input and Output (I/O) are easy to use, but only provide limited control of how data is processed. For more control, a set of functions modeled after the C standard library are also provided by Octave.

### 14.1 Basic Input and Output

### 14.1.1 Terminal Output

Since Octave normally prints the value of an expression as soon as it has been evaluated, the simplest of all I/O functions is a simple expression. For example, the following expression will display the value of 'pi'
pi

$$
\dashv \text { ans }=3.1416
$$

This works well as long as it is acceptable to have the name of the variable (or 'ans') printed along with the value. To print the value of a variable without printing its name, use the function disp.

The format command offers some control over the way Octave prints values with disp and through the normal echoing mechanism.

```
disp (x)
str = disp (x)
```

Display the value of $x$.
For example:

```
disp ("The value of pi is:"), disp (pi)
the value of pi is:
\dashv 3.1416
```

Note that the output from disp always ends with a newline.
If an output value is requested, disp prints nothing and returns the formatted output in a string.

See also: [fdisp], page 283.
list_in_columns (arg, width, prefix)
Return a string containing the elements of arg listed in columns with an overall maximum width of width and optional prefix prefix.
The argument arg must be a cell array of character strings or a character array.
If width is not specified or is an empty matrix, or less than or equal to zero, the width of the terminal screen is used. Newline characters are used to break the lines in the output string. For example:

```
list_in_columns ({"abc", "def", "ghijkkl", "mnop", "qrs", "tuv"}, 20)
    a abc mnop
        def qrs
        ghijkl tuv
```

whos ans
$\Rightarrow$
Variables in the current scope:

| Attr Name | Size | Bytes | Class |
| :---: | :---: | ---: | :--- |
| $========$ | $====$ | $=====$ | $=====$ |
| ans | $1 \times 37$ | 37 | char |

Total is 37 elements using 37 bytes
See also: [terminal_size], page 272.

```
terminal_size ()
```

Query or set the size of the terminal window. If called with no arguments, return a two-element row vector containing the current size of the terminal window in characters (rows and columns). If called with a two-element vector of integer values, set the terminal size and return the previous setting. Setting the size manually should not be needed when using readline for command-line editing.
See also: [list_in_columns], page 271.

```
format
format options
[format, formatspacing, uppercase] = format
```

Reset or specify the format of the output produced by disp and Octave's normal echoing mechanism.
This command only affects the display of numbers, but not how they are stored or computed. To change the internal representation from the default double use one of the conversion functions such as single, uint8, int64, etc.
By default, Octave displays 5 significant digits in a human readable form (option 'short', option 'lowercase', and option 'loose' format for matrices). If format is invoked without any options, this default format is restored.
Valid format options for floating point numbers are listed in the following table.
short Fixed point format with 5 significant figures (default).
long Fixed point format with 16 significant figures.
As with the 'short' format, Octave will switch to an exponential 'e' format if it is unable to format a matrix properly using the current format.
shorte
longe Exponential format. The number to be represented is split between a mantissa and an exponent (power of 10 ). The mantissa has 5 significant digits in the short format. In the long format, double values are displayed with 16 significant digits and single values are displayed with 8 . For example, with the 'shorte' format, pi is displayed as $3.1416 \mathrm{e}+00$. Optionally, the trailing 'e' can be split into a second argument.

```
shortg
```

longg Optimally choose between fixed point and exponential format based on the magnitude of the number. For example, with the 'shortg' format, pi . $[2 ; 4 ; 8 ; 16 ; 32]$ is displayed as ans $=$
9.8696
97.409
9488.5
$9.0032 \mathrm{e}+07$
$8.1058 \mathrm{e}+15$
Optionally, the trailing ' $g$ ' can be split into a second argument.
shorteng
longeng Identical to 'shorte' or 'longe' but displays the value using an engineering format, where the exponent is divisible by 3 . For example, with the 'shorteng' format, $10 *$ pi is displayed as $31.416 \mathrm{e}+00$. Optionally, the trailing 'eng' can be split into a second argument.
free
none Print output in free format, without trying to line up columns of matrices on the decimal point. This is a raw format equivalent to the C++ code std: : cout << variable. In general, the result is a presentation with 6 significant digits where unnecessary precision (such as trailing zeros for integers) is suppressed. Complex numbers are formatted as numeric pairs like this ' $(0.60419,0.60709)$ ' instead of like this ' $0.60419+0.60709 i$ '.

The following formats affect all numeric output (floating point and integer types).

```
"+"
"+" "chars"
plus
plus chars
```

Print a ' + ' symbol for matrix elements greater than zero, a '-' symbol for elements less than zero, and a space for zero matrix elements. This format can be useful for examining the sparsity structure of a large matrix. For very large matrices the function spy which plots the sparsity pattern will be clearer.

The optional argument chars specifies a list of 3 characters to use for printing values greater than zero, less than zero, and equal to zero. For example, with the format "+" "+-.", the matrix $[1,0,-1 ;-1,0,1]$ is displayed as

```
ans =
+.-
-.+
```

bank Print variable in a format appropriate for a currency (fixed format with two digits to the right of the decimal point). Only the real part of a variable is displayed, as the imaginary part makes no sense for a currency.
native-hex Print the hexadecimal representation of numbers as they are stored in memory. For example, on a workstation which stores 8 byte real values in IEEE format with the least significant byte first, the value of pi when printed in native-hex format is 400921fb54442d18.
hex The same as native-hex, but always print the most significant byte first.
native-bit Print the bit representation of numbers as stored in memory. For example, the value of pi is

01000000000010010010000111111011
01010100010001000010110100011000
(shown here in two 32 bit sections for typesetting purposes) when printed in native-bit format on a workstation which stores 8 byte real values in IEEE format with the least significant byte first.
bit The same as native-bit, but always print the most significant bits first.
rat Print a rational approximation, i.e., values are approximated as the ratio of small integers. For example, with the 'rat' format, pi is displayed as 355/113.

The following two options affect the display of scientific and hex notations.
lowercase (default)
Use a lowercase ' $e$ ' for the exponent character in scientific notation and lowercase ' $\mathrm{a}-\mathrm{f}$ ' for the hex digits representing 10-15.
uppercase
Use an uppercase ' $E$ ' for the exponent character in scientific notation and uppercase ' $\mathrm{A}-\mathrm{F}$ ' for the hex digits representing 10-15.

The following two options affect the display of all matrices.
compact Remove blank lines around column number labels and between matrices producing more compact output with more data per page.
loose (default)
Insert blank lines above and below column number labels and between matrices to produce a more readable output with less data per page.

If format is called with multiple competing options, the rightmost one is used. In case of an error the format remains unchanged.

If called with one to three output arguments, and no inputs, return the current format, format spacing, and uppercase preference.

See also: [fixed_point_format], page 54, [output_precision], page 53, [split_long_rows], page 54, [print_empty_dimensions], page 55, [rats], page 577.

### 14.1.1.1 Paging Screen Output

When running interactively, Octave normally sends all output directly to the Command Window. However, when using the CLI version of Octave this can create a problem because large volumes of data will stream by before you can read them. In such cases, it is better to use a paging program such as less or more which displays just one screenful at a time. With less (and some versions of more) you can also scan forward and backward, and search for specific items. The pager is enabled by the command more on.

Normally, no output is displayed by the pager until just before Octave is ready to print the top level prompt, or read from the standard input (for example, by using the fscanf or scanf functions). This means that there may be some delay before any output appears on your screen if you have asked Octave to perform a significant amount of work with a single command statement. The function fflush may be used to force output to be sent to the pager (or any other stream) immediately.

You can select the program to run as the pager with the PAGER function, and configure the pager itself with the PAGER_FLAGS function.
more
more on
more off
Turn output pagination on or off.
Without an argument, more toggles the current state.
The current state can be determined via page_screen_output.
See also: [page_screen_output], page 276, [page_output_immediately], page 276, [PAGER], page 275, [PAGER_FLAGS], page 275.
val = PAGER ()
old_val = PAGER (new_val)
PAGER (new_val, "local")
Query or set the internal variable that specifies the program to use to display terminal output on your system.
The default value is normally "less", "more", or "pg", depending on what programs are installed on your system. See Appendix E [Installation], page 1085.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [PAGER_FLAGS], page 275, [page_output_immediately], page 276, [more], page 275, [page_screen_output], page 276.
val = PAGER_FLAGS ()
old_val = PAGER_FLAGS (new_val)
PAGER_FLAGS (new_val, "local")
Query or set the internal variable that specifies the options to pass to the pager.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [PAGER], page 275, [more], page 275, [page_screen_output], page 276, [page_output_immediately], page 276.
val = page_screen_output ()
old_val = page_screen_output (new_val)
page_screen_output (new_val, "local")
Query or set the internal variable that controls whether output intended for the terminal window that is longer than one page is sent through a pager.
This allows you to view one screenful at a time. Some pagers (such as less-see Appendix E [Installation], page 1085) are also capable of moving backward on the output.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [more], page 275, [page_output_immediately], page 276, [PAGER], page 275, [PAGER_FLAGS], page 275.

```
val = page_output_immediately ()
```

old_val = page_output_immediately (new_val)
page_output_immediately (new_val, "local")
Query or set the internal variable that controls whether Octave sends output to the pager as soon as it is available.
When the value is false, Octave buffers its output and waits until just before the prompt is printed to flush it to the pager. This is the default.
When page_screen_output is false, this variable has no effect.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [page_screen_output], page 276, [more], page 275, [PAGER], page 275, [PAGER_FLAGS], page 275.
fflush (fid)
Flush output to file descriptor fid.
fflush returns 0 on success and an OS dependent error value ( -1 on Unix) on error.
Programming Note: Flushing is useful for ensuring that all pending output makes it to the screen before some other event occurs. For example, it is always a good idea to flush the standard output stream before calling input.
See also: [fopen], page 295, [fclose], page 296.

### 14.1.2 Terminal Input

Octave has three functions that make it easy to prompt users for input. The input and menu functions are normally used for managing an interactive dialog with a user, and the keyboard function is normally used for doing simple debugging.

```
ans = input (prompt)
ans = input (prompt, "s")
```

    Print prompt and wait for user input.
    For example,

```
input ("Pick a number, any number! ")
```

prints the prompt
Pick a number, any number!
and waits for the user to enter a value. The string entered by the user is evaluated as an expression, so it may be a literal constant, a variable name, or any other valid Octave code.
The number of return arguments, their size, and their class depend on the expression entered.

If you are only interested in getting a literal string value, you can call input with the character string "s" as the second argument. This tells Octave to return the string entered by the user directly, without evaluating it first.
Because there may be output waiting to be displayed by the pager, it is a good idea to always call fflush (stdout) before calling input. This will ensure that all pending output is written to the screen before your prompt.

See also: [yes_or_no], page 277, [kbhit], page 278, [pause], page 947, [menu], page 277, [listdlg], page 918.

```
choice = menu (title, opt1,...)
```

choice $=$ menu (title, $\{o p t 1, \ldots\}$ )

Display a menu with heading title and options opt1, ..., and wait for user input.
If the GUI is running, the menu is displayed graphically using listdlg. Otherwise, the title and menu options are printed on the console.
title is a string and the options may be input as individual strings or as a cell array of strings.

The return value choice is the number of the option selected by the user counting from 1. If the user aborts the dialog or makes an invalid selection then 0 is returned.

This function is useful for interactive programs. There is no limit to the number of options that may be passed in, but it may be confusing to present more than will fit easily on one screen.

See also: [input], page 276, [listdlg], page 918.

```
ans = yes_or_no ("prompt")
```

Ask the user a yes-or-no question.
Return logical true if the answer is yes or false if the answer is no.
Takes one argument, prompt, which is the string to display when asking the question. prompt should end in a space; yes-or-no adds the string ' (yes or no)' to it. The user must confirm the answer with RET and can edit it until it has been confirmed.

See also: [input], page 276.

For input, the normal command line history and editing functions are available at the prompt.

Octave also has a function that makes it possible to get a single character from the keyboard without requiring the user to type a carriage return.
kbhit ()
kbhit (1)
Read a single keystroke from the keyboard.
If called with an argument, don't wait for a keypress.
For example,

```
x = kbhit ();
```

will set $x$ to the next character typed at the keyboard as soon as it is typed.

```
x = kbhit (1);
```

is identical to the above example, but doesn't wait for a keypress, returning the empty string if no key is available.
See also: [input], page 276, [pause], page 947.

### 14.1.3 Simple File I/O

The save and load commands allow data to be written to and read from disk files in various formats. The default format of files written by the save command can be controlled using the functions save_default_options and save_precision.

As an example the following code creates a 3 -by- 3 matrix and saves it to the file 'myfile.mat'.
$\mathrm{A}=[1: 3 ; 4: 6 ; 7: 9] ;$
save myfile.mat A
Once one or more variables have been saved to a file, they can be read into memory using the load command.
load myfile.mat
A

| $\dashv A$ |  |  |  |
| :--- | :--- | :--- | :--- |
| $\dashv$ |  |  |  |
| $\dashv$ | 1 | 2 | 3 |
| $\dashv$ | 4 | 5 | 6 |
| $\dashv$ | 7 | 8 | 9 |

save file
save options file
save options file v1 v2 ...
save options file -struct STRUCT
save options file -struct STRUCT f1 f2 ...
save - v1 v2...
str = save ("-", "v1", "v2", ...)
Save the named variables $v 1, v 2, \ldots$, in the file file.
The special filename '-' may be used to return the content of the variables as a string. If no variable names are listed, Octave saves all the variables in the current scope.

Otherwise, full variable names or pattern syntax can be used to specify the variables to save. If the -struct modifier is used then the fields of the scalar struct are saved as if they were variables with the corresponding field names. The -struct option can be combined with specific field names $f 1, f 2, \ldots$ to write only certain fields to the file.
Valid options for the save command are listed in the following table. Options that modify the output format override the format specified by save_default_options.
If save is invoked using the functional form

```
save ("-option1", ..., "file", "v1", ...)
```

then the options, file, and variable name arguments ( $v 1, \ldots$ ) must be specified as character strings.
If called with a filename of "-", write the output to stdout if nargout is 0 , otherwise return the output in a character string.
-append Append to the destination instead of overwriting.
-ascii Save a matrix in a text file without a header or any other information. The matrix must be 2-D and only the real part of any complex value is written to the file. Numbers are stored in single-precision format and separated by spaces. Additional options for the -ascii format are
-double Store numbers in double-precision format.
-tabs Separate numbers with tabs.
-binary Save the data in Octave's binary data format.
-float-binary
Save the data in Octave's binary data format but using only single precision. Use this format only if you know that all the values to be saved can be represented in single precision.
-hdf5 Save the data in HDF5 format. (HDF5 is a free, portable, binary format developed by the National Center for Supercomputing Applications at the University of Illinois.) This format is only available if Octave was built with a link to the HDF5 libraries.
-float-hdf5
Save the data in HDF5 format but using only single precision. Use this format only if you know that all the values to be saved can be represented in single precision.
-text Save the data in Octave's text data format. (default)
-v7. 3
-V7. 3
-7.3 Octave does not yet implement matlab's v7.3 binary data format.
-v7
-V7
-7
-mat7-binary
Save the data in matlab's v7 binary data format.

```
-v6
-V6
-6
-mat
-mat-binary
```

Save the data in matlab's v6 binary data format.
-v4
-V4
-4
-mat4-binary
Save the data in matlab's v4 binary data format.
-zip
-z Use the gzip algorithm to compress the file. This works on files that are compressed with gzip outside of Octave, and gzip can also be used to convert the files for backward compatibility. This option is only available if Octave was built with a link to the zlib libraries.

The list of variables to save may use wildcard patterns (glob patterns) containing the following special characters:
? Match any single character.

* Match zero or more characters.
[ list] Match the list of characters specified by list. If the first character is ! or ${ }^{\wedge}$, match all characters except those specified by list. For example, the pattern [a-zA-Z] will match all lower and uppercase alphabetic characters.
Wildcards may also be used in the field name specifications when using the -struct modifier (but not in the struct name itself).

Except when using the matlab binary data file format or the '-ascii' format, saving global variables also saves the global status of the variable. If the variable is restored at a later time using 'load', it will be restored as a global variable.

## Example:

The command

```
save -binary data a b*
```

saves the variable 'a' and all variables beginning with 'b' to the file data in Octave's binary format.
See also: [load], page 281, [save_default_options], page 280, [save_header_format_string], page 281, [save_precision], page 281, [dlmread], page 284, [csvread], page 285, [fread], page 308 .

There are three functions that modify the behavior of save.

```
val = save_default_options ()
old_val = save_default_options (new_val)
```

save_default_options (new_val, "local")
Query or set the internal variable that specifies the default options for the save command, and defines the default format.
The default value is "-text" (Octave's own text-based file format). See the documentation of the save command for other choices.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [save], page 278, [save_header_format_string], page 281, [save_precision], page 281.
val = save_precision ()
old_val = save_precision (new_val)
save_precision (new_val, "local")
Query or set the internal variable that specifies the number of digits to keep when saving data in text format.
The default value is 17 which is the minimum necessary for the lossless saving and restoring of IEEE-754 double values; For IEEE-754 single values the minimum value is 9 . If file size is a concern, it is probably better to choose a binary format for saving data rather than to reduce the precision of the saved values.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [save_default_options], page 280.
val = save_header_format_string ()
old_val = save_header_format_string (new_val)
save_header_format_string (new_val, "local")
Query or set the internal variable that specifies the format string used for the comment line written at the beginning of text-format data files saved by Octave.
The format string is passed to strftime and must begin with the character '\#' and contain no newline characters. If the value of save_header_format_string is the empty string, the header comment is omitted from text-format data files. The default value is
"\# Created by Octave VERSION, \%a \%b \%d \%H:\%M:\%S \%Y \%Z [USER@HOST](mailto:USER@HOST)"
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [strftime], page 943, [save_default_options], page 280.
load file
load options file
load options file v1 v2 ...
S = load ("options", "file", "v1", "v2", ...)
load file options
load file options v1 v2...
S = load ("file", "options", "v1", "v2", ...)
Load the named variables $v 1, v 2, \ldots$, from the file file.
If no variables are specified then all variables found in the file will be loaded. As with save, the list of variables to extract can be full names or use a pattern syntax. The format of the file is automatically detected but may be overridden by supplying the appropriate option.
If load is invoked using the functional form

> load ("-option1", ..., "file", "v1", ...)
then the options, file, and variable name arguments (v1, ...) must be specified as character strings.
If a variable that is not marked as global is loaded from a file when a global symbol with the same name already exists, it is loaded in the global symbol table. Also, if a variable is marked as global in a file and a local symbol exists, the local symbol is moved to the global symbol table and given the value from the file.
If invoked with a single output argument, Octave returns data instead of inserting variables in the symbol table. If the data file contains only numbers (TAB- or space-delimited columns), a matrix of values is returned. Otherwise, load returns a structure with members corresponding to the names of the variables in the file.
The load command can read data stored in Octave's text and binary formats, and matlab's binary format. If compiled with zlib support, it can also load gzip-compressed files. It will automatically detect the type of file and do conversion from different floating point formats (currently only IEEE big and little endian, though other formats may be added in the future).
Valid options for load are listed in the following table.
-force This option is accepted for backward compatibility but is ignored. Octave now overwrites variables currently in memory with those of the same name found in the file.
-ascii Force Octave to assume the file contains columns of numbers in text format without any header or other information. Data in the file will be loaded as a single numeric matrix with the name of the variable derived from the name of the file.
-binary Force Octave to assume the file is in Octave's binary format.
-hdf5 Force Octave to assume the file is in HDF5 format. (HDF5 is a free, portable binary format developed by the National Center for Supercomputing Applications at the University of Illinois.) Note that Octave can only read HDF5 files that were created by itself with save. This format is only available if Octave was built with a link to the HDF5 libraries.
-import This option is accepted for backward compatibility but is ignored. Octave can now support multi-dimensional HDF data and automatically modifies variable names if they are invalid Octave identifiers.
-text Force Octave to assume the file is in Octave's text format.
-v7. 3
-V7. 3
-7.3 Octave does not yet implement matlab's v7.3 binary data format.
-v7
-V7
-7
-mat7-binary
Force Octave to assume the file is in MATLAB's version 7 binary format.
-v6
-V6
-6
-mat
-mat-binary
Force Octave to assume the file is in MATLAB's version 6 binary format.
$-\mathrm{v} 4$
-V4
-4
-mat4-binary
Force Octave to assume the file is in matlab's version 4 binary format.
See also: [save], page 278, [dlmwrite], page 284, [csvwrite], page 285, [fwrite], page 310.
str $=$ fileread (filename)
Read the contents of filename and return it as a string.
See also: [fread], page 308, [fscanf], page 304, [importdata], page 291, [textscan], page 287, [type], page 148.
native_float_format ()
Return the native floating point format as a string.
It is possible to write data to a file in a similar way to the disp function for writing data to the screen. The fdisp works just like disp except its first argument is a file pointer as created by fopen. As an example, the following code writes to data 'myfile.txt'.

```
fid = fopen ("myfile.txt", "w");
fdisp (fid, "3/8 is ");
fdisp (fid, 3/8);
fclose (fid);
```

See Section 14.2.1 [Opening and Closing Files], page 294, for details on how to use fopen and fclose.
fdisp (fid, x)
Display the value of $x$ on the stream fid.
For example:

```
fdisp (stdout, "The value of pi is:"), fdisp (stdout, pi)
the value of pi is:
\dashv3.1416
```

Note that the output from fdisp always ends with a newline.
See also: [disp], page 271.
Octave can also read and write matrices text files such as comma separated lists.
dlmwrite (file, M)
dlmwrite (file, M, delim, r, c)
dlmwrite (file, M, key, val...)
dlmwrite (file, M, "-append", ...)
dlmwrite (fid,...)
Write the numeric matrix $M$ to the text file file using a delimiter.
file should be a filename or a writable file ID given by fopen.
The parameter delim specifies the delimiter to use to separate values on a row. If no delimiter is specified the comma character ',' is used.
The value of $r$ specifies the number of delimiter-only lines to add to the start of the file.
The value of $c$ specifies the number of delimiters to prepend to each line of data.
If the argument "-append" is given, append to the end of file.
In addition, the following keyword value pairs may appear at the end of the argument list:
"append" Either "on" or "off". See "-append" above.
"delimiter"
See delim above.
"newline"
The character(s) to separate each row. Three special cases exist for this option. "unix" is changed into " $\backslash \mathrm{n} "$, " pc " is changed into " $\backslash \mathrm{r} \backslash \mathrm{n}$ ", and "mac" is changed into " $\backslash r$ ". Any other value is used directly as the newline separator.
"roffset"
See $r$ above.
"coffset"
See $c$ above.
"precision"
The precision to use when writing the file. It can either be a format string (as used by fprintf) or a number of significant digits.

```
dlmwrite ("file.csv", reshape (1:16, 4, 4));
dlmwrite ("file.tex", a, "delimiter", "&", "newline", "\n")
```

See also: [dlmread], page 284, [csvread], page 285, [csvwrite], page 285.

```
data = dlmread (file)
data = dlmread (file, sep)
data = dlmread (file, sep, r0, c0)
data = dlmread (file, sep, range)
```

data $=$ dlmread (..., "emptyvalue", EMPTYVAL)
Read numeric data from the text file file which uses the delimiter sep between data values.
If sep is not defined the separator between fields is determined from the file itself.
The optional scalar arguments r0 and $c 0$ define the starting row and column of the data to be read. These values are indexed from zero, i.e., the first data row corresponds to an index of zero.
The range parameter specifies exactly which data elements are read. The first form of the parameter is a 4 -element vector containing the upper left and lower right corners [ $R 0, C 0, R 1, C 1]$ where the indices are zero-based. To specify the last column-the equivalent of end when indexing-use the specifier Inf. Alternatively, a spreadsheet style form such as "A2 . .Q15" or "T1:AA5" can be used. The lowest alphabetical index ' $A$ ' refers to the first column. The lowest row index is 1 .
file should be a filename or a file id given by fopen. In the latter case, the file is read until end of file is reached.

The "emptyvalue" option may be used to specify the value used to fill empty fields. The default is zero. Note that any non-numeric values, such as text, are also replaced by the "emptyvalue".
See also: [csvread], page 285, [textscan], page 287, [dlmwrite], page 284.
csvwrite (filename, x)
csvwrite (filename, $x$, dlm_opt1, ...)
Write the numeric matrix $x$ to the file filename in comma-separated-value (CSV) format.
This function is equivalent to

```
dlmwrite (filename, x, ",", dlm_opt1, ...)
```

Any optional arguments are passed directly to dlmwrite (see [dlmwrite], page 284).
See also: [csvread], page 285, [dlmwrite], page 284, [dlmread], page 284.
$x=\operatorname{csvread}$ (filename)
$\mathrm{x}=\mathrm{csvread}$ (filename, dlm_opt1, ...)
Read the comma-separated-value (CSV) file filename into the matrix $x$.
Note: only CSV files containing numeric data can be read.
This function is equivalent to

```
x = dlmread (filename, "," , dlm_opt1, ...)
```

Any optional arguments are passed directly to dlmread (see [dlmread], page 284).
See also: [dlmread], page 284, [textscan], page 287, [csvwrite], page 285, [dlmwrite], page 284.

Formatted data from can be read from, or written to, text files as well.
[a, ...] = textread (filename)
[a, ...] = textread (filename, format)
[a, ...] = textread (filename, format, n)
[a, ...] = textread (filename, format, prop1, value1, ...)
[a, ...] = textread (filename, format, n, prop1, value1, ...)
This function is obsolete. Use textscan instead.
Read data from a text file.
The file filename is read and parsed according to format. The function behaves like strread except it works by parsing a file instead of a string. See the documentation of strread for details.

In addition to the options supported by strread, this function supports two more:

- "headerlines": The first value number of lines of filename are skipped.
- "endofline": Specify a single character or " $\backslash \mathrm{r} \backslash \mathrm{n}$ ". If no value is given, it will be inferred from the file. If set to "" (empty string) EOLs are ignored as delimiters.

The optional input $n$ (format repeat count) specifies the number of times the format string is to be used or the number of lines to be read, whichever happens first while reading. The former is equivalent to requesting that the data output vectors should be of length $N$. Note that when reading files with format strings referring to multiple lines, $n$ should rather be the number of lines to be read than the number of format string uses.

If the format string is empty (not just omitted) and the file contains only numeric data (excluding headerlines), textread will return a rectangular matrix with the number of columns matching the number of numeric fields on the first data line of the file. Empty fields are returned as zero values.

Examples:

```
Assume a data file like:
1 a 2 b
3 c 4 d
5 e
[a, b] = textread (f, "%f %s")
returns two columns of data, one with doubles, the other a
cellstr array:
a = [1; 2; 3; 4; 5]
b = {"a"; "b"; "c"; "d"; "e"}
[a, b] = textread (f, "%f %s", 3)
(read data into two culumns, try to use the format string
three times)
returns
a = [1; 2; 3]
b = {"a"; "b"; "c"}
```

```
With a data file like:
1
a
2
b
```

```
[a, b] = textread (f, "%f %s", 2)
returns a = 1 and b = {"a"}; i.e., the format string is used
only once because the format string refers to 2 lines of the
data file. To obtain 2x1 data output columns, specify N = 4
(number of data lines containing all requested data) rather
than 2.
```

See also: [textscan], page 287, [load], page 281, [dlmread], page 284, [fscanf], page 304, [strread], page 102.
$C=$ textscan (fid, format)
$C=$ textscan (fid, format, repeat)
$C=$ textscan (fid, format, param, value, ...)
$C=$ textscan (fid, format, repeat, param, value, ...)
$C=$ textscan (str, ...)
[C, position, errmsg] = textscan (...)
Read data from a text file or string.
The string str or file associated with fid is read from and parsed according to format. The function is an extension of strread and textread. Differences include: the ability to read from either a file or a string, additional options, and additional format specifiers.
The input is interpreted as a sequence of words, delimiters (such as whitespace), and literals. The characters that form delimiters and whitespace are determined by the options. The format consists of format specifiers interspersed between literals. In the format, whitespace forms a delimiter between consecutive literals, but is otherwise ignored.
The output $C$ is a cell array where the number of columns is determined by the number of format specifiers.
The first word of the input is matched to the first specifier of the format and placed in the first column of the output; the second is matched to the second specifier and placed in the second column and so forth. If there are more words than specifiers then the process is repeated until all words have been processed or the limit imposed by repeat has been met (see below).
The string format describes how the words in str should be parsed. As in fscanf, any (non-whitespace) text in the format that is not one of these specifiers is considered a literal. If there is a literal between two format specifiers then that same literal must appear in the input stream between the matching words.
The following specifiers are valid:
\%f
\%f64
\%n The word is parsed as a number and converted to double.
\%f32 The word is parsed as a number and converted to single (float).
\%d
\% d8
\%d16
\%d32
\%d64
The word is parsed as a number and converted to int8, int16, int32, or int64. If no size is specified then int 32 is used.
\%u
\%u8
\%u16
\%u32
\%u64
$\%$ s The word is parsed as a string ending at the last character before whitespace, an end-of-line, or a delimiter specified in the options.
$\% \mathrm{q} \quad$ The word is parsed as a "quoted string". If the first character of the string is a double quote (") then the string includes everything until a matching double quote - including whitespace, delimiters, and end-of-line characters. If a pair of consecutive double quotes appears in the input, it is replaced in the output by a single double quote. For examples, the input "He said ""Hello"" " would return the value 'He said "Hello"'.
$\%$ The next character of the input is read. This includes delimiters, whitespace, and end-of-line characters.
$\%[. .$.
\% [^...] In the first form, the word consists of the longest run consisting of only characters between the brackets. Ranges of characters can be specified by a hyphen; for example, $\%[0-9 \mathrm{a}-\mathrm{zA}-\mathrm{Z}]$ matches all alphanumeric characters (if the underlying character set is ASCII). Since matlab treats hyphens literally, this expansion only applies to alphanumeric characters. To include '-' in the set, it should appear first or last in the brackets; to include ']', it should be the first character. If the first character is ${ }^{\prime} \rightarrow$, then the word consists of characters not listed.
$\% \mathrm{~N} . . \quad$ For $\% \mathrm{~s}, \% \mathrm{c} \% \mathrm{~d}, \% \mathrm{f}, \% \mathrm{n}, \% \mathrm{u}$, an optional width can be specified as $\% \mathrm{Ns}$, etc. where N is an integer $>1$. For $\% \mathrm{c}$, this causes exactly N characters to be read instead of a single character. For the other specifiers, it is an upper bound on the number of characters read; normal delimiters can cause fewer characters to be read. For complex numbers, this limit applies to the real and imaginary components individually. For \%f and \%n, format specifiers like $\% \mathrm{~N} . \mathrm{Mf}$ are allowed, where M is an upper bound on number of characters after the decimal point to be considered; subsequent digits are skipped. For example, the specifier $\% 8.2 \mathrm{f}$ would read 12.345 e 6 as 1.234 e 7 .
$\% * \ldots \quad$ The word specified by the remainder of the conversion specifier is skipped.
literals In addition the format may contain literal character strings; these will be skipped during reading. If the input string does not match this literal, the processing terminates.

Parsed words corresponding to the first specifier are returned in the first output argument and likewise for the rest of the specifiers.
By default, if there is only one input argument, format is "\%f". This means that numbers are read from the input into a single column vector. If format is explicitly empty ("") then textscan will return data in a number of columns matching the number of fields on the first data line of the input. Either of these is suitable only when the input is exclusively numeric.
For example, the string

```
str = "\
Bunny Bugs 5.5\n\
Duck Daffy -7.5e-5\n\
Penguin Tux 6"
```

can be read using

```
a = textscan (str, "%s %s %f");
```

The optional numeric argument repeat can be used for limiting the number of items read:
-1 Read all of the string or file until the end (default).
N Read until the first of two conditions occurs: 1) the format has been processed N times, or 2) N lines of the input have been processed. Zero $(0)$ is an acceptable value for repeat. Currently, end-of-line characters inside $\% \mathrm{q}, \% \mathrm{c}$, and $\%[\ldots] \$$ conversions do not contribute to the line count. This is incompatible with matlab and may change in future.

The behavior of textscan can be changed via property/value pairs. The following properties are recognized:
"BufSize"
This specifies the number of bytes to use for the internal buffer. A modest speed improvement may be obtained by setting this to a large value when reading a large file, especially if the input contains long strings. The default is 4096, or a value dependent on $n$ if that is specified.
"CollectOutput"
A value of 1 or true instructs textscan to concatenate consecutive columns of the same class in the output cell array. A value of 0 or false (default) leaves output in distinct columns.

## "CommentStyle"

Specify parts of the input which are considered comments and will be skipped. value is the comment style and can be either (1) A string or $1 \times 1$ cell string, to skip everything to the right of it; (2) A cell array of two strings, to skip everything between the first and second strings. Comments are only parsed where whitespace is accepted and do not act as delimiters.
"Delimiter"
If value is a string, any character in value will be used to split the input into words. If value is a cell array of strings, any string in the array will be used to split the input into words. (default value $=$ any whitespace.)
"EmptyValue"
Value to return for empty numeric values in non-whitespace delimited data. The default is NaN. When the data type does not support NaN (int32 for example), then the default is zero.
"EndOfLine"
value can be either an empty or one character specifying the end-of-line character, or the pair " $\backslash r \backslash n$ " (CRLF). In the latter case, any of " $\backslash r$ ", " $\backslash n$ " or " $\backslash r \backslash n$ " is counted as a (single) newline. If no value is given, " $\backslash \mathrm{r} \backslash \mathrm{n}$ " is used.
"HeaderLines"
The first value number of lines of fid are skipped. Note that this does not refer to the first non-comment lines, but the first lines of any type.
"MultipleDelimsAsOne"
If value is nonzero, treat a series of consecutive delimiters, without whitespace in between, as a single delimiter. Consecutive delimiter series need not be vertically aligned. Without this option, a single delimiter before the end of the line does not cause the line to be considered to end with an empty value, but a single delimiter at the start of a line causes the line to be considered to start with an empty value.
"TreatAsEmpty"
Treat single occurrences (surrounded by delimiters or whitespace) of the string(s) in value as missing values.
"ReturnOnError"
If set to numerical 1 or true, return normally as soon as an error is encountered, such as trying to read a string using $\%$ f. If set to 0 or false, return an error and no data.
"Whitespace"
Any character in value will be interpreted as whitespace and trimmed; The default value for whitespace is " $\backslash \mathrm{b} \backslash r \backslash n \backslash t$ " (note the space). Unless whitespace is set to "" (empty) AND at least one "\%s" format conversion specifier is supplied, a space is always part of whitespace.

When the number of words in str or fid doesn't match an exact multiple of the number of format conversion specifiers, textscan's behavior depends on whether the last character of the string or file is an end-of-line as specified by the EndOfLine option:
last character $=$ end-of-line
Data columns are padded with empty fields, NaN or 0 (for integer fields) so that all columns have equal length
last character is not end-of-line
Data columns are not padded; textscan returns columns of unequal length

The second output position provides the location, in characters from the beginning of the file or string, where processing stopped.

See also: [dlmread], page 284, [fscanf], page 304, [load], page 281, [strread], page 102, [textread], page 285.

The importdata function has the ability to work with a wide variety of data.
$A=$ importdata (fname)
$A=$ importdata (fname, delimiter)
$A=$ importdata (fname, delimiter, header_rows)
[A, delimiter] = importdata (...)
[A, delimiter, header_rows] = importdata (...) Import data from the file fname.
Input parameters:

- fname The name of the file containing data.
- delimiter The character separating columns of data. Use \t for tab. (Only valid for ASCII files)
- header_rows The number of header rows before the data begins. (Only valid for ASCII files)

Different file types are supported:

- ASCII table

Import ASCII table using the specified number of header rows and the specified delimiter.

- Image file
- matlab file
- Spreadsheet files (depending on external software)
- WAV file

See also: [textscan], page 287, [dlmread], page 284, [csvread], page 285, [load], page 281.

After importing, the data may need to be transformed before further analysis. The rescale function can shift and normalize a data set to a specified range.
$B=$ rescale $(A)$
$B=\operatorname{rescale}(A, l, u)$
$B=$ rescale (..., "inputmin", inmin)
$B=$ rescale (..., "inputmax", inmax)
Scale matrix elements to a specified range of values.
When called with a single matrix argument $A$, rescale elements to occupy the interval $[0,1]$.

The optional inputs $[1, u]$ will scale $A$ to the interval with lower bound $l$ and upper bound $u$.
The optional input "inputmin" replaces all elements less than the specified value inmin with inmin. Similarly, the optional input "inputmax" replaces all elements greater than the specified value inmax with inmax. If unspecified the minimum and maximum are taken from the data itself (inmin $=\min (A(:))$ and inmax $=\max (A(:))$.

Programming Notes: The applied formula is

$$
B=l+\frac{A-\text { inmin }}{\text { inmax }- \text { inmin }} \cdot(u-l)
$$

The class of the output matrix $B$ is single if the input $A$ is single, but otherwise is of class double for inputs which are of double, integer, or logical type.
See also: [bounds], page 758, [min], page 559, [max], page 558.

### 14.1.3.1 Saving Data on Unexpected Exits

If Octave for some reason exits unexpectedly it will by default save the variables available in the workspace to a file in the current directory. By default this file is named 'octave-workspace' and can be loaded into memory with the load command. While the default behavior most often is reasonable it can be changed through the following functions.

```
val = crash_dumps_octave_core ()
```

old_val = crash_dumps_octave_core (new_val)
crash_dumps_octave_core (new_val, "local")

Query or set the internal variable that controls whether Octave tries to save all current variables to the file octave-workspace if it crashes or receives a hangup, terminate or similar signal.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [octave_core_file_limit], page 293, [octave_core_file_name], page 294, [octave_core_file_options], page 293.
val = sighup_dumps_octave_core ()
old_val = sighup_dumps_octave_core (new_val)
sighup_dumps_octave_core (new_val, "local")
Query or set the internal variable that controls whether Octave tries to save all current variables to the file octave-workspace if it receives a hangup signal.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

```
val = sigquit_dumps_octave_core ()
old_val = sigquit_dumps_octave_core (new_val)
```

sigquit_dumps_octave_core (new_val, "local")
Query or set the internal variable that controls whether Octave tries to save all current variables to the file octave-workspace if it receives a quit signal.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

```
val = sigterm_dumps_octave_core ()
old_val = sigterm_dumps_octave_core (new_val)
sigterm_dumps_octave_core (new_val, "local")
```

Query or set the internal variable that controls whether Octave tries to save all current variables to the file octave-workspace if it receives a terminate signal.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
val = octave_core_file_options ()
old_val = octave_core_file_options (new_val)
octave_core_file_options (new_val, "local")
Query or set the internal variable that specifies the options used for saving the workspace data if Octave aborts.
The value of octave_core_file_options should follow the same format as the options for the save function. The default value is Octave's binary format.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [crash_dumps_octave_core], page 292, [octave_core_file_name], page 294, [octave_core_file_limit], page 293.

```
val = octave_core_file_limit ()
old_val = octave_core_file_limit (new_val)
octave_core_file_limit (new_val, "local")
```

Query or set the internal variable that specifies the maximum amount of memory that Octave will save when writing a crash dump file.
The limit is measured in kilobytes and is applied to the top-level workspace. The name of the crash dump file is specified by octave_core_file_name.
If octave_core_file_options flags specify a binary format, then octave_core_file_limit will be approximately the maximum size of the file. If a text file format is used, then the file could be much larger than the limit. The default value is -1 (unlimited).
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [crash_dumps_octave_core], page 292, [octave_core_file_name], page 294, [octave_core_file_options], page 293.

```
val = octave_core_file_name ()
```

old_val = octave_core_file_name (new_val)
octave_core_file_name (new_val, "local")

Query or set the internal variable that specifies the name of the file used for saving data from the top-level workspace if Octave aborts.
The default value is "octave-workspace"
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [crash_dumps_octave_core], page 292, [octave_core_file_name], page 294, [octave_core_file_options], page 293.

### 14.2 C-Style I/O Functions

Octave's C-style input and output functions provide most of the functionality of the C programming language's standard I/O library. The argument lists for some of the input functions are slightly different, however, because Octave has no way of passing arguments by reference.

In the following, file refers to a filename and fid refers to an integer file number, as returned by fopen.

There are three files that are always available. Although these files can be accessed using their corresponding numeric file ids, you should always use the symbolic names given in the table below, since it will make your programs easier to understand.
stdin ()
Return the numeric value corresponding to the standard input stream.
When Octave is used interactively, stdin is filtered through the command line editing functions.

See also: [stdout], page 294, [stderr], page 294.
stdout ()
Return the numeric value corresponding to the standard output stream.
Data written to the standard output may be filtered through the pager.
See also: [stdin], page 294, [stderr], page 294, [page_screen_output], page 276.
stderr ()
Return the numeric value corresponding to the standard error stream.
Even if paging is turned on, the standard error is not sent to the pager. It is useful for error messages and prompts.
See also: [stdin], page 294, [stdout], page 294.

### 14.2.1 Opening and Closing Files

When reading data from a file it must be opened for reading first, and likewise when writing to a file. The fopen function returns a pointer to an open file that is ready to be read or written. Once all data has been read from or written to the opened file it should be closed.

The fclose function does this. The following code illustrates the basic pattern for writing to a file, but a very similar pattern is used when reading a file.

```
filename = "myfile.txt";
fid = fopen (filename, "w");
# Do the actual I/O here...
fclose (fid);
```

fid $=$ fopen (name)
fid $=$ fopen (name, mode)
fid $=$ fopen (name, mode, arch)
fid $=$ fopen (name, mode, arch, encoding)
[fid, msg] = fopen (...)
fid_list $=$ fopen ("all")
[file, mode, arch, encoding] = fopen (fid)

Open a file for low-level I/O or query open files and file descriptors.
The first form of the fopen function opens the named file with the specified mode (read-write, read-only, etc.), architecture interpretation (IEEE big endian, IEEE little endian, etc.) and file encoding, and returns an integer value that may be used to refer to the file later. If an error occurs, fid is set to -1 and $m s g$ contains the corresponding system error message. The mode is a one or two character string that specifies whether the file is to be opened for reading, writing, or both. The encoding is a character string with a valid encoding identifier. This encoding is used when strings are read from or written to the file. By default, the same encoding specified for reading .m files is used for interpreting user files.
The second form of the fopen function returns a vector of file ids corresponding to all the currently open files, excluding the stdin, stdout, and stderr streams.
The third form of the fopen function returns information about the open file given its file id.
For example,

```
myfile = fopen ("splat.dat", "r", "ieee-le");
```

opens the file splat. dat for reading. If necessary, binary numeric values will be read assuming they are stored in IEEE format with the least significant bit first, and then converted to the native representation.
Opening a file that is already open simply opens it again and returns a separate file id. It is not an error to open a file several times, though writing to the same file through several different file ids may produce unexpected results.
The possible values of mode are 'r' (default)

Open a file for reading.
'w' Open a file for writing. The previous contents are discarded.
'a' Open or create a file for writing at the end of the file.
' $r+$ ' Open an existing file for reading and writing.
'w+' Open a file for reading or writing. The previous contents are discarded.
'a+' Open or create a file for reading or writing at the end of the file.
Append a " t " to the mode string to open the file in text mode or a "b" to open in binary mode. On Windows systems, text mode reading and writing automatically converts linefeeds to the appropriate line end character for the system (carriage-return linefeed on Windows). The default when no mode is specified is binary.
Additionally, you may append a "z" to the mode string to open a gzipped file for reading or writing. For this to be successful, you must also open the file in binary mode.

The parameter arch is a string specifying the default data format for the file. Valid values for arch are:

```
"native" or "n" (default)
```

The format of the current machine.
"ieee-be" or "b"
IEEE big endian format.
"ieee-le" or "1"
IEEE little endian format.
When opening a new file that does not yet exist, permissions will be set to 0666 umask.
Compatibility Note: Octave opens files using buffered I/O. Small writes are accumulated until an internal buffer is filled, and then everything is written in a single operation. This is very efficient and improves performance. MATLAB, however, opens files using flushed I/O where every write operation is immediately performed. If the write operation must be performed immediately after data has been written then the write should be followed by a call to fflush to flush the internal buffer.

See also: [fclose], page 296, [fgets], page 298, [fgetl], page 297, [fscanf], page 304, [fread], page 308, [fputs], page 297, [fdisp], page 283, [fprintf], page 299, [fwrite], page 310, [fskipl], page 298, [fseek], page 313, [frewind], page 314, [ftell], page 313, [feof], page 312, [ferror], page 313, [fclear], page 313, [fflush], page 276, [freport], page 313, [umask], page 955 .
fclose (fid)
fclose ("all")
status = fclose ("all")
Close the file specified by the file descriptor fid.
If successful, fclose returns 0 , otherwise, it returns -1 . The second form of the fclose call closes all open files except stdin, stdout, stderr, and any FIDs associated with gnuplot.
See also: [fopen], page 295, [fflush], page 276, [freport], page 313.
is_valid_file_id (fid)
Return true if fid refers to an open file.
See also: [freport], page 313, [fopen], page 295.

### 14.2.2 Simple Output

Once a file has been opened for writing a string can be written to the file using the fputs function. The following example shows how to write the string 'Free Software is needed for Free Science' to the file 'free.txt'.
filename = "free.txt";
fid $=$ fopen (filename, "w");
fputs (fid, "Free Software is needed for Free Science");
fclose (fid);
fputs (fid, string)
status = fputs (fid, string)
Write the string string to the file with file descriptor fid.
The string is written to the file with no additional formatting. Use fdisp instead to automatically append a newline character appropriate for the local machine.
Return a non-negative number on success or EOF on error.
See also: [fdisp], page 283, [fprintf], page 299, [fwrite], page 310, [fopen], page 295.
A function much similar to fputs is available for writing data to the screen. The puts function works just like fputs except it doesn't take a file pointer as its input.

```
puts (string)
status = puts (string)
```

Write a string to the standard output with no formatting.
The string is written verbatim to the standard output. Use disp to automatically append a newline character appropriate for the local machine.
Return a non-negative number on success and EOF on error.
See also: [fputs], page 297, [disp], page 271.

### 14.2.3 Line-Oriented Input

To read from a file it must be opened for reading using fopen. Then a line can be read from the file using fgetl as the following code illustrates

```
fid = fopen ("free.txt");
txt = fgetl (fid)
    Free Software is needed for Free Science
fclose (fid);
```

This of course assumes that the file 'free.txt' exists and contains the line 'Free Software is needed for Free Science'.
str $=$ fgetl (fid)
str = fgetl (fid, len)
Read characters from a file, stopping after a newline, or EOF, or len characters have been read.
The characters read, excluding the possible trailing newline, are returned as a string.
If len is omitted, fgetl reads until the next newline character.
If there are no more characters to read, fgetl returns -1 .

To read a line and return the terminating newline, see [fgets], page 298.
See also: [fgets], page 298, [fscanf], page 304, [fread], page 308, [fopen], page 295.

```
str = fgets (fid)
```

str $=$ fgets (fid, len)

Read characters from a file, stopping after a newline, or EOF, or len characters have been read.
The characters read, including the possible trailing newline, are returned as a string. If len is omitted, fgets reads until the next newline character.
If there are no more characters to read, fgets returns -1 .
To read a line and discard the terminating newline, see [fgetl], page 297.
See also: [fputs], page 297, [fgetl], page 297, [fscanf], page 304, [fread], page 308, [fopen], page 295.

```
nlines = fskipl (fid)
nlines = fskipl (fid, count)
nlines = fskipl (fid, Inf)
```

Read and skip count lines from the file specified by the file descriptor fid.
fskipl discards characters until an end-of-line is encountered exactly count-times, or until the end-of-file marker is found.
If count is omitted, it defaults to 1 . count may also be Inf, in which case lines are skipped until the end of the file. This form is suitable for counting the number of lines in a file.
Returns the number of lines skipped (end-of-line sequences encountered).
See also: [fgetl], page 297, [fgets], page 298, [fscanf], page 304, [fopen], page 295.

### 14.2.4 Formatted Output

This section describes how to call printf and related functions.
The following functions are available for formatted output. They are modeled after the C language functions of the same name, but they interpret the format template differently in order to improve the performance of printing vector and matrix values.

Implementation Note: For compatibility with matlab, escape sequences in the template string (e.g., "\n" => newline) are expanded even when the template string is defined with single quotes.
printf (template,...)
Print optional arguments under the control of the template string template to the stream stdout and return the number of characters printed.
See the Formatted Output section of the GNU Octave manual for a complete description of the syntax of the template string.
Implementation Note: For compatibility with matlab, escape sequences in the template string (e.g., " $\backslash \mathrm{n} " ~=>~ n e w l i n e) ~ a r e ~ e x p a n d e d ~ e v e n ~ w h e n ~ t h e ~ t e m p l a t e ~ s t r i n g ~ i s ~$ defined with single quotes.
See also: [fprintf], page 299, [sprintf], page 299, [scanf], page 305.
fprintf (fid, template,...)
fprintf (template,...)
numbytes = fprintf (...)
This function is equivalent to printf, except that the output is written to the file descriptor fid instead of stdout.
If fid is omitted, the output is written to stdout making the function exactly equivalent to printf.
The optional output returns the number of bytes written to the file.
Implementation Note: For compatibility with matlab, escape sequences in the template string (e.g., " $\backslash \mathrm{n} " ~=>$ newline) are expanded even when the template string is defined with single quotes.
See also: [fputs], page 297, [fdisp], page 283, [fwrite], page 310, [fscanf], page 304, [printf], page 298, [sprintf], page 299, [fopen], page 295.
sprintf (template, ...)
This is like printf, except that the output is returned as a string.
Unlike the C library function, which requires you to provide a suitably sized string as an argument, Octave's sprintf function returns the string, automatically sized to hold all of the items converted.
Implementation Note: For compatibility with matlab, escape sequences in the template string (e.g., " $\backslash \mathrm{n} "=>$ newline) are expanded even when the template string is defined with single quotes.
See also: [printf], page 298, [fprintf], page 299, [sscanf], page 305.
The printf function can be used to print any number of arguments. The template string argument you supply in a call provides information not only about the number of additional arguments, but also about their types and what style should be used for printing them.

Ordinary characters in the template string are simply written to the output stream as-is, while conversion specifications introduced by a ' $\%$ ' character in the template cause subsequent arguments to be formatted and written to the output stream. For example,

```
pct = 37;
filename = "foo.txt";
printf ("Processed %d%% of '%s'.\nPlease be patient.\n",
    pct, filename);
```

produces output like
Processed $37 \%$ of 'foo.txt'.
Please be patient.
This example shows the use of the '\%d' conversion to specify that a scalar argument should be printed in decimal notation, the ' $\%$ s' conversion to specify printing of a string argument, and the '\%\%' conversion to print a literal ' $\%$ ' character.

There are also conversions for printing an integer argument as an unsigned value in octal, decimal, or hexadecimal radix ( $\%$ \% ${ }^{\prime}$ ', $\% \%^{\prime}$ ', or ' $\% \mathrm{x}^{\prime}$, respectively) ; or as a character value ( ${ }^{\prime} \% c^{\prime}$ ).

Floating-point numbers can be printed in normal, fixed-point notation using the $\%$ \% ${ }^{\prime}$ conversion or in exponential notation using the '\%e' conversion. The ' $\% \mathrm{~g}$ ' conversion uses either ' $\%$ e' or ' $\%$ f ' format, depending on what is more appropriate for the magnitude of the particular number.

You can control formatting more precisely by writing modifiers between the ' $\%$ ' and the character that indicates which conversion to apply. These slightly alter the ordinary behavior of the conversion. For example, most conversion specifications permit you to specify a minimum field width and a flag indicating whether you want the result left- or right-justified within the field.

The specific flags and modifiers that are permitted and their interpretation vary depending on the particular conversion. They're all described in more detail in the following sections.

### 14.2.5 Output Conversion for Matrices

When given a matrix value, Octave's formatted output functions cycle through the format template until all the values in the matrix have been printed. For example:

```
printf ("%4.2f %10.2e %8.4g\n", hilb (3));
    \dashv1.00 5.00e-01 0.3333
    \dashv 0.50 3.33e-01 0.25
    \dashv 0.33 2.50e-01 0.2
```

If more than one value is to be printed in a single call, the output functions do not return to the beginning of the format template when moving on from one value to the next. This can lead to confusing output if the number of elements in the matrices are not exact multiples of the number of conversions in the format template. For example:

```
printf ("%4.2f %10.2e %8.4g\n", [1, 2], [3, 4]);
    \dashv1.00 2.00e+00 3
    \dashv4.00
```

If this is not what you want, use a series of calls instead of just one.

### 14.2.6 Output Conversion Syntax

This section provides details about the precise syntax of conversion specifications that can appear in a printf template string.

Characters in the template string that are not part of a conversion specification are printed as-is to the output stream.

The conversion specifications in a printf template string have the general form:

```
% flags width [ . precision ] type conversion
```

For example, in the conversion specifier '\%-10.81d', the '-' is a flag, ' 10 ' specifies the field width, the precision is ' 8 ', the letter ' 1 ' is a type modifier, and ' $d$ ' specifies the conversion style. (This particular type specifier says to print a numeric argument in decimal notation, with a minimum of 8 digits left-justified in a field at least 10 characters wide.)

In more detail, output conversion specifications consist of an initial '\%' character followed in sequence by:

- Zero or more flag characters that modify the normal behavior of the conversion specification.
- An optional decimal integer specifying the minimum field width. If the normal conversion produces fewer characters than this, the field is padded with spaces to the specified width. This is a minimum value; if the normal conversion produces more characters than this, the field is not truncated. Normally, the output is right-justified within the field.
You can also specify a field width of ' $*$ '. This means that the next argument in the argument list (before the actual value to be printed) is used as the field width. The value is rounded to the nearest integer. If the value is negative, this means to set the '-' flag (see below) and to use the absolute value as the field width.
- An optional precision to specify the number of digits to be written for the numeric conversions. If the precision is specified, it consists of a period ('.') followed optionally by a decimal integer (which defaults to zero if omitted).
You can also specify a precision of ' $*$ '. This means that the next argument in the argument list (before the actual value to be printed) is used as the precision. The value must be an integer, and is ignored if it is negative.
- An optional type modifier character. This character is ignored by Octave's printf function, but is recognized to provide compatibility with the C language printf.
- A character that specifies the conversion to be applied.

The exact options that are permitted and how they are interpreted vary between the different conversion specifiers. See the descriptions of the individual conversions for information about the particular options that they use.

### 14.2.7 Table of Output Conversions

Here is a table summarizing what all the different conversions do:
${ }^{`} \% \mathrm{~d}$ ', ${ }^{`} \% \mathrm{i}$ ' Print an integer as a signed decimal number. See Section 14.2 .8 [Integer Con-
 different when used with scanf for input (see Section 14.2.13 [Table of Input Conversions], page 306).
'\%o' Print an integer as an unsigned octal number. See Section 14.2.8 [Integer Conversions], page 302, for details.
‘\%u’ Print an integer as an unsigned decimal number. See Section 14.2.8 [Integer Conversions], page 302, for details.
${ }^{6} \% \mathrm{x}$ ', $\% \mathrm{\%}$ ' $\quad$ Print an integer as an unsigned hexadecimal number. ' $\% \mathrm{x}$ ' uses lowercase letters and '\%X' uses uppercase. See Section 14.2.8 [Integer Conversions], page 302, for details.
(\%f $\quad$ Print a floating-point number in normal (fixed-point) notation. See Section 14.2.9 [Floating-Point Conversions], page 303, for details.
'\%e', '\%E' Print a floating-point number in exponential notation. '\%e' uses lowercase letters and '\%E’ uses uppercase. See Section 14.2.9 [Floating-Point Conversions], page 303 , for details.
'\%g', '\%G' Print a floating-point number in either normal (fixed-point) or exponential notation, whichever is more appropriate for its magnitude. ' $\%$ ' ' uses lowercase letters and '\%G’ uses uppercase. See Section 14.2.9 [Floating-Point Conversions], page 303 , for details.
'\%c' Print a single character. See Section 14.2.10 [Other Output Conversions], page 303.
‘\%s' Print a string. See Section 14.2.10 [Other Output Conversions], page 303.
(\%\%' Print a literal '\%' character. See Section 14.2.10 [Other Output Conversions], page 303.

If the syntax of a conversion specification is invalid, unpredictable things will happen, so don't do this. In particular, matlab allows a bare percentage sign '\%' with no subsequent conversion character. Octave will emit an error and stop if it sees such code. When the string variable to be processed cannot be guaranteed to be free of potential format codes it is better to use the two argument form of any of the printf functions and set the format string to \%s. Alternatively, for code which is not required to be backwards-compatible with matlab the Octave function puts or disp can be used.

```
printf (strvar); # Unsafe if strvar contains format codes
printf ("%s", strvar); # Safe
puts (strvar); # Safe
```

If there aren't enough function arguments provided to supply values for all the conversion specifications in the template string, or if the arguments are not of the correct types, the results are unpredictable. If you supply more arguments than conversion specifications, the extra argument values are simply ignored; this is sometimes useful.

### 14.2.8 Integer Conversions

 specifications. These conversions print integers in various formats.

The ' $\% \mathrm{~d}$ ' and ' $\% \mathrm{i}$ ' conversion specifications both print an numeric argument as a signed decimal number; while '\% $\%$ ', ' $\% \mathrm{u}$ ', and ${ }^{\prime} \%$ ' print the argument as an unsigned octal, decimal, or hexadecimal number (respectively). The ' $\% \mathrm{X}$ ' conversion specification is just like ' $\% \mathrm{x}$ ' except that it uses the characters 'ABCDEF' as digits instead of 'abcdef'.

The following flags are meaningful:
'-' Left-justify the result in the field (instead of the normal right-justification).
' + ' For the signed '\% d ' and ' $\% \mathrm{i}$ ' conversions, print a plus sign if the value is positive.
‘, For the signed $\% \mathrm{~m}$ ' and ' $\% \mathrm{i}$ ' conversions, if the result doesn't start with a plus or minus sign, prefix it with a space character instead. Since the ' + ' flag ensures that the result includes a sign, this flag is ignored if you supply both of them.
'\#' For the ' $\%$ ' ' conversion, this forces the leading digit to be ' 0 ', as if by increasing the precision. For ' $\% \mathrm{x}$ ' or ' $\% \mathrm{X}$ ', this prefixes a leading ' Ox ' or ' OX ' (respectively) to the result. This doesn't do anything useful for the ' $\% \mathrm{~d}$ ', ' $\% \mathrm{i}$ ', or ${ }^{~} \% \mathrm{\%}$ ' conversions.
' 0 ' Pad the field with zeros instead of spaces. The zeros are placed after any indication of sign or base. This flag is ignored if the '-' flag is also specified, or if a precision is specified.

If a precision is supplied, it specifies the minimum number of digits to appear; leading zeros are produced if necessary. If you don't specify a precision, the number is printed with as many digits as it needs. If you convert a value of zero with an explicit precision of zero, then no characters at all are produced.

### 14.2.9 Floating-Point Conversions

This section discusses the conversion specifications for floating-point numbers: the ' $\% \mathrm{f}$ ', '\%e', ‘\%E', '\% $\mathrm{g}^{\prime}$, and $\% \mathrm{G}$ ' conversions.

The '\%f' conversion prints its argument in fixed-point notation, producing output of the form $[-] d d d . d d d$, where the number of digits following the decimal point is controlled by the precision you specify.

The '\%e' conversion prints its argument in exponential notation, producing output of the form $[-] d . d d d e[+\mid-] d d$. Again, the number of digits following the decimal point is controlled by the precision. The exponent always contains at least two digits. The '\%E' conversion is similar but the exponent is marked with the letter ' $E$ ' instead of ' $e$ '.

The ' $\% \mathrm{~g}$ ' and '\%G' conversions print the argument in the style of ' $\%$ ' or ${ }^{\prime} \% \mathrm{E}$ ' (respectively) if the exponent would be less than -4 or greater than or equal to the precision; otherwise they use the '\%f' style. Trailing zeros are removed from the fractional portion of the result and a decimal-point character appears only if it is followed by a digit.

The following flags can be used to modify the behavior:
'-' Left-justify the result in the field. Normally the result is right-justified.
‘+' Always include a plus or minus sign in the result.
، If the result doesn't start with a plus or minus sign, prefix it with a space instead. Since the ' + ' flag ensures that the result includes a sign, this flag is ignored if you supply both of them.
'\#' Specifies that the result should always include a decimal point, even if no digits follow it. For the ' $\% \mathrm{~g}$ ' and ' $\% \mathrm{G}$ ' conversions, this also forces trailing zeros after the decimal point to be left in place where they would otherwise be removed.
' 0 ' Pad the field with zeros instead of spaces; the zeros are placed after any sign. This flag is ignored if the '-' flag is also specified.

The precision specifies how many digits follow the decimal-point character for the '\%f', ${ }^{\prime} \%$ e', and ' $\% \mathrm{E}$ ’ conversions. For these conversions, the default precision is 6 . If the precision is explicitly 0 , this suppresses the decimal point character entirely. For the ' $\% \mathrm{~g}$ ' and ' $\% \mathrm{G}$ ' conversions, the precision specifies how many significant digits to print. Significant digits are the first digit before the decimal point, and all the digits after it. If the precision is 0 or not specified for ' $\% \mathrm{~g}$ ' or ' $\% \mathrm{G}$ ', it is treated like a value of 1 . If the value being printed cannot be expressed precisely in the specified number of digits, the value is rounded to the nearest number that fits.

### 14.2.10 Other Output Conversions

This section describes miscellaneous conversions for printf.

The '\%c' conversion prints a single character. The '-' flag can be used to specify leftjustification in the field, but no other flags are defined, and no precision or type modifier can be given. For example:

```
printf ("%c%c%c%c%c", "h", "e", "l", "l", "o");
```

prints 'hello'.
The '\%s' conversion prints a string. The corresponding argument must be a string. A precision can be specified to indicate the maximum number of characters to write; otherwise characters in the string up to but not including the terminating null character are written to the output stream. The '-' flag can be used to specify left-justification in the field, but no other flags or type modifiers are defined for this conversion. For example:
printf ("\%3s\%-6s", "no", "where");
prints ' nowhere ' (note the leading and trailing spaces).

### 14.2.11 Formatted Input

Octave provides the scanf, fscanf, and sscanf functions to read formatted input. There are two forms of each of these functions. One can be used to extract vectors of data from a file, and the other is more 'C-like'.
[val, count, errmsg] = fscanf (fid, template, size)
[v1, v2, ..., count, errmsg] = fscanf (fid, template, "C")
In the first form, read from fid according to template, returning the result in the matrix val.

The optional argument size specifies the amount of data to read and may be one of
Inf Read as much as possible, returning a column vector.
nr $\quad$ Read up to nr elements, returning a column vector.
[nr, Inf] Read as much as possible, returning a matrix with nr rows. If the number of elements read is not an exact multiple of $n r$, the last column is padded with zeros.
[nr, nc] Read up to $n r * n c$ elements, returning a matrix with $n r$ rows. If the number of elements read is not an exact multiple of nr, the last column is padded with zeros.

If size is omitted, a value of Inf is assumed.
A string is returned if template specifies only character conversions.
The number of items successfully read is returned in count.
If an error occurs, errmsg contains a system-dependent error message.
In the second form, read from fid according to template, with each conversion specifier in template corresponding to a single scalar return value. This form is more "C-like", and also compatible with previous versions of Octave. The number of successful conversions is returned in count
See the Formatted Input section of the GNU Octave manual for a complete description of the syntax of the template string.
See also: [fgets], page 298, [fgetl], page 297, [fread], page 308, [scanf], page 305, [sscanf], page 305, [fopen], page 295.
[val, count, errmsg] = scanf (template, size)
[v1, v2, ..., count, errmsg] = scanf (template, "C")
This is equivalent to calling fscanf with fid $=$ stdin.
It is currently not useful to call scanf in interactive programs.
See also: [fscanf], page 304, [sscanf], page 305, [printf], page 298.
[val, count, errmsg, pos] = sscanf (string, template, size)
[v1, v2, ..., count, errmsg] = sscanf (string, template, "C")
This is like fscanf, except that the characters are taken from the string string instead of from a stream.
Reaching the end of the string is treated as an end-of-file condition. In addition to the values returned by fscanf, the index of the next character to be read is returned in pos.

See also: [fscanf], page 304, [scanf], page 305, [sprintf], page 299.
Calls to scanf are superficially similar to calls to printf in that arbitrary arguments are read under the control of a template string. While the syntax of the conversion specifications in the template is very similar to that for printf, the interpretation of the template is oriented more towards free-format input and simple pattern matching, rather than fixedfield formatting. For example, most scanf conversions skip over any amount of "white space" (including spaces, tabs, and newlines) in the input file, and there is no concept of precision for the numeric input conversions as there is for the corresponding output conversions. Ordinarily, non-whitespace characters in the template are expected to match characters in the input stream exactly. For example, note that sscanf parses the string and whitespace differently when using mixed numeric and string output types:

```
teststr = "1 is a lonely number";
sscanf (teststr, "%s is a %s")
=> 1lonelynumber
sscanf (teststr, "%g is a %s")
#
        1
    108
        111
        110
        101
        108
        121
[a, b, c] = sscanf ("1 is a lonely number", "%g is a %s %s", "C")
a = 1
=> b = lonely
c = number
```

When a matching failure occurs, scanf returns immediately, leaving the first nonmatching character as the next character to be read from the stream, and scanf returns all the items that were successfully converted.

The formatted input functions are not used as frequently as the formatted output functions. Partly, this is because it takes some care to use them properly. Another reason is that it is difficult to recover from a matching error.

The specific flags and modifiers that are permitted in the template string and their interpretation are all described in more detail in the following sections.

### 14.2.12 Input Conversion Syntax

A scanf template string is a string that contains ordinary multibyte characters interspersed with conversion specifications that start with ' $\%$ '.

Any whitespace character in the template causes any number of whitespace characters in the input stream to be read and discarded. The whitespace characters that are matched need not be exactly the same whitespace characters that appear in the template string. For example, write ', ' in the template to recognize a comma with optional whitespace before and after.

Other characters in the template string that are not part of conversion specifications must match characters in the input stream exactly; if this is not the case, a matching failure occurs.

The conversion specifications in a scanf template string have the general form:

```
% flags width type conversion
```

In more detail, an input conversion specification consists of an initial '\%' character followed in sequence by:

- An optional flag character '*', which says to ignore the text read for this specification. When scanf finds a conversion specification that uses this flag, it reads input as directed by the rest of the conversion specification, but it discards this input, does not return any value, and does not increment the count of successful assignments.
- An optional decimal integer that specifies the maximum field width. Reading of characters from the input stream stops either when this maximum is reached or when a non-matching character is found, whichever happens first. Most conversions discard initial whitespace characters, and these discarded characters don't count towards the maximum field width. Conversions that do not discard initial whitespace are explicitly documented.
- An optional type modifier character. This character is ignored by Octave's scanf function, but is recognized to provide compatibility with the C language scanf.
- A character that specifies the conversion to be applied.

The exact options that are permitted and how they are interpreted vary between the different conversion specifiers. See the descriptions of the individual conversions in Section 14.2.13 [Table of Input Conversions], page 306, for information about the particular options that they allow.

### 14.2.13 Table of Input Conversions

Here is a table that summarizes the various conversion specifications:
‘\% d' Matches an optionally signed integer written in decimal. See Section 14.2.14 [Numeric Input Conversions], page 307.
‘\%i' Matches an optionally signed integer in any of the formats that the C language defines for specifying an integer constant. See Section 14.2.14 [Numeric Input Conversions], page 307.
‘\% ${ }^{\prime}$ Matches an unsigned integer written in octal radix. See Section 14.2.14 [Numeric Input Conversions], page 307.
‘\%u' Matches an unsigned integer written in decimal radix. See Section 14.2.14 [Numeric Input Conversions], page 307.
${ }^{6} \% \mathrm{x}$ ', ${ }^{6} \% \mathrm{X}$ ’ Matches an unsigned integer written in hexadecimal radix. See Section 14.2.14 [Numeric Input Conversions], page 307.

Matches an optionally signed floating-point number. See Section 14.2.14 [Numeric Input Conversions], page 307.
'\%s' Matches a string containing only non-whitespace characters. See Section 14.2.15 [String Input Conversions], page 308.
‘\% $\%$ ' Matches a string of one or more characters; the number of characters read is controlled by the maximum field width given for the conversion. See Section 14.2.15 [String Input Conversions], page 308.
‘\%\%' This matches a literal ${ }^{\circ} \%$ ' character in the input stream. No corresponding argument is used.

If the syntax of a conversion specification is invalid, the behavior is undefined. If there aren't enough function arguments provided to supply addresses for all the conversion specifications in the template strings that perform assignments, or if the arguments are not of the correct types, the behavior is also undefined. On the other hand, extra arguments are simply ignored.

### 14.2.14 Numeric Input Conversions

This section describes the scanf conversions for reading numeric values.
The '\%d' conversion matches an optionally signed integer in decimal radix.
The ' $\% \mathrm{i}$ ' conversion matches an optionally signed integer in any of the formats that the C language defines for specifying an integer constant.

For example, any of the strings ' 10 ', ' $0 x a$ ', or ' 012 ' could be read in as integers under the ' $\%$ ' ' conversion. Each of these specifies a number with decimal value 10.

The ' $\%$ ', ' $\%$ u', and ' $\%$ x' conversions match unsigned integers in octal, decimal, and hexadecimal radices, respectively.

The ' $\% \mathrm{X}$ ' conversion is identical to the ' $\%$ x' conversion. They both permit either uppercase or lowercase letters to be used as digits.

By default, integers are read as 32 -bit quantities. With the ' h ' modifier, 16 -bit integers are used, and with the ' 1 ' modifier, 64 -bit integers are used.

The ${ }^{\prime} \% e^{\prime},{ }^{\prime} \% \mathrm{f}$ ', ${ }^{\prime} \% \mathrm{~g}$ ', ${ }^{\prime} \% \mathrm{E}$ ', and ${ }^{\prime} \% \mathrm{G}$ ' conversions match optionally signed floating-point numbers. All five conversion specifications behave identically, and will read in numerical values of any floating point display style.

### 14.2.15 String Input Conversions

This section describes the scanf input conversions for reading string and character values: ' $\% s^{\prime}$ ' and ' $\% c^{\prime}$ '.

The ' $\%$ c' conversion is the simplest: it matches a fixed number of characters, always. The maximum field with says how many characters to read; if you don't specify the maximum, the default is 1 . This conversion does not skip over initial whitespace characters. It reads precisely the next $n$ characters, and fails if it cannot get that many.

The '\%s' conversion matches a string of non-whitespace characters. It skips and discards initial whitespace, but stops when it encounters more whitespace after having read something.

For example, reading the input:

```
hello, world
```

with the conversion $\% 10 \mathrm{c}$ ' produces " hello, wo", but reading the same input with the conversion '\%10s' produces "hello,".

### 14.2.16 Binary I/O

Octave can read and write binary data using the functions fread and fwrite, which are patterned after the standard C functions with the same names. They are able to automatically swap the byte order of integer data and convert among the supported floating point formats as the data are read.

```
val = fread (fid)
val = fread (fid, size)
val = fread (fid, size, precision)
val = fread (fid, size, precision, skip)
val = fread (fid, size, precision, skip, arch)
[val, count] = fread (...)
```

    Read binary data from the file specified by the file descriptor fid.
    The optional argument size specifies the amount of data to read and may be one of
    Inf Read as much as possible, returning a column vector.
    \(n r \quad\) Read up to \(n r\) elements, returning a column vector.
    [ \(n r\), Inf] Read as much as possible, returning a matrix with \(n r\) rows. If the number
                        of elements read is not an exact multiple of \(n r\), the last column is padded
                        with zeros.
    [ \(n r, n c\) ] Read up to \(n r * n c\) elements, returning a matrix with \(n r\) rows. If the
                number of elements read is not an exact multiple of \(n r\), the last column
                is padded with zeros.
    If size is omitted, a value of Inf is assumed.
    The optional argument precision is a string specifying the type of data to read and may be one of
"uint8" (default)
8 -bit unsigned integer.

```
"int8"
"integer*1"
            8-bit signed integer.
"uint16"
"ushort"
"unsigned short"
            16-bit unsigned integer.
"int16"
"integer*2"
"short" 16-bit signed integer.
"uint"
"uint32"
"unsigned int"
"ulong"
"unsigned long"
            32-bit unsigned integer.
"int"
"int32"
"integer*4"
"long" 32-bit signed integer.
"uint64" 64-bit unsigned integer.
"int64"
"integer*8"
    64-bit signed integer.
"single"
"float"
"float32"
"real*4" 32-bit floating point number.
"double"
"float64"
"real*8" 64-bit floating point number.
"char"
"char*1" 8-bit single character.
"uchar"
"unsigned char"
    8-bit unsigned character.
"schar"
"signed char"
    8-bit signed character.
```

The default precision is "uint8".

The precision argument may also specify an optional repeat count. For example, ' $32 *$ single' causes fread to read a block of 32 single precision floating point numbers. Reading in blocks is useful in combination with the skip argument.
The precision argument may also specify a type conversion. For example, 'int16=>int32' causes fread to read 16-bit integer values and return an array of 32-bit integer values. By default, fread returns a double precision array. The special form '*TYPE' is shorthand for 'TYPE=>TYPE'.
The conversion and repeat counts may be combined. For example, the specification ' $32 *$ single=>single' causes fread to read blocks of single precision floating point values and return an array of single precision values instead of the default array of double precision values.
The optional argument skip specifies the number of bytes to skip after each element (or block of elements) is read. If it is not specified, a value of 0 is assumed. If the final block read is not complete, the final skip is omitted. For example,

```
fread (f, 10, "3*single=>single", 8)
```

will omit the final 8 -byte skip because the last read will not be a complete block of 3 values.
The optional argument arch is a string specifying the data format for the file. Valid values are

```
"native" or "n"
```

The format of the current machine.
"ieee-be" or "b"
IEEE big endian.
"ieee-le" or "1"
IEEE little endian.
If no arch is given the value used in the call to fopen which created the file descriptor is used. Otherwise, the value specified with fread overrides that of fopen and determines the data format.
The output argument val contains the data read from the file.
The optional return value count contains the number of elements read.
See also: [fwrite], page 310, [fgets], page 298, [fgetl], page 297, [fscanf], page 304, [fopen], page 295.

```
fwrite (fid, data)
fwrite (fid, data, precision)
fwrite (fid, data, precision, skip)
fwrite (fid, data, precision, skip, arch)
count = fwrite (...)
```

Write data in binary form to the file specified by the file descriptor fid, returning the number of values count successfully written to the file.
The argument data is a matrix of values that are to be written to the file. The values are extracted in column-major order.

The remaining arguments precision, skip, and arch are optional, and are interpreted as described for fread.
The behavior of fwrite is undefined if the values in data are too large to fit in the specified precision.
See also: [fread], page 308, [fputs], page 297, [fprintf], page 299, [fopen], page 295.

### 14.2.17 Temporary Files

Sometimes one needs to write data to a file that is only temporary. This is most commonly used when an external program launched from within Octave needs to access data. When Octave exits all temporary files will be deleted, so this step need not be executed manually.
[fid, name, msg] = mkstemp ("template")
[fid, name, msg] = mkstemp ("template", delete)
Return the file descriptor fid corresponding to a new temporary file with a unique name created from template.

The last six characters of template must be "XXXXXX" and these are replaced with a string that makes the filename unique. The file is then created with mode read/write and permissions that are system dependent (on GNU/Linux systems, the permissions will be 0600 for versions of glibc 2.0.7 and later). The file is opened in binary mode and with the 0_EXCL flag.
If the optional argument delete is supplied and is true, the file will be deleted automatically when Octave exits.
If successful, fid is a valid file ID, name is the name of the file, and msg is an empty string. Otherwise, fid is -1 , name is empty, and msg contains a system-dependent error message.
See also: [tempname], page 311, [tempdir], page 312, [ $\mathrm{P}_{-}$tmpdir], page 312, [tmpfile], page 311, [fopen], page 295.
[fid, msg] = tmpfile ()
Return the file ID corresponding to a new temporary file with a unique name.
The file is opened in binary read/write ( $\mathrm{w}+\mathrm{b} \mathrm{b}$ ) mode and will be deleted automatically when it is closed or when Octave exits.
If successful, fid is a valid file ID and msg is an empty string. Otherwise, fid is -1 and msg contains a system-dependent error message.
See also: [tempname], page 311, [mkstemp], page 311, [tempdir], page 312, [P_tmpdir], page 312.

```
fname = tempname ()
fname = tempname (dir)
fname = tempname (dir, prefix)
```

Return a unique temporary filename as a string.
If prefix is omitted, a value of "oct-" is used.
If dir is also omitted, the default directory for temporary files (P_tmpdir) is used. If dir is provided, it must exist, otherwise the default directory for temporary files is used.

Programming Note: Because the named file is not opened by tempname, it is possible, though relatively unlikely, that it will not be available by the time your program attempts to open it. If this is a concern, see [tmpfile], page 311.

See also: [mkstemp], page 311, [tempdir], page 312, [P_tmpdir], page 312, [tmpfile], page 311.
dir $=$ tempdir ()
Return the name of the host system's directory for temporary files.
The directory name is taken first from the environment variable TMPDIR. If that does not exist the system default returned by P_tmpdir is used.

See also: [P_tmpdir], page 312, [tempname], page 311, [mkstemp], page 311, [tmpfile], page 311.

P_tmpdir ()
Return the name of the host system's default directory for temporary files.
Programming Note: The value returned by P_tmpdir is always the default location. This value may not agree with that returned from tempdir if the user has overridden the default with the TMPDIR environment variable.

See also: [tempdir], page 312, [tempname], page 311, [mkstemp], page 311, [tmpfile], page 311 .

### 14.2.18 End of File and Errors

Once a file has been opened its status can be acquired. As an example the feof functions determines if the end of the file has been reached. This can be very useful when reading small parts of a file at a time. The following example shows how to read one line at a time from a file until the end has been reached.

```
filename = "myfile.txt";
fid = fopen (filename, "r");
while (! feof (fid) )
    text_line = fgetl (fid);
endwhile
fclose (fid);
```

Note that in some situations it is more efficient to read the entire contents of a file and then process it, than it is to read it line by line. This has the potential advantage of removing the loop in the above code.

## status $=$ feof (fid)

Return 1 if an end-of-file condition has been encountered for the file specified by file descriptor fid and 0 otherwise.
Note that feof will only return 1 if the end of the file has already been encountered, not if the next read operation will result in an end-of-file condition.
See also: [fread], page 308, [frewind], page 314, [fseek], page 313, [fclear], page 313, [fopen], page 295.
$m s g=$ ferror (fid)
[msg, err] $=$ ferror (fid)
[...] $=$ ferror (fid, "clear")
Query the error status of the stream specified by file descriptor fid.
If an error condition exists then return a string msg describing the error. Otherwise, return an empty string "".
The second input "clear" is optional. If supplied, the error state on the stream will be cleared.
The optional second output is a numeric indication of the error status. err is 1 if an error condition has been encountered and 0 otherwise.

Note that ferror indicates if an error has already occurred, not whether the next operation will result in an error condition.

See also: [fclear], page 313, [fopen], page 295.
fclear (fid)
Clear the stream state for the file specified by the file descriptor fid.
See also: [ferror], page 313, [fopen], page 295.
freport ()
Print a list of which files have been opened, and whether they are open for reading, writing, or both.
For example:

| freport () |  |  |  |  |
| ---: | :--- | :--- | :--- | :--- |
| $\dashv$ | number | mode | arch | name |
| $\dashv$ | ------ | ---- | --- | ---- |
| $\dashv$ | 0 | $r$ | ieee-le | stdin |
| $\dashv$ | 1 | w | ieee-le | stdout |
| -1 | 2 | w | ieee-le | stderr |
| $\dashv$ | 3 | $r$ | ieee-le | myfile |

See also: [fopen], page 295, [fclose], page 296, [is_valid_file_id], page 296.

### 14.2.19 File Positioning

Three functions are available for setting and determining the position of the file pointer for a given file.

```
pos = ftell (fid)
```

Return the position of the file pointer as the number of characters from the beginning of the file specified by file descriptor fid.

See also: [fseek], page 313, [frewind], page 314, [feof], page 312, [fopen], page 295.

```
fseek (fid, offset)
fseek (fid, offset, origin)
status = fseek (...)
```

Set the file pointer to the location offset within the file fid.

The pointer is positioned offset characters from the origin, which may be one of the predefined variables SEEK_SET (beginning), SEEK_CUR (current position), or SEEK_END (end of file) or strings "bof", "cof", or "eof". If origin is omitted, SEEK_SET is assumed. offset may be positive, negative, or zero but not all combinations of origin and offset can be realized.
fseek returns 0 on success and -1 on error.
See also: [fskipl], page 298, [frewind], page 314, [ftell], page 313, [fopen], page 295, [SEEK_SET], page 314, [SEEK_CUR], page 314, [SEEK_END], page 314.

## SEEK_SET ()

Return the numerical value to pass to fseek to position the file pointer relative to the beginning of the file.
See also: [SEEK_CUR], page 314, [SEEK_END], page 314, [fseek], page 313.
SEEK_CUR ()
Return the numerical value to pass to fseek to position the file pointer relative to the current position.

See also: [SEEK_SET], page 314, [SEEK_END], page 314, [fseek], page 313.
SEEK_END ()
Return the numerical value to pass to fseek to position the file pointer relative to the end of the file.

See also: [SEEK_SET], page 314, [SEEK_CUR], page 314, [fseek], page 313.
frewind (fid)
status = frewind (fid)
Move the file pointer to the beginning of the file specified by file descriptor fid.
frewind returns 0 for success, and -1 if an error is encountered. It is equivalent to fseek (fid, 0, SEEK_SET).

See also: [fseek], page 313, [ftell], page 313, [fopen], page 295.
The following example stores the current file position in the variable marker, moves the pointer to the beginning of the file, reads four characters, and then returns to the original position.

```
marker = ftell (myfile);
frewind (myfile);
fourch = fgets (myfile, 4);
fseek (myfile, marker, SEEK_SET);
```


## 15 Plotting

### 15.1 Introduction to Plotting

Earlier versions of Octave provided plotting through the use of gnuplot. This capability is still available. But, newer versions of Octave offer more modern plotting capabilities using OpenGL. Which plotting system is used is controlled by the graphics_toolkit function. See Section 15.4.8 [Graphics Toolkits], page 507.

The function call graphics_toolkit ("qt") selects the Qt/OpenGL system, graphics_toolkit ("fltk") selects the FLTK/OpenGL system, and graphics_toolkit ("gnuplot") selects the gnuplot system. The three systems may be used selectively through the use of the graphics_toolkit property of the graphics handle for each figure. This is explained in Section 15.3 [Graphics Data Structures], page 421.

Caution: The OpenGL-based toolkits use single precision variables internally which limits the maximum value that can be displayed to approximately $10^{38}$. If your data contains larger values you must use the gnuplot toolkit which supports values up to $10^{308}$. Similarly, single precision variables can accurately represent only $6-9$ base 10 digits. If your data contains very fine differences (approximately 1e-8) these cannot be resolved with the OpenGL-based graphics toolkits and the gnuplot toolkit is required.

Note: The gnuplot graphics toolkit uses the third party program gnuplot for plotting. The communication from Octave to gnuplot is done via a one-way pipe. This has implications for performance and functionality. Performance is significantly slower because the entire data set, which could be many megabytes, must be passed to gnuplot over the pipe. Functionality is negatively affected because the pipe is one-way from Octave to gnuplot. Octave has no way of knowing about user interactions with the plot window (be it resizing, moving, closing, or anything else). It is recommended not to interact with (or close) a gnuplot window if you will access the figure from Octave later on.

### 15.2 High-Level Plotting

Octave provides simple means to create many different types of two- and three-dimensional plots using high-level functions.

If you need more detailed control, see Section 15.3 [Graphics Data Structures], page 421, and Section 15.4 [Advanced Plotting], page 491.

### 15.2.1 Two-Dimensional Plots

The plot function allows you to create simple x-y plots with linear axes. For example,

```
x = -10:0.1:10;
plot (x, sin (x));
xlabel ("x");
ylabel ("sin (x)");
title ("Simple 2-D Plot");
```

displays a sine wave shown in Figure 15.1. On most systems, this command will open a separate plot window to display the graph.


Figure 15.1: Simple Two-Dimensional Plot.

```
plot (y)
plot (x, y)
plot (x, y, fmt)
plot (..., property, value,...)
plot (x1, y1, ..., xn, yn)
plot (hax,...)
h = plot (...)
Produce 2-D plots.
```

Many different combinations of arguments are possible. The simplest form is plot ( $y$ )
where the argument is taken as the set of $y$ coordinates and the $x$ coordinates are taken to be the range 1:numel ( $y$ ).
If more than one argument is given, they are interpreted as
plot (y, property, value, ...)
or
plot (x, y, property, value, ...)
or
plot (x, y, fmt, ...)
and so on. Any number of argument sets may appear. The $x$ and $y$ values are interpreted as follows:

- If a single data argument is supplied, it is taken as the set of $y$ coordinates and the $x$ coordinates are taken to be the indices of the elements, starting with 1 .
- If $x$ and $y$ are scalars, a single point is plotted.
- squeeze() is applied to arguments with more than two dimensions, but no more than two singleton dimensions.
- If both arguments are vectors, the elements of $y$ are plotted versus the elements of $x$.
- If $x$ is a vector and $y$ is a matrix, then the columns (or rows) of $y$ are plotted versus $x$. (using whichever combination matches, with columns tried first.)
- If the $x$ is a matrix and $y$ is a vector, $y$ is plotted versus the columns (or rows) of $x$. (using whichever combination matches, with columns tried first.)
- If both arguments are matrices, the columns of $y$ are plotted versus the columns of $x$. In this case, both matrices must have the same number of rows and columns and no attempt is made to transpose the arguments to make the number of rows match.

Multiple property-value pairs may be specified, but they must appear in pairs. These arguments are applied to the line objects drawn by plot. Useful properties to modify are "linestyle", "linewidth", "color", "marker", "markersize", "markeredgecolor", "markerfacecolor". The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450.
The fmt format argument can also be used to control the plot style. It is a string composed of four optional parts: "<linestyle><marker><color><;displayname;>". When a marker is specified, but no linestyle, only the markers are plotted. Similarly, if a linestyle is specified, but no marker, then only lines are drawn. If both are specified then lines and markers will be plotted. If no fmt and no property/value pairs are given, then the default plot style is solid lines with no markers and the color determined by the "colororder" property of the current axes.
Format arguments:
linestyle

$$
\begin{array}{ll}
\therefore- & \text { Use solid lines (default). } \\
\therefore--, & \text { Use dashed lines. } \\
\ddots: \prime & \text { Use dotted lines. } \\
\therefore-. & \text { Use dash-dotted lines. }
\end{array}
$$

marker

$$
\begin{aligned}
& \text { '+' crosshair } \\
& \text { 'o' circle } \\
& \text { '*' star } \\
& \text { '.' point } \\
& \text { ' } x \text { ' cross } \\
& \text { 'I' vertical line } \\
& \text { '_' horizontal line } \\
& \text { 's' square } \\
& \text { 'd' diamond } \\
& \text { (a) upward-facing triangle } \\
& \text { 'v' downward-facing triangle } \\
& \text { '>' right-facing triangle } \\
& \text { '<' left-facing triangle } \\
& \text { 'p' pentagram } \\
& \text { ' } \mathrm{h} \text { ' hexagram }
\end{aligned}
$$

color

| 'k', "black" | blacK |
| :--- | :--- |
| 'r', "red" | Red |
| 'g', "green" | Green |
| 'b', "blue" | Blue |
| 'y', "yellow" | Yellow |
| 'm', "magenta" | Magenta |
| 'c', "cyan" | Cyan |
| 'w', "white" | White |

"; displayname;"
The text between semicolons is used to set the "displayname" property which determines the label used for the plot legend.

The fmt argument may also be used to assign legend labels. To do so, include the desired label between semicolons after the formatting sequence described above, e.g., "+b;Data Series 3;". Note that the last semicolon is required and Octave will generate an error if it is left out.
Here are some plot examples:

```
plot (x, y, "or", x, y2, x, y3, "m", x, y4, "+")
```

This command will plot y with red circles, y2 with solid lines, y3 with solid magenta lines, and y4 with points displayed as ' + '.

```
plot (b, "*", "markersize", 10)
```

This command will plot the data in the variable b , with points displayed as ' $*$ ' and a marker size of 10 .

```
t = 0:0.1:6.3;
plot (t, cos(t), "-;cos(t);", t, sin(t), "-b;sin(t);");
```

This will plot the cosine and sine functions and label them accordingly in the legend. If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a vector of graphics handles to the created line objects.
To save a plot, in one of several image formats such as PostScript or PNG, use the print command.
See also: [axis], page 348, [box], page 394, [grid], page 394, [hold], page 403, [legend], page 390, [title], page 389, [xlabel], page 393, [ylabel], page 393, [xlim], page 351, [ylim], page 351, [ezplot], page 354, [errorbar], page 333, [fplot], page 353, [line], page 424, [plot3], page 374, [polar], page 337, [loglog], page 320, [semilogx], page 319, [semilogy], page 320, [subplot], page 398.

The plotyy function may be used to create a plot with two independent y axes.

```
plotyy (x1, y1, x2, y2)
plotyy (..., fun)
plotyy (..., fun1, fun2)
plotyy (hax, ...)
```

[ax, h1, h2] = plotyy (...)
Plot two sets of data with independent y -axes and a common x -axis.
The arguments $x 1$ and $y 1$ define the arguments for the first plot and $x 1$ and $y 2$ for the second.
By default the arguments are evaluated with feval (@plot, x, y). However the type of plot can be modified with the fun argument, in which case the plots are generated by feval (fun, $x, y$ ). fun can be a function handle, an inline function, or a string of a function name.
The function to use for each of the plots can be independently defined with fun1 and fun2.

The first argument hax can be an axes handle to the principal axes in which to plot the $x 1$ and $y 1$ data. It can also be a two-element vector with the axes handles to the primary and secondary axes (see output ax).
The return value ax is a vector with the axes handles of the two $y$-axes. $h 1$ and $h 2$ are handles to the objects generated by the plot commands.

```
x = 0:0.1:2*pi;
y1 = sin (x);
y2 = exp (x - 1);
ax = plotyy (x, y1, x - 1, y2, @plot, @semilogy);
xlabel ("X");
ylabel (ax(1), "Axis 1");
ylabel (ax(2), "Axis 2");
```

When using plotyy in conjunction with subplot make sure to call subplot first and pass the resulting axes handle to plotyy. Do not call subplot with any of the axes handles returned by plotyy or the other axes will be removed.

See also: [plot], page 316, [subplot], page 398.
The functions semilogx, semilogy, and loglog are similar to the plot function, but produce plots in which one or both of the axes use log scales.

```
semilogx (y)
semilogx (x, y)
semilogx (x, y, property, value,...)
semilogx (x, y, fmt)
semilogx (hax,...)
h = semilogx (...)
Produce a 2-D plot using a logarithmic scale for the x -axis.
```

See the documentation of plot for a description of the arguments that semilogx will accept.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.
See also: [plot], page 316, [semilogy], page 320, [loglog], page 320.

```
semilogy ( \(y\) )
semilogy ( \(x, y\) )
semilogy ( \(x, y\), property, value, ...)
semilogy ( \(x, y, f m t\) )
semilogy (h, ...)
\(h=\) semilogy (...)
```

    Produce a 2-D plot using a logarithmic scale for the y-axis.
    See the documentation of plot for a description of the arguments that semilogy will accept.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a graphics handle to the created plot.
See also: [plot], page 316, [semilogx], page 319, [loglog], page 320.

```
loglog (y)
loglog (x, y)
loglog (x, y, prop, value,...)
loglog (x, y, fmt)
loglog (hax,...)
h = loglog (...)
Produce a 2-D plot using logarithmic scales for both axes.
```

See the documentation of plot for a description of the arguments that loglog will accept.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a graphics handle to the created plot.
See also: [plot], page 316, [semilogx], page 319, [semilogy], page 320.

The functions bar, barh, stairs, and stem are useful for displaying discrete data. For example,

```
randn ("state", 1);
hist (randn (10000, 1), 30);
xlabel ("Value");
ylabel ("Count");
title ("Histogram of 10,000 normally distributed random numbers");
```

produces the histogram of 10,000 normally distributed random numbers shown in Figure 15.2. Note that, randn ("state", 1); initializes the random number generator for randn to a known value so that the returned values are reproducible; This guarantees that the figure produced is identical to the one in this manual.

Histogram of 10,000 normally distributed random numbers


Figure 15.2: Histogram.
bar (y)
bar (x, y)
bar (..., w)
bar (..., style)
bar (..., prop, val, ...)
bar (hax, ...)
h = bar (..., prop, val, ...)
Produce a bar graph from two vectors of X-Y data.
If only one argument is given, $y$, it is taken as a vector of Y values and the X coordinates are the range 1 : numel ( $y$ ).
The optional input $w$ controls the width of the bars. A value of 1.0 will cause each bar to exactly touch any adjacent bars. The default width is 0.8 .
If $y$ is a matrix, then each column of $y$ is taken to be a separate bar graph plotted on the same graph. By default the columns are plotted side-by-side. This behavior can be changed by the style argument which can take the following values:
"grouped" (default)
Side-by-side bars with a gap between bars and centered over the Xcoordinate.
"stacked"
Bars are stacked so that each X value has a single bar composed of multiple segments.
"hist" Side-by-side bars with no gap between bars and centered over the Xcoordinate.
"histc" Side-by-side bars with no gap between bars and left-aligned to the Xcoordinate.

Optional property/value pairs are passed directly to the underlying patch objects. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of handles to the created "bar series" hggroups with one handle per column of the variable $y$. This series makes it possible to change a common element in one bar series object and have the change reflected in the other "bar series". For example,

```
h = bar (rand (5, 10));
set (h(1), "basevalue", 0.5);
```

changes the position on the base of all of the bar series.
The following example modifies the face and edge colors using property/value pairs.

```
bar (randn (1, 100), "facecolor", "r", "edgecolor", "b");
```

The default color for bars is taken from the axes' "ColorOrder" property. The default color for bars when a histogram option ("hist", "histc" is used is the "Colormap" property of either the axes or figure. The color of bars can also be set manually using the "facecolor" property as shown below.

```
h = bar (rand (10, 3));
set (h(1), "facecolor", "r")
set (h(2), "facecolor", "g")
set (h(3), "facecolor", "b")
```

See also: [barh], page 322, [hist], page 323, [pie], page 338, [plot], page 316, [patch], page 425 .
barh (y)
barh ( $x, y$ )
barh (..., w)
barh (..., style)
barh (..., prop, val, ...)
barh (hax, ...)
$h=\operatorname{barh}(\ldots$, prop, val, ...)
Produce a horizontal bar graph from two vectors of X-Y data.
If only one argument is given, it is taken as a vector of Y values and the X coordinates are the range 1 :numel ( $y$ ).
The optional input $w$ controls the width of the bars. A value of 1.0 will cause each bar to exactly touch any adjacent bars. The default width is 0.8 .
If $y$ is a matrix, then each column of $y$ is taken to be a separate bar graph plotted on the same graph. By default the columns are plotted side-by-side. This behavior can be changed by the style argument which can take the following values:
"grouped" (default)
Side-by-side bars with a gap between bars and centered over the Ycoordinate.
"stacked"
Bars are stacked so that each Y value has a single bar composed of multiple segments.
"hist" Side-by-side bars with no gap between bars and centered over the Ycoordinate.
"histc" Side-by-side bars with no gap between bars and left-aligned to the Ycoordinate.

Optional property/value pairs are passed directly to the underlying patch objects. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created bar series hggroup. For a description of the use of the bar series, see [bar], page 321.
See also: [bar], page 321, [hist], page 323, [pie], page 338, [plot], page 316, [patch], page 425.
hist (y)
hist (y, nbins)
hist ( $y, x$ )
hist (y, x, norm)
hist (..., prop, val, ...)
hist (hax, ...)
[nn, xx] = hist (...)
Produce histogram counts or plots.
With one vector input argument, $y$, plot a histogram of the values with 10 bins. The range of the histogram bins is determined by the range of the data (difference between maximum and minimum value in $y$ ). Extreme values are lumped into the first and last bins. If $y$ is a matrix then plot a histogram where each bin contains one bar per input column of $y$.
If the optional second argument is a scalar, nbins, it defines the number of bins.
If the optional second argument is a vector, $x$, it defines the centers of the bins. The width of the bins is determined from the adjacent values in the vector. The total number of bins is numel ( $x$ ).

If a third argument norm is provided, the histogram is normalized. In case norm is a positive scalar, the resulting bars are normalized to norm. If norm is a vector of positive scalars of length columns ( $y$ ), then the resulting bar of $y(:, i)$ is normalized to norm(i).

```
[nn, xx] = hist (rand (10, 3), 5, [1 2 3]);
sum (nn, 1)
a ans =
```

    123
    The histogram's appearance may be modified by specifying property/value pairs to the underlying patch object. For example, the face and edge color may be modified:
hist (randn (1, 100), 25, "facecolor", "r", "edgecolor", "b");
The full list of patch properties is documented at Section 15.3.3.8 [Patch Properties], page 457. property. If not specified, the default colors for the histogram are taken from the "Colormap" property of the axes or figure.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If an output is requested then no plot is made. Instead, return the values nn (numbers of elements) and $x x$ (bin centers) such that bar ( $x x, n n$ ) will plot the histogram.

See also: [histc], page 784, [bar], page 321, [pie], page 338, [rose], page 329.
stemleaf (x, caption)
stemleaf (x, caption, stem_sz)
plotstr $=$ stemleaf (...)
Compute and display a stem and leaf plot of the vector $x$.
The input $x$ should be a vector of integers. Any non-integer values will be converted to integer by $x=f i x(x)$. By default each element of $x$ will be plotted with the last digit of the element as a leaf value and the remaining digits as the stem. For example, 123 will be plotted with the stem ' 12 ' and the leaf ' 3 '. The second argument, caption, should be a character array which provides a description of the data. It is included as a heading for the output.
The optional input stem_sz sets the width of each stem. The stem width is determined by $10^{\wedge}$ (stem_sz+1). The default stem width is 10 .
The output of stemleaf is composed of two parts: a "Fenced Letter Display," followed by the stem-and-leaf plot itself. The Fenced Letter Display is described in Exploratory Data Analysis. Briefly, the entries are as shown:


The stem-and-leaf plot shows on each line the stem value followed by the string made up of the leaf digits. If the stem_sz is not 1 the successive leaf values are separated by ",".
With no return argument, the plot is immediately displayed. If an output argument is provided, the plot is returned as an array of strings.
The leaf digits are not sorted. If sorted leaf values are desired, use $x$ s $=\operatorname{sort}(x)$ before calling stemleaf (xs).

The stem and leaf plot and associated displays are described in: Chapter 3, Exploratory Data Analysis by J. W. Tukey, Addison-Wesley, 1977.
See also: [hist], page 323, [printd], page 325.
printd (obj, filename)
out_file = printd (...)
Convert any object acceptable to disp into the format selected by the suffix of filename.
If the return argument out_file is given, the name of the created file is returned.
This function is intended to facilitate manipulation of the output of functions such as stemleaf.

See also: [stemleaf], page 324.

```
stairs (y)
```

stairs ( $x, y$ )
stairs (..., style)
stairs (..., prop, val, ...)
stairs (hax,...)
h = stairs (...)
[xstep, ystep] = stairs (...)

Produce a stairstep plot.
The arguments $x$ and $y$ may be vectors or matrices. If only one argument is given, it is taken as a vector of Y values and the X coordinates are taken to be the indices of the elements ( $x=1$ : numel $(y)$ ).
The style to use for the plot can be defined with a line style style of the same format as the plot command.
Multiple property/value pairs may be specified, but they must appear in pairs. The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If one output argument is requested, return a graphics handle to the created plot. If two output arguments are specified, the data are generated but not plotted. For example,

```
stairs (x, y);
```

and

```
[xs, ys] = stairs (x, y);
plot (xs, ys);
```

are equivalent.
See also: [bar], page 321, [hist], page 323, [plot], page 316, [stem], page 325.
stem (y)
stem ( $x, y$ )
stem (..., linespec)
stem (..., "filled")

```
stem (..., prop, val, ...)
stem (hax,...)
h = stem (...)
Plot a 2-D stem graph.
```

If only one argument is given, it is taken as the $y$-values and the $x$-coordinates are taken from the indices of the elements.
If $y$ is a matrix, then each column of the matrix is plotted as a separate stem graph. In this case $x$ can either be a vector, the same length as the number of rows in $y$, or it can be a matrix of the same size as $y$.
The default color is "b" (blue), the default line style is "-", and the default marker is "o". The line style can be altered by the linespec argument in the same manner as the plot command. If the "filled" argument is present the markers at the top of the stems will be filled in. For example,

```
x = 1:10;
y = 2*x;
stem (x, y, "r");
```

plots 10 stems with heights from 2 to 20 in red;
Optional property/value pairs may be specified to control the appearance of the plot. The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a handle to a "stem series" hggroup. The single hggroup handle has all of the graphical elements comprising the plot as its children; This allows the properties of multiple graphics objects to be changed by modifying just a single property of the "stem series" hggroup.
For example,

```
x = [0:10]';
y = [sin(x), cos(x)]
h = stem (x, y);
set (h(2), "color", "g");
set (h(1), "basevalue", -1)
```

changes the color of the second "stem series" and moves the base line of the first.
Stem Series Properties
linestyle The linestyle of the stem. (Default: "-")
linewidth The width of the stem. (Default: 0.5)
color The color of the stem, and if not separately specified, the marker. (Default: "b" [blue])
marker The marker symbol to use at the top of each stem. (Default: "०")
markeredgecolor
The edge color of the marker. (Default: "color" property)
markerfacecolor
The color to use for "filling" the marker. (Default: "none" [unfilled])
markersize
The size of the marker. (Default: 6)
baseline The handle of the line object which implements the baseline. Use set with the returned handle to change graphic properties of the baseline.
basevalue The y -value where the baseline is drawn. (Default: 0 )
See also: [stem3], page 327, [bar], page 321, [hist], page 323, [plot], page 316, [stairs], page 325 .

```
stem3 (x, y, z)
stem3 (..., linespec)
stem3 (..., "filled")
stem3 (..., prop, val, ...)
stem3 (hax,...)
h = stem3 (...)
Plot a 3-D stem graph.
```

Stems are drawn from the height $z$ to the location in the $x-y$ plane determined by $x$ and $y$. The default color is "b" (blue), the default line style is " - ", and the default marker is "०".
The line style can be altered by the linespec argument in the same manner as the plot command. If the "filled" argument is present the markers at the top of the stems will be filled in.
Optional property/value pairs may be specified to control the appearance of the plot. The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450. If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a handle to the "stem series" hggroup containing the line and marker objects used for the plot. See [stem], page 325, for a description of the "stem series" object.
Example:

```
theta = 0:0.2:6;
stem3 (cos (theta), sin (theta), theta);
```

plots 31 stems with heights from 0 to 6 lying on a circle.
Implementation Note: Color definitions with RGB-triples are not valid.
See also: [stem], page 325, [bar], page 321, [hist], page 323, [plot], page 316.

```
scatter (x, y)
```

scatter ( $x, y, s$ )
scatter ( $x, y, s, c$ )
scatter (..., style)
scatter (..., "filled")
scatter (..., prop, val, ...)
scatter (hax, ...)
$h=\operatorname{scatter}(\ldots)$

Draw a 2-D scatter plot.

A marker is plotted at each point defined by the coordinates in the vectors $x$ and $y$. The size of the markers is determined by $s$, which can be a scalar or a vector of the same length as $x$ and $y$. If $s$ is not given, or is an empty matrix, then a default value of 36 square points is used (The marker size itself is sqrt (s)).
The color of the markers is determined by $c$, which can be a string defining a fixed color; a 3 -element vector giving the red, green, and blue components of the color; a vector of the same length as $x$ that gives a scaled index into the current colormap; or an Nx3 matrix defining the RGB color of each marker individually.
The marker to use can be changed with the style argument; it is a string defining a marker in the same manner as the plot command. If no marker is specified it defaults to "o" or circles. If the argument "filled" is given then the markers are filled.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created scatter object.
Example:

```
x = randn (100, 1);
y = randn (100, 1);
scatter (x, y, [], sqrt (x.^2 + y.^2));
```

Programming Note: The full list of properties is documented at Section 15.3.3.9 [Scatter Properties], page 461.
See also: [scatter3], page 382, [patch], page 425, [plot], page 316.
plotmatrix ( $x, y$ )
plotmatrix (x)
plotmatrix (..., style)
plotmatrix (hax, ...)
[h, ax, bigax, p, pax] = plotmatrix (...)
Scatter plot of the columns of one matrix against another.
Given the arguments $x$ and $y$ that have a matching number of rows, plotmatrix plots a set of axes corresponding to

```
plot (x(:, i), y(:, j))
```

When called with a single argument $x$ this is equivalent to
plotmatrix ( $x, x$ )
except that the diagonal of the set of axes will be replaced with the histogram hist ( $x(:, i)$ ).
The marker to use can be changed with the style argument, that is a string defining a marker in the same manner as the plot command.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ provides handles to the individual graphics objects in the scatter plots, whereas ax returns the handles to the scatter plot axes objects.
bigax is a hidden axes object that surrounds the other axes, such that the commands xlabel, title, etc., will be associated with this hidden axes.

Finally, $p$ returns the graphics objects associated with the histogram and pax the corresponding axes objects.
Example:

```
plotmatrix (randn (100, 3), "g+")
```

See also: [scatter], page 327, [plot], page 316.

```
pareto (y)
pareto (y, x)
pareto (hax,...)
h = pareto (...)
    Draw a Pareto chart.
```

A Pareto chart is a bar graph that arranges information in such a way that priorities for process improvement can be established; It organizes and displays information to show the relative importance of data. The chart is similar to the histogram or bar chart, except that the bars are arranged in decreasing magnitude from left to right along the x -axis.
The fundamental idea (Pareto principle) behind the use of Pareto diagrams is that the majority of an effect is due to a small subset of the causes. For quality improvement, the first few contributing causes (leftmost bars as presented on the diagram) to a problem usually account for the majority of the result. Thus, targeting these "major causes" for elimination results in the most cost-effective improvement scheme.
Typically only the magnitude data $y$ is present in which case $x$ is taken to be the range 1 : length ( $y$ ). If $x$ is given it may be a string array, a cell array of strings, or a numerical vector.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a 2 -element vector with a graphics handle for the created bar plot and a second handle for the created line plot.
An example of the use of pareto is

```
Cheese = {"Cheddar", "Swiss", "Camembert", ...
    "Munster", "Stilton", "Blue"};
Sold = [105, 30, 70, 10, 15, 20];
pareto (Sold, Cheese);
```

See also: [bar], page 321, [barh], page 322, [hist], page 323, [pie], page 338, [plot], page 316.
rose (th)
rose (th, nbins)
rose (th, bins)
rose (hax, ...)
$h=$ rose (...)
[thout rout] = rose (...)
Plot an angular histogram.
With one vector argument, th, plot the histogram with 20 angular bins. If th is a matrix then each column of $t h$ produces a separate histogram.

If nbins is given and is a scalar, then the histogram is produced with nbin bins. If bins is a vector, then the center of each bin is defined by the values in bins and the number of bins is given by the number of elements in bins.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the line objects representing each histogram.
If two output arguments are requested then no plot is made and the polar vectors necessary to plot the histogram are returned instead.
Example

```
[th, r] = rose ([2*randn(1e5,1), pi + 2*randn(1e5,1)]);
```

polar (th, r);

Programming Note: When specifying bin centers with the bins input, the edges for bins 2 to $\mathrm{N}-1$ are spaced so that bins(i) is centered between the edges. The final edge is drawn halfway between bin N and bin 1. This guarantees that all input th will be placed into one of the bins, but also means that for some combinations bin 1 and bin N may not be centered on the user's given values.
See also: [hist], page 323, [polar], page 337.
The contour, contourf and contourc functions produce two-dimensional contour plots from three-dimensional data.

```
contour (z)
contour (z,vn)
contour (x, y, z)
contour (x, y, z, vn)
contour (..., style)
contour (hax,...)
[c,h] = contour (...)
```

Create a 2-D contour plot.
Plot level curves (contour lines) of the matrix z, using the contour matrix $c$ computed by contourc from the same arguments; see the latter for their interpretation.
The appearance of contour lines can be defined with a line style style in the same manner as plot. Only line style and color are used; Any markers defined by style are ignored.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional output contains the contour levels in contourc format.
The optional return value $h$ is a graphics handle to the hggroup comprising the contour lines.
Example:

```
x = 0:3;
y = 0:2;
z = y' * x;
contour (x, y, z, 2:3)
```

See also: [ezcontour], page 355, [contourc], page 331, [contourf], page 331, [contour3], page 332, [clabel], page 393, [meshc], page 359, [surfc], page 361, [caxis], page 350, [colormap], page 870, [plot], page 316.

```
contourf (z)
contourf (z, vn)
contourf (x, y, z)
contourf (x, y, z, vn)
contourf (..., style)
contourf (hax,...)
[c,h] = contourf (...)
```

Create a 2-D contour plot with filled intervals.
Plot level curves (contour lines) of the matrix $z$ and fill the region between lines with colors from the current colormap.
The level curves are taken from the contour matrix $c$ computed by contourc for the same arguments; see the latter for their interpretation.

The appearance of contour lines can be defined with a line style style in the same manner as plot. Only line style and color are used; Any markers defined by style are ignored.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional output contains the contour levels in contourc format.
The optional return value $h$ is a graphics handle to the hggroup comprising the contour lines.
The following example plots filled contours of the peaks function.

```
[x, y, z] = peaks (50);
contourf (x, y, z, -7:9)
```

See also: [ezcontourf], page 355, [contour], page 330, [contourc], page 331, [contour3], page 332, [clabel], page 393, [meshc], page 359, [surfc], page 361, [caxis], page 350, [colormap], page 870, [plot], page 316.
$c=$ contourc (z)
$c=$ contourc ( $z, v n$ )
$c=$ contourc ( $x, y, z$ )
$c=$ contourc ( $x, y, z, v n$ )
[c, lev] = contourc (...)
Compute contour lines (isolines of constant Z value).
The matrix $z$ contains height values above the rectangular grid determined by $x$ and $y$. If only a single input $z$ is provided then $x$ is taken to be 1 :columns ( $z$ ) and $y$ is taken to be 1 :rows ( $z$ ). The minimum data size is 2 x 2 .
The optional input vn is either a scalar denoting the number of contour lines to compute or a vector containing the Z values where lines will be computed. When vn is a vector the number of contour lines is numel (vn). However, to compute a single contour line at a given value use $v n=[v a l, v a l]$. If $v n$ is omitted it defaults to 10 .

The return value $c$ is a 2 xn matrix containing the contour lines in the following format

```
c = [lev1, x1, x2, ..., levn, x1, x2, ...
    len1, y1, y2, ..., lenn, y1, y2, ...]
```

in which contour line $n$ has a level (height) of levn and length of lenn.
The optional return value lev is a vector with the Z values of the contour levels.
Example:

```
x = 0:2;
y = x;
z = x' * y;
c = contourc (x, y, z, 2:3)
    c =
\begin{tabular}{llllllll}
2.0000 & 1.0000 & 1.0000 & 2.0000 & 2.0000 & 3.0000 & 1.5000 & 2.0000 \\
4.0000 & 2.0000 & 2.0000 & 1.0000 & 1.0000 & 2.0000 & 2.0000 & 1.5000
\end{tabular}
```

See also: [contour], page 330, [contourf], page 331, [contour3], page 332, [clabel], page 393 .

```
contour3 (z)
contour3 (z, vn)
contour3 (x, y, z)
contour3 (x, y, z, vn)
contour3 (..., style)
contour3 (hax,...)
[c, h] = contour3 (...)
```

Create a 3-D contour plot.
contour3 plots level curves (contour lines) of the matrix $z$ at a Z level corresponding to each contour. This is in contrast to contour which plots all of the contour lines at the same Z level and produces a 2-D plot.

The level curves are taken from the contour matrix $c$ computed by contourc for the same arguments; see the latter for their interpretation.

The appearance of contour lines can be defined with a line style style in the same manner as plot. Only line style and color are used; Any markers defined by style are ignored.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional output $c$ are the contour levels in contourc format.
The optional return value $h$ is a graphics handle to the hggroup comprising the contour lines.

Example:

```
contour3 (peaks (19));
colormap cool;
hold on;
surf (peaks (19), "facecolor", "none", "edgecolor", "black");
```

See also: [contour], page 330, [contourc], page 331, [contourf], page 331, [clabel], page 393, [meshc], page 359, [surfc], page 361, [caxis], page 350, [colormap], page 870, [plot], page 316.

The errorbar, semilogxerr, semilogyerr, and loglogerr functions produce plots with error bar markers. For example,

```
rand ("state", 2);
\(\mathrm{x}=0: 0.1: 10 ;\)
\(y=\sin (x)\);
lerr \(=0.1\).* rand (size (x));
uerr \(=0.1\).* rand (size (x));
errorbar (x, y, lerr, uerr);
axis ([0, 10, -1.1, 1.1]);
xlabel ("x");
ylabel ("sin (x)");
title ("Errorbar plot of \(\sin (x) ")\);
```

produces the figure shown in Figure 15.3.


Figure 15.3: Errorbar plot.

```
errorbar (y, ey)
errorbar (y,...,fmt)
errorbar (x, y, ey)
errorbar (x, y, err, fmt)
errorbar (x, y, lerr, uerr, fmt)
errorbar (x, y, ex, ey, fmt)
errorbar (x, y, lx, ux, ly, uy, fmt)
errorbar (x1, y1,..., fmt, xn, yn,...)
errorbar (hax,...)
```

$h=\operatorname{errorbar}(. .$.
Create a $2-\mathrm{D}$ plot with errorbars.
Many different combinations of arguments are possible. The simplest form is
errorbar ( $y$, ey)
where the first argument is taken as the set of $y$ coordinates, the second argument ey are the errors around the $y$ values, and the $x$ coordinates are taken to be the indices of the elements (1:numel (y)).
The general form of the function is
errorbar (x, y, err1, ..., fmt, ...)
After the $x$ and $y$ arguments there can be 1,2 , or 4 parameters specifying the error values depending on the nature of the error values and the plot format fmt.
err (scalar)
When the error is a scalar all points share the same error value. The errorbars are symmetric and are drawn from data-err to data+err. The fimt argument determines whether err is in the x -direction, y -direction (default), or both.
err (vector or matrix)
Each data point has a particular error value. The errorbars are symmetric and are drawn from $\operatorname{data}(\mathrm{n})-\operatorname{err}(\mathrm{n})$ to $\operatorname{data}(\mathrm{n})+\operatorname{err}(\mathrm{n})$.
lerr, uerr (scalar)
The errors have a single low-side value and a single upper-side value. The errorbars are not symmetric and are drawn from data-lerr to data+uerr.
lerr, uerr (vector or matrix)
Each data point has a low-side error and an upper-side error. The errorbars are not symmetric and are drawn from data(n)-lerr(n) to $\operatorname{data}(\mathrm{n})+\operatorname{uerr}(\mathrm{n})$.

Any number of data sets $(x 1, y 1, x 2, y 2, \ldots)$ may appear as long as they are separated by a format string fmt.
If $y$ is a matrix, $x$ and the error parameters must also be matrices having the same dimensions. The columns of $y$ are plotted versus the corresponding columns of $x$ and errorbars are taken from the corresponding columns of the error parameters.
If fmt is missing, the yerrorbars ("~") plot style is assumed.
If the fmt argument is supplied then it is interpreted, as in normal plots, to specify the line style, marker, and color. In addition, fmt may include an errorbar style which must precede the ordinary format codes. The following errorbar styles are supported:

$$
\begin{array}{ll}
‘ \sim, & \text { Set yerrorbars plot style (default). } \\
‘>, & \text { Set xerrorbars plot style. } \\
‘ \sim>, & \text { Set xyerrorbars plot style. } \\
‘ \#^{\sim}, & \text { Set yboxes plot style. } \\
‘ \# ' & \text { Set xboxes plot style. }
\end{array}
$$

'\#~>' Set xyboxes plot style.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a handle to the hggroup object representing the data plot and errorbars.
Note: For compatibility with matlab a line is drawn through all data points. However, most scientific errorbar plots are a scatter plot of points with errorbars. To accomplish this, add a marker style to the fmt argument such as ". ". Alternatively, remove the line by modifying the returned graphic handle with set (h, "linestyle", "none").
Examples:

```
errorbar (x, y, ex, ">.r")
```

produces an xerrorbar plot of $y$ versus $x$ with $x$ errorbars drawn from $x$-ex to $x+e x$. The marker "." is used so no connecting line is drawn and the errorbars appear in red.

```
errorbar (x, y1, ey, "~",
    x, y2, ly, uy)
```

produces yerrorbar plots with y1 and y2 versus x. Errorbars for y1 are drawn from y1-ey to $y 1+e y$, errorbars for $y 2$ from y2-ly to $y 2+u y$.

```
errorbar (x, y, lx, ux,
    ly, uy, "~>")
```

produces an xyerrorbar plot of $y$ versus $x$ in which $x$ errorbars are drawn from $x$-lx to $x+u x$ and $y$ errorbars from $y$-ly to $y+u y$.
See also: [semilogxerr], page 335, [semilogyerr], page 336, [loglogerr], page 336, [plot], page 316.

```
semilogxerr (y, ey)
semilogxerr (y, .., fmt)
semilogxerr (x, y, ey)
semilogxerr (x, y, err, fmt)
semilogxerr (x, y, lerr, uerr, fmt)
semilogxerr (x, y, ex, ey, fmt)
semilogxerr (x, y, lx, ux, ly, uy, fmt)
semilogxerr (x1, y1,..., fmt, xn, yn,...)
semilogxerr (hax,...)
h = semilogxerr (...)
```

Produce 2-D plots using a logarithmic scale for the x -axis and errorbars at each data point.
Many different combinations of arguments are possible. The most common form is

```
semilogxerr (x, y, ey, fmt)
```

which produces a semi-logarithmic plot of $y$ versus $x$ with errors in the $y$-scale defined by ey and the plot format defined by fmt. See [errorbar], page 333, for available formats and additional information.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [errorbar], page 333, [semilogyerr], page 336, [loglogerr], page 336.

```
semilogyerr (y, ey)
semilogyerr (y, .., fmt)
semilogyerr (x, y, ey)
semilogyerr (x, y, err, fmt)
semilogyerr (x, y, lerr, uerr, fmt)
semilogyerr (x, y, ex, ey, fmt)
semilogyerr (x, y, lx, ux, ly, uy, fmt)
semilogyerr (x1, y1, .., fmt, xn, yn,...)
semilogyerr (hax,...)
h = semilogyerr (...)
```

Produce 2-D plots using a logarithmic scale for the $y$-axis and errorbars at each data point.
Many different combinations of arguments are possible. The most common form is

```
semilogyerr (x, y, ey, fmt)
```

which produces a semi-logarithmic plot of $y$ versus $x$ with errors in the $y$-scale defined by ey and the plot format defined by fmt. See [errorbar], page 333, for available formats and additional information.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [errorbar], page 333, [semilogxerr], page 335, [loglogerr], page 336.

```
loglogerr (y, ey)
loglogerr (y,...,fmt)
loglogerr (x, y, ey)
loglogerr (x, y, err, fmt)
loglogerr (x, y, lerr, uerr, fmt)
loglogerr (x, y, ex, ey, fmt)
loglogerr (x, y, lx, ux, ly, uy, fmt)
loglogerr (x1, y1, ..., fmt, xn, yn,...)
loglogerr (hax,...)
h = loglogerr (...)
```

Produce 2-D plots on a double logarithm axis with errorbars.
Many different combinations of arguments are possible. The most common form is

```
loglogerr (x, y, ey, fmt)
```

which produces a double logarithm plot of $y$ versus $x$ with errors in the $y$-scale defined by ey and the plot format defined by fmt. See [errorbar], page 333, for available formats and additional information.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [errorbar], page 333, [semilogxerr], page 335, [semilogyerr], page 336.

Finally, the polar function allows you to easily plot data in polar coordinates. However, the display coordinates remain rectangular and linear. For example,
polar ( $0: 0.1: 10 * \mathrm{pi}, 0: 0.1: 10 * \mathrm{pi})$;
title ("Example polar plot from 0 to $10 *$ pi");
produces the spiral plot shown in Figure 15.4.


Figure 15.4: Polar plot.

```
polar (theta, rho)
polar (theta, rho, fmt)
polar (cplx)
polar (cplx, fmt)
polar (hax, ...)
h = polar (...)
Create a 2-D plot from polar coordinates theta and rho.
```

The input theta is assumed to be radians and is converted to degrees for plotting. If you have degrees then you must convert (see [cart2pol], page 578) to radians before passing the data to this function.
If a single complex input cplx is given then the real part is used for theta and the imaginary part is used for rho.
The optional argument fmt specifies the line format in the same way as plot.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.
Implementation Note: The polar axis is drawn using line and text objects encapsulated in an hggroup. The hggroup properties are linked to the original axes object such that altering an appearance property, for example fontname, will update the polar axis. Two new properties are added to the original axes-rtick, ttick-which
replace xtick, ytick. The first is a list of tick locations in the radial (rho) direction; The second is a list of tick locations in the angular (theta) direction specified in degrees, i.e., in the range 0-359.

See also: [rose], page 329, [compass], page 344, [plot], page 316, [cart2pol], page 578.

```
pie (x)
pie (..., explode)
pie (..., labels)
pie (hax,...)
h = pie (...)
```

Plot a 2-D pie chart.
When called with a single vector argument, produce a pie chart of the elements in $x$. The size of the ith slice is the percentage that the element xi represents of the total sum of $x$ : pct $=x(i) /$ sum ( $x$ ).
The optional input explode is a vector of the same length as $x$ that, if nonzero, "explodes" the slice from the pie chart.
The optional input labels is a cell array of strings of the same length as $x$ specifying the label for each slice.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a list of handles to the patch and text objects generating the plot.
Note: If sum ( $x$ ) $\leq 1$ then the elements of $x$ are interpreted as percentages directly and are not normalized by sum (x). Furthermore, if the sum is less than 1 then there will be a missing slice in the pie plot to represent the missing, unspecified percentage.
See also: [pie3], page 338, [bar], page 321, [hist], page 323, [rose], page 329.

```
pie3 (x)
pie3 (..., explode)
pie3 (..., labels)
pie3 (hax,...)
h = pie3 (...)
Plot a 3-D pie chart.
```

Called with a single vector argument, produces a 3-D pie chart of the elements in $x$. The size of the ith slice is the percentage that the element xi represents of the total sum of $x$ : pct $=x(i) / \operatorname{sum}(x)$.
The optional input explode is a vector of the same length as $x$ that, if nonzero, "explodes" the slice from the pie chart.
The optional input labels is a cell array of strings of the same length as $x$ specifying the label for each slice.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a list of graphics handles to the patch, surface, and text objects generating the plot.

Note: If sum $(x) \leq 1$ then the elements of $x$ are interpreted as percentages directly and are not normalized by sum ( x ). Furthermore, if the sum is less than 1 then there will be a missing slice in the pie plot to represent the missing, unspecified percentage.
See also: [pie], page 338, [bar], page 321, [hist], page 323, [rose], page 329.

```
quiver (u, v)
quiver (x, y, u,v)
quiver (..., s)
quiver (..., style)
quiver (..., "filled")
quiver (hax, ...)
h = quiver (...)
```

Plot a 2-D vector field with arrows.
Plot the ( $u, v$ ) components of a vector field at the grid points defined by $(x, y)$. If the grid is uniform then $x$ and $y$ can be specified as vectors and meshgrid is used to create the 2-D grid.
If $x$ and $y$ are not given they are assumed to be ( $1: m, 1: n$ ) where $[m, n]=$ size (u).

The optional input $s$ is a scalar defining a scaling factor to use for the arrows of the field relative to the mesh spacing. A value of 1.0 will result in the longest vector exactly filling one grid square. A value of 0 disables all scaling. The default value is 0.9 .

The style to use for the plot can be defined with a line style style of the same format as the plot command. If a marker is specified then the markers are drawn at the origin of the vectors (which are the grid points defined by $x$ and $y$ ). When a marker is specified, the arrowhead is not drawn. If the argument "filled" is given then the markers are filled.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to a quiver object. A quiver object regroups the components of the quiver plot (body, arrow, and marker), and allows them to be changed together.
Example:

```
[x, y] = meshgrid (1:2:20);
h = quiver (x, y, sin (2*pi*x/10), sin (2*pi*y/10));
set (h, "maxheadsize", 0.33);
```

See also: [quiver3], page 339, [compass], page 344, [feather], page 345, [plot], page 316.

```
quiver3 (x, y, z, u,v, w)
quiver3 (z, u, v, w)
quiver3 (..., s)
quiver3 (..., style)
quiver3 (..., "filled")
quiver3 (hax,...)
```

$h=$ quiver3 (...)
Plot a 3-D vector field with arrows.
Plot the $(u, v, w)$ components of a vector field at the grid points defined by $(x, y$, $z$ ). If the grid is uniform then $x, y$, and $z$ can be specified as vectors and meshgrid is used to create the $3-\mathrm{D}$ grid.
If $x$ and $y$ are not given they are assumed to be $(1: m, 1: n)$ where $[m, n]=$ size (u).

The optional input $s$ is a scalar defining a scaling factor to use for the arrows of the field relative to the mesh spacing. A value of 1.0 will result in the longest vector exactly filling one grid cube. A value of 0 disables all scaling. The default value is 0.9 .

The style to use for the plot can be defined with a line style style of the same format as the plot command. If a marker is specified then the markers are drawn at the origin of the vectors (which are the grid points defined by $x, y, z$ ). When a marker is specified, the arrowhead is not drawn. If the argument "filled" is given then the markers are filled.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to a quiver object. A quiver object regroups the components of the quiver plot (body, arrow, and marker), and allows them to be changed together.

```
[x, y, z] = peaks (25);
surf (x, y, z);
hold on;
[u, v, w] = surfnorm (x, y, z / 10);
h = quiver3 (x, y, z, u, v, w);
set (h, "maxheadsize", 0.33);
```

See also: [quiver], page 339, [compass], page 344, [feather], page 345, [plot], page 316.

```
streamribbon (x, y, z, u, v, w, sx, sy, sz)
streamribbon (u, v, w, sx, sy, sz)
streamribbon (xyz, x, y, z, anlr_spd, lin_spd)
streamribbon (xyz, anlr_spd, lin_spd)
streamribbon (xyz, anlr_rot)
streamribbon (..., width)
streamribbon (hax,...)
h = streamribbon (...)
```

Calculate and display streamribbons.
The streamribbon is constructed by rotating a normal vector around a streamline according to the angular rotation of the vector field.
The vector field is given by [ $u, v, w$ ] and is defined over a rectangular grid given by [ $x, y, z]$. The streamribbons start at the seed points [sx, sy, sz].
streamribbon can be called with a cell array that contains pre-computed streamline data. To do this, xyz must be created with the stream3 function. lin_spd is the linear speed of the vector field and can be calculated from $[u, v, w]$ by the square root of
the sum of the squares. The angular speed anlr_spd is the projection of the angular velocity onto the velocity of the normalized vector field and can be calculated with the curl command. This option is useful if you need to alter the integrator step size or the maximum number of streamline vertices.
Alternatively, ribbons can be created from an array of vertices xyz of a path curve. anlr_rot contains the angles of rotation around the edges between adjacent vertices of the path curve.
The input parameter width sets the width of the streamribbons.
Streamribbons are colored according to the total angle of rotation along the ribbon.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the plot objects created for each streamribbon.
Example:

```
[x, y, z] = meshgrid (0:0.2:4, -1:0.2:1, -1:0.2:1);
u = - x + 10;
v = 10 * z.*x;
w = - 10 * y.*x;
streamribbon (x, y, z, u, v, w, [0, 0], [0, 0.6], [0, 0]);
view (3);
```

See also: [streamline], page 343, [stream3], page 344, [streamtube], page 341, [ostreamtube], page 342.
streamtube ( $x, y, z, u, v, w, s x, s y, s z$ )
streamtube (u, v, w, sx, sy, sz)
streamtube (xyz, x, y, z, div)
streamtube (xyz, div)
streamtube (xyz, dia)
streamtube (..., options)
streamtube (hax, ...)
$h=$ streamtube (...)
Plot tubes scaled by the divergence along streamlines.
streamtube draws tubes whose diameter is scaled by the divergence of a vector field. The vector field is given by [ $u, v, w]$ and is defined over a rectangular grid given by $[x, y, z]$. The tubes start at the seed points [sx, sy, sz] and are plot along streamlines.
streamtube can also be called with a cell array containing pre-computed streamline data. To do this, xyz must be created with the stream3 command. div is used to scale the tubes. In order to plot tubes scaled by the vector field divergence, div must be calculated with the divergence command.
A tube diameter of zero corresponds to the smallest scaling value along the streamline and the largest tube diameter corresponds to the largest scaling value.
It is also possible to draw a tube along an arbitrary array of vertices xyz. The tube diameter can be specified by the vertex array dia or by a constant.

The input parameter options is a 2-D vector of the form [scale, n]. The first parameter scales the tube diameter (default 1). The second parameter specifies the number of vertices that are used to construct the tube circumference (default 20).
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a graphics handle to the plot objects created for each tube.

See also: [stream3], page 344, [streamline], page 343, [streamribbon], page 340, [ostreamtube], page 342.
ostreamtube ( $x, y, z, u, v, w, s x, s y, s z)$
ostreamtube (u, v, w, sx, sy, sz)
ostreamtube (xyz, x, y, z, u, v, w)
ostreamtube (..., options)
ostreamtube (hax, ...)
h = ostreamtube (...)
Calculate and display streamtubes.
Streamtubes are approximated by connecting circular crossflow areas along a streamline. The expansion of the flow is determined by the local crossflow divergence.
The vector field is given by [ $u, v, w]$ and is defined over a rectangular grid given by $[x, y, z]$. The streamtubes start at the seed points [sx, sy, sz].
The tubes are colored based on the local vector field strength.
The input parameter options is a 2-D vector of the form [scale, n]. The first parameter scales the start radius of the streamtubes (default 1 ). The second parameter specifies the number of vertices that are used to construct the tube circumference (default 20).
ostreamtube can be called with a cell array containing pre-computed streamline data. To do this, xyz must be created with the stream3 function. This option is useful if you need to alter the integrator step size or the maximum number of vertices of the streamline.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the plot objects created for each streamtube.

Example:

```
[x, y, z] = meshgrid (-1:0.1:1, -1:0.1:1, -3:0.1:0);
u = -x / 10 - y;
v = x - y / 10;
w = - ones (size (x)) / 10;
ostreamtube (x, y, z, u, v, w, 1, 0, 0);
```

See also: [stream3], page 344, [streamline], page 343, [streamribbon], page 340, [streamtube], page 341.
streamline ( $x, y, z, u, v, w, s x, s y, s z$ )
streamline (u, v, w, sx, sy, sz)
streamline (..., options)
streamline (hax, ...)
$h=$ streamline (...)
Plot streamlines of 2-D or 3-D vector fields.
Plot streamlines of a 2-D or 3-D vector field given by [u, v] or [u,v,w]. The vector field is defined over a rectangular grid given by $[x, y]$ or $[x, y, z]$. The streamlines start at the seed points [sx, sy] or [sx, sy, sz].
The input parameter options is a 2-D vector of the form [stepsize, max_vertices]. The first parameter specifies the step size used for trajectory integration (default 0.1). A negative value is allowed which will reverse the direction of integration. The second parameter specifies the maximum number of segments used to create a streamline (default 10,000).

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a graphics handle to the hggroup comprising the field lines.

Example:

```
[x, y] = meshgrid (-1.5:0.2:2, -1:0.2:2);
u = - x / 4 - y;
v = x - y / 4;
streamline (x, y, u, v, 1.7, 1.5);
```

See also: [stream2], page 343, [stream3], page 344, [streamribbon], page 340, [streamtube], page 341, [ostreamtube], page 342.
$x y=\operatorname{stream} 2(x, y, u, v, s x, s y)$
xy = stream2 (u, v, sx, sy)
xy = stream2 (..., options)
Compute 2-D streamline data.
Calculates streamlines of a vector field given by [u, v]. The vector field is defined over a rectangular grid given by $[x, y]$. The streamlines start at the seed points [sx, sy]. The returned value xy contains a cell array of vertex arrays. If the starting point is outside the vector field, [] is returned.

The input parameter options is a 2-D vector of the form [stepsize, max_vertices]. The first parameter specifies the step size used for trajectory integration (default 0.1). A negative value is allowed which will reverse the direction of integration. The second parameter specifies the maximum number of segments used to create a streamline (default 10,000).
The return value $x y$ is a nverts x 2 matrix containing the coordinates of the field line segments.
Example:

```
[x, y] = meshgrid (0:3);
u = 2 * x;
v = y;
xy = stream2 (x, y, u, v, 1.0, 0.5);
```

See also: [streamline], page 343, [stream3], page 344 .
$x y z=\operatorname{stream} 3(x, y, z, u, v, w, s x, s y, s z)$
xyz $=$ stream3 (u, v, w, sx, sy, sz)
xyz = stream3 (..., options)
Compute 3-D streamline data.
Calculate streamlines of a vector field given by [ $u, v, w$ ]. The vector field is defined over a rectangular grid given by $[x, y, z]$. The streamlines start at the seed points [sx, sy, sz]. The returned value xyz contains a cell array of vertex arrays. If the starting point is outside the vector field, [] is returned.

The input parameter options is a 2-D vector of the form [stepsize, max_vertices]. The first parameter specifies the step size used for trajectory integration (default 0.1). A negative value is allowed which will reverse the direction of integration. The second parameter specifies the maximum number of segments used to create a streamline (default 10,000).
The return value $x y z$ is a nverts $x 3$ matrix containing the coordinates of the field line segments.
Example:

```
[x, y, z] = meshgrid (0:3);
u = 2 * x;
v = y;
w = 3 * z;
xyz = stream3 (x, y, z, u, v, w, 1.0, 0.5, 0.0);
```

See also: [stream2], page 343, [streamline], page 343, [streamribbon], page 340, [streamtube], page 341, [ostreamtube], page 342.

```
compass (u,v)
compass (z)
compass (..., style)
compass (hax,...)
h = compass (...)
```

Plot the ( $u, v$ ) components of a vector field emanating from the origin of a polar plot.
The arrow representing each vector has one end at the origin and the tip at $[u(i)$, $v(i)]$. If a single complex argument $z$ is given, then $u=r e a l(z)$ and $v=i m a g(z)$.
The style to use for the plot can be defined with a line style style of the same format as the plot command.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a vector of graphics handles to the line objects representing the drawn vectors.

```
a = toeplitz ([1;randn(9,1)], [1,randn(1,9)]);
compass (eig (a));
```

See also: [polar], page 337, [feather], page 345, [quiver], page 339, [rose], page 329, [plot], page 316.
feather (u, v)
feather ( $z$ )
feather (..., style)
feather (hax, ...)
$h=$ feather (...)
Plot the ( $u, v$ ) components of a vector field emanating from equidistant points on the x -axis.
If a single complex argument $z$ is given, then $u=$ real ( $z$ ) and $v=i m a g(z)$.
The style to use for the plot can be defined with a line style style of the same format as the plot command.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the line objects representing the drawn vectors.

```
phi = [0 : 15 : 360] * pi/180;
feather (sin (phi), cos (phi));
```

See also: [plot], page 316, [quiver], page 339, [compass], page 344.

```
pcolor (x, y, c)
pcolor (c)
pcolor (hax,...)
h = pcolor (...)
```

Produce a 2-D density plot.
A pcolor plot draws rectangles with colors from the matrix $c$ over the two-dimensional region represented by the matrices $x$ and $y . x$ and $y$ are the coordinates of the mesh's vertices and are typically the output of meshgrid. If $x$ and $y$ are vectors, then a typical vertex is $(x(\mathrm{j}), y(\mathrm{i}), c(\mathrm{i}, \mathrm{j}))$. Thus, columns of $c$ correspond to different $x$ values and rows of $c$ correspond to different $y$ values.
The values in $c$ are scaled to span the range of the current colormap. Limits may be placed on the color axis by the command caxis, or by setting the clim property of the parent axis.
The face color of each cell of the mesh is determined by interpolating the values of $c$ for each of the cell's vertices; Contrast this with imagesc which renders one cell for each element of $c$.
shading modifies an attribute determining the manner by which the face color of each cell is interpolated from the values of $c$, and the visibility of the cells' edges. By default the attribute is "faceted", which renders a single color for each cell's face with the edge visible.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
See also: [caxis], page 350, [shading], page 382, [meshgrid], page 373, [contour], page 330, [imagesc], page 866.
area (y)
area ( $x, y$ )
area (..., lvl)
area (..., prop, val, ...)
area (hax, ...)
$h=\operatorname{area}(. .$.
Area plot of the columns of $y$.
This plot shows the contributions of each column value to the row sum. It is functionally similar to plot ( $x$, cumsum ( $y, 2$ ) ), except that the area under the curve is shaded.
If the $x$ argument is omitted it defaults to 1:rows ( $y$ ). A value lvl can be defined that determines where the base level of the shading under the curve should be defined. The default level is 0 .
Additional property/value pairs are passed directly to the underlying patch object. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the hggroup object comprising the area patch objects. The "BaseValue" property of the hggroup can be used to adjust the level where shading begins.
Example: Verify identity $\sin ^{\wedge} 2+\cos ^{\wedge} 2=1$

```
t = linspace (0, 2*pi, 100)';
y = [sin(t).^2, cos(t).^2];
area (t, y);
legend ("sin^2", "cos^2", "location", "NorthEastOutside");
```

See also: [plot], page 316, [patch], page 425.
fill ( $x, y, c$ )
fill (x1, y1, c1, x2, y2, c2)
fill (..., prop, val)
fill (hax, ...)
$h=$ fill (...)
Create one or more filled 2-D polygons.
The inputs $x$ and $y$ are the coordinates of the polygon vertices. If the inputs are matrices then the rows represent different vertices and each column produces a different polygon. fill will close any open polygons before plotting.
The input $c$ determines the color of the polygon. The simplest form is a single color specification such as a plot format or an RGB-triple. In this case the polygon(s) will
have one unique color. If $c$ is a vector or matrix then the color data is first scaled using caxis and then indexed into the current colormap. A row vector will color each polygon (a column from matrices $x$ and $y$ ) with a single computed color. A matrix $c$ of the same size as $x$ and $y$ will compute the color of each vertex and then interpolate the face color between the vertices.
Multiple property/value pairs for the underlying patch object may be specified, but they must appear in pairs. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the created patch objects.
Example: red square

```
vertices = [0 0
    10
    11
    0 1];
fill (vertices(:,1), vertices(:,2), "r");
axis ([-0.5 1.5, -0.5 1.5])
axis equal
```

See also: [patch], page 425, [fill3], page 347, [caxis], page 350, [colormap], page 870.

```
fill3 (x, y, z, c)
fill3 (x1, y1, z1, c1, x2, y2, z2, c2)
fill3 (..., prop, val)
fill3 (hax, ...)
h = fill3 (...)
Create one or more filled 3-D polygons.
```

The inputs $x, y$, and $z$ are the coordinates of the polygon vertices. If the inputs are matrices then the rows represent different vertices and each column produces a different polygon. fill3 will close any open polygons before plotting.
The input $c$ determines the color of the polygon. The simplest form is a single color specification such as a plot format or an RGB-triple. In this case the polygon(s) will have one unique color. If $c$ is a vector or matrix then the color data is first scaled using caxis and then indexed into the current colormap. A row vector will color each polygon (a column from matrices $x, y$, and $z$ ) with a single computed color. A matrix $c$ of the same size as $x, y$, and $z$ will compute the color of each vertex and then interpolate the face color between the vertices.
Multiple property/value pairs for the underlying patch object may be specified, but they must appear in pairs. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the created patch objects.

Example: oblique red rectangle

```
vertices = [0 0 0
        1 1 0
        1 1 1
        0 0 1];
fill3 (vertices(:,1), vertices(:,2), vertices(:,3), "r");
axis ([-0.5 1.5, -0.5 1.5, -0.5 1.5]);
axis ("equal");
grid ("on");
view (-80, 25);
```

See also: [patch], page 425, [fill], page 346, [caxis], page 350, [colormap], page 870.

```
comet (y)
```

comet ( $x, y$ )
comet ( $x, y, p$ )
comet (hax, ...)

Produce a simple comet style animation along the trajectory provided by the input coordinate vectors ( $x, y$ ).
If $x$ is not specified it defaults to the indices of $y$.
The speed of the comet may be controlled by $p$, which represents the time each point is displayed before moving to the next one. The default for $p$ is $5 /$ numel (y).
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [comet3], page 348.

```
comet3 (z)
comet3 (x, y, z)
comet3 (x, y, z, p)
comet3 (hax,...)
```

Produce a simple comet style animation along the trajectory provided by the input coordinate vectors ( $x, y, z$ ).
If only $z$ is specified then $x, y$ default to the indices of $z$.
The speed of the comet may be controlled by $p$, which represents the time each point is displayed before moving to the next one. The default for $p$ is $5 /$ numel ( $z$ ).
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [comet], page 348.

### 15.2.1.1 Axis Configuration

The axis function may be used to change the axis limits of an existing plot and various other axis properties, such as the aspect ratio and the appearance of tic marks. By default, high level plotting functions such as plot reset axes properties. Any customization of properties, for example by calling axis, xlim, etc., should happen after the plot is done or, alternatively, after calling the [hold function], page 403.

```
axis ()
axis ([x_lo x_hi])
axis ([x_lo x_hi y_lo y_hi])
axis ([x_lo x_hi y_lo y_hi z_lo z_hi])
axis ([x_lo x_hi y_lo y_hi z_lo z_hi c_lo c_hi])
axis (option)
axis (option1, option2,...)
axis (hax,...)
limits = axis ()
```

Set axis limits and appearance.
The argument limits should be a 2 -, 4 -, 6 -, or 8 -element vector. The first and second elements specify the lower and upper limits for the x -axis. The third and fourth specify the limits for the $y$-axis, the fifth and sixth specify the limits for the $z$-axis, and the seventh and eighth specify the limits for the color axis. The special values -Inf and Inf may be used to indicate that the limit should be automatically computed based on the data in the axes.
Without any arguments, axis turns autoscaling on.
With one output argument, limits = axis returns the current axis limits.
The vector argument specifying limits is optional, and additional string arguments may be used to specify various axis properties.
The following options control the aspect ratio of the axes.
"equal" Force x -axis unit distance to equal y -axis (and z -axis) unit distance.
"square" Force a square axis aspect ratio.
"vis3d" Set aspect ratio modes ("DataAspectRatio", "PlotBoxAspectRatio") to "manual" for rotation without stretching.
"normal"
"fill" Restore default automatically computed aspect ratios.
The following options control the way axis limits are interpreted.
"auto"
"auto[xyz]"
"auto [xyz]"
Set nice auto-computed limits around the data for all axes, or only the specified axes.
"manual" Fix the current axes limits.
"tight" Fix axes to the limits of the data.
"image" Equivalent to "tight" and "equal".
The following options affect the appearance of tick marks.
"tic"
"tic[xyz]"
"tic [xyz]"
Turn tick marks on for all axes, or turn them on for the specified axes and off for the remainder.
"label"
"label[xyz]"
"label [xyz]"
Turn tick labels on for all axes, or turn them on for the specified axes and off for the remainder.
"nolabel"
Turn tick labels off for all axes.
Note: If there are no tick marks for an axes then there can be no labels.
The following options affect the direction of increasing values on the axes.
"xy" Default y-axis, larger values are near the top.
"ij" Reverse y-axis, smaller values are near the top.
The following options affects the visibility of the axes.
"on" Make the axes visible.
"off" Hide the axes.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
Example 1: set X/Y limits and force a square aspect ratio

```
axis ([1, 2, 3, 4], "square");
```

Example 2: enable tick marks on all axes, enable tick mark labels only on the y-axis axis ("tic", "labely");
See also: [xlim], page 351, [ylim], page 351, [zlim], page 351, [caxis], page 350, [daspect], page 383, [pbaspect], page 384, [box], page 394, [grid], page 394.

Similarly the axis limits of the colormap can be changed with the caxis function.

```
caxis ([cmin cmax])
```

caxis ("auto")
caxis ("manual")
caxis (hax, ...)
limits = caxis ()

Query or set color axis limits for plots.
The limits argument should be a 2 -element vector specifying the lower and upper limits to assign to the first and last value in the colormap. Data values outside this range are clamped to the first and last colormap entries.
If the "auto" option is given then automatic colormap limits are applied. The automatic algorithm sets cmin to the minimum data value and cmax to the maximum data value. If "manual" is specified then the "climmode" property is set to "manual" and the numeric values in the "clim" property are used for limits.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
Called without arguments the current color axis limits are returned.

Programming Note: The color axis affects the display of image, patch, and surface graphics objects, but only if the "cdata" property has indexed data and the "cdatamapping" property is set to "scaled". Graphic objects with true color cdata, or "direct" cdatamapping are not affected.

See also: [colormap], page 870, [axis], page 348.
The xlim, ylim, and zlim functions may be used to get or set individual axis limits. Each has the same form.

```
xlimits = xlim ()
xmode = xlim ("mode")
xlim ([x_lo x_hi])
xlim ("auto")
xlim ("manual")
xlim (hax,...)
Query or set the limits of the x -axis for the current plot.
```

Called without arguments xlim returns the x-axis limits of the current plot.
With the input query "mode", return the current x-limit calculation mode which is either "auto" or "manual".
If passed a 2-element vector [x_lo x_hi], the limits of the x-axis are set to these values and the mode is set to "manual". The special values -Inf and Inf can be used to indicate that either the lower axis limit or upper axis limit should be automatically calculated.
The current plotting mode can be changed by using either "auto" or "manual" as the argument.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
Programming Note: The xlim function operates by modifying the "xlim" and "xlimmode" properties of an axes object. These properties can be directly inspected and altered with get/set.
See also: [ylim], page 351, [zlim], page 351, [axis], page 348, [set], page 430, [get], page 430, [gca], page 429.

The xticks, yticks, zticks, rticks, and thetaticks functions may be used to get or set the tick mark locations and modes on the respective axis. Each has the same form, although mode options are not currently available for rticks, and thetaticks.

```
tickval = xticks
mode = xticks ("mode")
xticks (tickval)
xticks ("auto")
xticks ("manual")
... = xticks (hax,...)
```

Query or set the tick values on the x -axis of the current axis.
When called without an argument, return the current tick locations as specified in the "xtick" axes property. These locations can be changed by calling xticks with a vector of tick values. Note: ascending order is not required.

When called with argument "mode", xticks returns the current value of the axes property "xtickmode". This property can be changed by calling xticks with either "auto" (algorithm determines tick positions) or "manual" (tick values remain fixed regardless of axes resizing or rotation). Note: Specifying xtick values will also set the property "xtickmode" to "manual".
If the first argument hax is an axes handle, then operate on this axis rather than the current axes returned by gca.
Requesting a return value when calling xticks to set a property value will result in an error.
See also: [xticklabels], page 352, [yticks], page 351, [zticks], page 351, [rticks], page 351, [thetaticks], page 351, [get], page 430, [set], page 430.

The xticklabels, yticklabels, and zticklabels functions may be used to get or set the label assigned to each tick location and the labeling mode on the respective axis. Each has the same form.

```
tickval = xticklabels
mode = xticklabels ("mode")
xticklabels (tickval)
xticklabels ("auto")
xticklabels ("manual")
... = xticklabels (hax,...)
    Query or set the tick labels on the x-axis of the current axis.
```

When called without an argument, return a cell array of strings of the current tick labels as specified in the "xticklabel" axes property. These labels can be changed by calling xticklabels with a cell array of strings. Note: a vector of numbers will be mapped to a cell array of strings. If fewer labels are specified than the current number of ticks, blank labels will be appended to the array.
When called with argument "mode", xticklabels returns the current value of the axes property "xticklabelmode". This property can be changed by calling xticklabels with either "auto" (algorithm determines tick labels) or "manual" (tick labels remain fixed). Note: Specifying xticklabel values will also set the "xticklabelmode" and "xticks" properties to "manual".
If the first argument hax is an axes handle, then operate on this axis rather than the current axes returned by gca.
Requesting a return value when calling xticklabels to set a property value will result in an error.

See also: [xticks], page 351, [yticklabels], page 352, [zticklabels], page 352, [get], page 430, [set], page 430.

The xtickangle, ytickangle, and ztickangle functions may be used to get or set the rotation angle of labels for the respective axis. Each has the same form.

```
angle = xtickangle ()
angle = xtickangle (hax)
xtickangle (angle)
```

xtickangle (hax, angle)
Query or set the rotation angle of the tick labels on the x -axis of the current axes.
When called without an argument, return the rotation angle in degrees of the tick labels as specified in the axes property "XTickLabelRotation". When called with a numeric scalar angle, rotate the tick labels counterclockwise to angle degrees.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
Programming Note: Requesting a return value while also setting a specified rotation will result in an error.
See also: [ytickangle], page 352, [ztickangle], page 352, [get], page 430, [set], page 430.

### 15.2.1.2 Two-dimensional Function Plotting

Octave can plot a function from a function handle or string defining the function without the user needing to explicitly create the data to be plotted. The function fplot also generates two-dimensional plots with linear axes using a function name and limits for the range of the x -coordinate instead of the x and y data. For example,
fplot (@sin, [-10, 10], 201);
produces a plot that is equivalent to the one above, but also includes a legend displaying the name of the plotted function.

```
fplot (fn)
fplot (fn, limits)
fplot (..., tol)
fplot (..., n)
fplot (..., fmt)
fplot (..., property, value,...)
fplot (hax,...)
[x, y] = fplot (...)
```

Plot a function fn within the range defined by limits.
fn is a function handle, inline function, or string containing the name of the function to evaluate.
The limits of the plot are of the form [xlo, xhi] or [xlo, xhi, ylo, yhi]. If no limits are specified the default is $[-5,5]$.
The next three arguments are all optional and any number of them may be given in any order.
tol is the relative tolerance to use for the plot and defaults to $2 \mathrm{e}-3(.2 \%)$.
$n$ is the minimum number of points to use. When $n$ is specified, the maximum stepsize will be (xhi-xlo) / $n$. More than $n$ points may still be used in order to meet the relative tolerance requirement.
The fmt argument specifies the linestyle to be used by the plot command.
Multiple property-value pairs may also be specified, but they must appear in pairs. These arguments are applied to the line objects drawn by plot.
The full list of line properties is documented at Section 15.3.3.5 [Line Properties], page 450.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
With no output arguments, the results are immediately plotted. With two output arguments, the 2-D plot data is returned. The data can subsequently be plotted manually with plot ( $x, y$ ).
Example:
fplot (@cos, [0, 2*pi])
fplot ("[cos(x), sin(x)]", [0, 2*pi])
Programming Notes:
fplot works best with continuous functions. Functions with discontinuities are unlikely to plot well. This restriction may be removed in the future.
fplot performance is better when the function accepts and returns a vector argument. Consider this when writing user-defined functions and use element-by-element operators such as.$*$, ./, etc.
See also: [ezplot], page 354, [plot], page 316.
Other functions that can create two-dimensional plots directly from a function include ezplot, ezcontour, ezcontourf and ezpolar.
ezplot (f)
ezplot (f2v)
ezplot (fx, fy)
ezplot (..., dom)
ezplot (..., n)
ezplot (hax, ...)
$h=$ ezplot (...)
Plot the 2-D curve defined by the function $f$.
The function $f$ may be a string, inline function, or function handle and can have either one or two variables. If $f$ has one variable, then the function is plotted over the domain $-2 *$ pi $<x<2 *$ pi with 500 points.
If $f 2 v$ is a function of two variables then the implicit function $f(x, y)=0$ is calculated over the meshed domain $-2 * \mathrm{pi}<=\mathrm{x} \mid \mathrm{y}<=2 * \mathrm{pi}$ with 60 points in each dimension.
For example:

```
ezplot (@(x, y) x.^2 - y.^2 - 1)
```

If two functions are passed as inputs then the parametric function

$$
\begin{aligned}
& x=f x(t) \\
& y=f y(t)
\end{aligned}
$$

is plotted over the domain $-2 *$ pi $<=t<=2 *$ pi with 500 points.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$, or $t$ for a parametric plot. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in plotting the function.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a vector of graphics handles to the created line objects.
See also: [plot], page 316, [ezplot3], page 384, [ezpolar], page 356, [ezcontour], page 355, [ezcontourf], page 355, [ezmesh], page 385, [ezmeshc], page 386, [ezsurf], page 386, [ezsurfc], page 387.
ezcontour (f)
ezcontour (..., dom)
ezcontour (..., n)
ezcontour (hax, ...)
$h=$ ezcontour (...)
Plot the contour lines of a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 *$ pi $<=x \mid y<=2 *$ pi with 60 points in each dimension.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.
Example:

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezcontour (f, [-3, 3]);
```

See also: [contour], page 330, [ezcontourf], page 355, [ezplot], page 354, [ezmeshc], page 386, [ezsurfc], page 387.

```
ezcontourf (f)
ezcontourf (..., dom)
ezcontourf (..., n)
ezcontourf (hax,...)
h = ezcontourf (...)
```

Plot the filled contour lines of a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 * \mathrm{pi}<=\mathrm{x} \mid \mathrm{y}<=2 * \mathrm{pi}$ with 60 points in each dimension.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.

Example:

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezcontourf (f, [-3, 3]);
```

See also: [contourf], page 331, [ezcontour], page 355, [ezplot], page 354, [ezmeshc], page 386, [ezsurfc], page 387.
ezpolar (f)
ezpolar (..., dom)
ezpolar (..., n)
ezpolar (hax, ...)
h = ezpolar (...)
Plot a 2-D function in polar coordinates.
The function $f$ is a string, inline function, or function handle with a single argument. The expected form of the function is rho $=f($ theta $)$. By default the plot is over the domain $0<=$ theta <= $2 *$ pi with 500 points.
If dom is a two element vector, it represents the minimum and maximum values of theta.
$n$ is a scalar defining the number of points to use in plotting the function.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.
Example:

```
ezpolar (@(t) sin (5/4 * t), [0, 8*pi]);
```

See also: [polar], page 337, [ezplot], page 354.

### 15.2.1.3 Two-dimensional Geometric Shapes

```
rectangle ()
rectangle (..., "Position", pos)
rectangle (..., "Curvature", curv)
rectangle (..., "EdgeColor", ec)
rectangle (..., "FaceColor", fc)
rectangle (hax,...)
h = rectangle (...)
```

Draw a rectangular patch defined by pos and curv.
The variable pos(1:2) defines the lower left-hand corner of the patch and pos(3:4) defines its width and height. By default, the value of pos is $[0,0,1,1]$.
The variable curv defines the curvature of the sides of the rectangle and may be a scalar or two-element vector with values between 0 and 1 . A value of 0 represents no curvature of the side, whereas a value of 1 means that the side is entirely curved into the arc of a circle. If curv is a two-element vector, then the first element is the curvature along the x -axis of the patch and the second along y -axis.
If curv is a scalar, it represents the curvature of the shorter of the two sides of the rectangle and the curvature of the other side is defined by

```
min (pos(1:2)) / max (pos(1:2)) * curv
```

Additional property/value pairs are passed to the underlying patch command. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457. If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created rectangle object.
See also: [patch], page 425, [line], page 424, [cylinder], page 388, [ellipsoid], page 389, [sphere], page 388.

### 15.2.2 Three-Dimensional Plots

The function mesh produces mesh surface plots. For example,

```
tx = ty = linspace (-8, 8, 41)';
[xx, yy] = meshgrid (tx, ty);
r = sqrt (xx .^ 2 + yy .^ 2) + eps;
tz = sin (r) ./ r;
mesh (tx, ty, tz);
xlabel ("tx");
ylabel ("ty");
zlabel ("tz");
title ("3-D Sombrero plot");
```

produces the familiar "sombrero" plot shown in Figure 15.5. Note the use of the function meshgrid to create matrices of X and Y coordinates to use for plotting the Z data. The ndgrid function is similar to meshgrid, but works for N -dimensional matrices.

## 3-D Sombrero plot



Figure 15.5: Mesh plot.
The meshc function is similar to mesh, but also produces a plot of contours for the surface.

The plot3 function displays arbitrary three-dimensional data, without requiring it to form a surface. For example,

```
\(\mathrm{t}=0: 0.1: 10 * \mathrm{pi}\);
r = linspace (0, 1, numel (t));
\(z=\) linspace ( 0,1 , numel ( \(t\) )) ;
plot3 (r.*sin (t), r.*cos (t), z);
xlabel ("r.*sin (t)");
ylabel ("r.*cos (t)");
zlabel ("z");
title ("plot3 display of 3-D helix");
```

displays the spiral in three dimensions shown in Figure 15.6.


Figure 15.6: Three-dimensional spiral.
Finally, the view function changes the viewpoint for three-dimensional plots.

```
mesh (x, y, z)
mesh (z)
mesh (..., c)
mesh (..., prop, val, ...)
mesh (hax, ...)
h = mesh (...)
```

Plot a 3-D wireframe mesh.
The wireframe mesh is plotted using rectangles. The vertices of the rectangles [x, $y]$ are typically the output of meshgrid. over a 2-D rectangular region in the $x-y$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns $(z), y=1$ : rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.

The color of the mesh is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.

Optionally, the color of the mesh can be specified independently of $z$ by supplying a color matrix, $c$.
Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
See also: [ezmesh], page 385, [meshc], page 359, [meshz], page 359, [trimesh], page 828, [contour], page 330, [surf], page 360, [surface], page 426, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350 .

```
meshc ( \(x, y, z\) )
meshc ( \(z\) )
meshc (..., c)
meshc (..., prop, val, ...)
meshc (hax, ...)
\(h=m e s h c(. .\).
```

Plot a 3-D wireframe mesh with underlying contour lines.
The wireframe mesh is plotted using rectangles. The vertices of the rectangles [x, $y]$ are typically the output of meshgrid. over a 2 -D rectangular region in the $x-y$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ : rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The color of the mesh is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally the color of the mesh can be specified independently of $z$ by supplying a color matrix, $c$.
Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a 2 -element vector with a graphics handle to the created surface object and to the created contour plot.
See also: [ezmeshc], page 386, [mesh], page 358, [meshz], page 359, [contour], page 330, [surfc], page 361, [surface], page 426, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.

```
meshz (x, y, z)
meshz (z)
meshz (..., c)
meshz (..., prop, val, ...)
meshz (hax,...)
```

$h=\operatorname{meshz}(.$.
Plot a 3-D wireframe mesh with a surrounding curtain.
The wireframe mesh is plotted using rectangles. The vertices of the rectangles $[x$, $y]$ are typically the output of meshgrid. over a $2-\mathrm{D}$ rectangular region in the $\mathrm{x}-\mathrm{y}$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ :rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The color of the mesh is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally the color of the mesh can be specified independently of $z$ by supplying a color matrix, $c$.

Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
See also: [mesh], page 358, [meshc], page 359, [contour], page 330, [surf], page 360, [surface], page 426, [waterfall], page 383, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.
hidden
hidden on
hidden off
mode $=$ hidden (...)
Control mesh hidden line removal.
When called with no argument the hidden line removal state is toggled.
When called with one of the modes "on" or "off" the state is set accordingly.
The optional output argument mode is the current state.
Hidden Line Removal determines what graphic objects behind a mesh plot are visible. The default is for the mesh to be opaque and lines behind the mesh are not visible. If hidden line removal is turned off then objects behind the mesh can be seen through the faces (openings) of the mesh, although the mesh grid lines are still opaque.
See also: [mesh], page 358, [meshc], page 359, [meshz], page 359, [ezmesh], page 385, [ezmeshc], page 386, [trimesh], page 828, [waterfall], page 383.
surf $(x, y, z)$
surf ( $z$ )
surf (..., c)
surf (..., prop, val, ...)
surf (hax,...)
$h=\operatorname{surf}(.$.
Plot a 3-D surface mesh.
The surface mesh is plotted using shaded rectangles. The vertices of the rectangles $[x, y]$ are typically the output of meshgrid. over a $2-\mathrm{D}$ rectangular region in the $\mathrm{x}-\mathrm{y}$
plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ :rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The color of the surface is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally, the color of the surface can be specified independently of $z$ by supplying a color matrix, $c$.
Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Note: The exact appearance of the surface can be controlled with the shading command or by using set to control surface object properties.
See also: [ezsurf], page 386, [surfc], page 361, [surfl], page 362, [surfnorm], page 363, [trisurf], page 828, [contour], page 330, [mesh], page 358, [surface], page 426, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.
surfc $(x, y, z)$
surfc ( $z$ )
surfc (..., c)
surfc (..., prop, val, ...)
surfc (hax, ...)
$h=\operatorname{surfc}(. .$.
Plot a 3-D surface mesh with underlying contour lines.
The surface mesh is plotted using shaded rectangles. The vertices of the rectangles $[x, y]$ are typically the output of meshgrid. over a 2-D rectangular region in the $x-y$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ : rows ( $z$ ). Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The color of the surface is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally, the color of the surface can be specified independently of $z$ by supplying a color matrix, $c$.
Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.

Note: The exact appearance of the surface can be controlled with the shading command or by using set to control surface object properties.

See also: [ezsurfc], page 387, [surf], page 360, [surff], page 362, [surfnorm], page 363, [trisurf], page 828, [contour], page 330, [mesh], page 358, [surface], page 426, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.

```
surfl (z)
surfl (x, y, z)
surfl (..., lsrc)
surfl (x, y, z, lsrc, P)
surfl (..., "cdata")
surfl (..., "light")
surfl (hax,...)
h = surfl (...)
    Plot a 3-D surface using shading based on various lighting models.
```

The surface mesh is plotted using shaded rectangles. The vertices of the rectangles $[x, y]$ are typically the output of meshgrid. over a 2-D rectangular region in the $x-y$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ :rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The default lighting mode "cdata", changes the cdata property of the surface object to give the impression of a lighted surface.

The alternate mode "light" creates a light object to illuminate the surface.
The light source location may be specified using lsrc which can be a 2 -element vector [azimuth, elevation] in degrees, or a 3 -element vector $[\mathrm{lx}, \mathrm{ly}, \mathrm{lz}]$. The default value is rotated 45 degrees counterclockwise to the current view.
The material properties of the surface can specified using a 4 -element vector $P=$ [AM D SP exp] which defaults to $p=\left[\begin{array}{lll}0.55 & 0.6 & 0.4 \\ 10\end{array}\right]$.
"AM" strength of ambient light
"D" strength of diffuse reflection
"SP" strength of specular reflection
"EXP" specular exponent
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Example:

```
colormap (bone (64));
surfl (peaks);
shading interp;
```

See also: [diffuse], page 370, [specular], page 370, [surf], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.
surfnorm ( $x, y, z$ )
surfnorm ( $z$ )
surfnorm (..., prop, val, ...)
surfnorm (hax, ...)
[nx, ny, nz] = surfnorm (...)
Find the vectors normal to a meshgridded surface.
If $x$ and $y$ are vectors, then a typical vertex is $(x(j), y(i), z(i, j))$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values. If only a single input $z$ is given then $x$ is taken to be 1 :columns ( $z$ ) and $y$ is 1:rows (z).

If no return arguments are requested, a surface plot with the normal vectors to the surface is plotted.
Any property/value input pairs are assigned to the surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If output arguments are requested then the components of the normal vectors are returned in $n x, n y$, and $n z$ and no plot is made. The normal vectors are unnormalized (magnitude $!=1$ ). To normalize, use

```
len = sqrt (nx.^2 + ny.^2 + nz.^2);
nx ./= len; ny ./= len; nz ./= len;
```

An example of the use of surfnorm is
surfnorm (peaks (25));
Algorithm: The normal vectors are calculated by taking the cross product of the diagonals of each of the quadrilateral faces in the meshgrid to find the normal vectors at the center of each face. Next, for each meshgrid point the four nearest normal vectors are averaged to obtain the final normal to the surface at the meshgrid point. For surface objects, the "VertexNormals" property contains equivalent information, except possibly near the boundary of the surface where different interpolation schemes may yield slightly different values.
See also: [isonormals], page 365, [quiver3], page 339, [surf], page 360, [meshgrid], page 373.

```
fv = isosurface (v, isoval)
fv = isosurface (v)
fv = isosurface (x, y, z, v, isoval)
fv = isosurface (x, y, z, v)
fvc = isosurface (..., col)
fv = isosurface (..., "noshare")
fv = isosurface (..., "verbose")
[f, v] = isosurface (...)
[f, v, c] = isosurface (...)
isosurface (...)
```

Calculate isosurface of 3-D volume data.

An isosurface connects points with the same value and is analogous to a contour plot, but in three dimensions.
The input argument $v$ is a three-dimensional array that contains data sampled over a volume.
The input isoval is a scalar that specifies the value for the isosurface. If isoval is omitted or empty, a "good" value for an isosurface is determined from $v$.
When called with a single output argument isosurface returns a structure array $f_{v}$ that contains the fields faces and vertices computed at the points $[x, y, z]=$ meshgrid ( $1: 1,1: m, 1: n$ ) where $[1, m, n]=$ size (v). The output $f v$ can be used directly as input to the patch function.
If called with additional input arguments $x, y$, and $z$ that are three-dimensional arrays with the same size as $v$ or vectors with lengths corresponding to the dimensions of $v$, then the volume data is taken at the specified points. If $x, y$, or $z$ are empty, the grid corresponds to the indices ( $1: \mathrm{n}$ ) in the respective direction (see [meshgrid], page 373 ).
The optional input argument col, which is a three-dimensional array of the same size as $v$, specifies coloring of the isosurface. The color data is interpolated, as necessary, to match isoval. The output structure array, in this case, has the additional field facevertexcdata.
If given the string input argument "noshare", vertices may be returned multiple times for different faces. The default behavior is to eliminate vertices shared by adjacent faces.
The string input argument "verbose" is supported for MATLAB compatibility, but has no effect.

Any string arguments must be passed after the other arguments.
If called with two or three output arguments, return the information about the faces $f$, vertices $v$, and color data $c$ as separate arrays instead of a single structure array.
If called with no output argument, the isosurface geometry is directly plotted with the patch command and a light object is added to the axes if not yet present.
For example,

```
[x, y, z] = meshgrid (1:5, 1:5, 1:5);
v = rand (5, 5, 5);
isosurface (x, y, z, v, .5);
```

will directly draw a random isosurface geometry in a graphics window.
An example of an isosurface geometry with different additional coloring:

```
N = 15; # Increase number of vertices in each direction
iso = .4; # Change isovalue to . }1\mathrm{ to display a sphere
lin = linspace (0, 2, N);
[x, y, z] = meshgrid (lin, lin, lin);
v = abs ((x-.5).^2 + (y-.5).^2 + (z-.5). ^2);
figure ();
subplot (2,2,1); view (-38, 20);
[f, vert] = isosurface (x, y, z, v, iso);
p = patch ("Faces", f, "Vertices", vert, "EdgeColor", "none");
pbaspect ([1 1 1 1]);
```

```
isonormals (x, y, z, v, p)
set (p, "FaceColor", "green", "FaceLighting", "gouraud");
light ("Position", [1 1 5]);
subplot (2,2,2); view (-38, 20);
p = patch ("Faces", f, "Vertices", vert, "EdgeColor", "blue");
pbaspect ([1 1 1 1]);
isonormals (x, y, z, v, p)
set (p, "FaceColor", "none", "EdgeLighting", "gouraud");
light ("Position", [1 1 5]);
subplot (2,2,3); view (-38, 20);
[f, vert, c] = isosurface (x, y, z, v, iso, y);
p = patch ("Faces", f, "Vertices", vert, "FaceVertexCData", c, ...
    "FaceColor", "interp", "EdgeColor", "none");
pbaspect ([llll
isonormals (x, y, z, v, p)
set (p, "FaceLighting", "gouraud");
light ("Position", [1 1 5]);
subplot (2,2,4); view (-38, 20);
p = patch ("Faces", f, "Vertices", vert, "FaceVertexCData", c, ...
    "FaceColor", "interp", "EdgeColor", "blue");
pbaspect ([llll
isonormals (x, y, z, v, p)
set (p, "FaceLighting", "gouraud");
light ("Position", [1 1 5]);
```

See also: [isonormals], page 365, [isocolors], page 367, [isocaps], page 366, [smooth3], page 367, [reducevolume], page 368, [reducepatch], page 369, [patch], page 425.

```
vn = isonormals (val, vert)
vn = isonormals (val, hp)
vn = isonormals (x, y, z, val, vert)
vn = isonormals (x, y, z, val, hp)
vn = isonormals (..., "negate")
isonormals (val, hp)
isonormals (x, y, z, val, hp)
isonormals (..., "negate")
```

Calculate normals to an isosurface.
The vertex normals vn are calculated from the gradient of the 3-dimensional array val (size: lxmxn) containing the data for an isosurface geometry. The normals point towards smaller values in val.

If called with one output argument vn, and the second input argument vert holds the vertices of an isosurface, then the normals vn are calculated at the vertices vert on a grid given by $[x, y, z]=$ meshgrid ( $1: 1,1: m, 1: n$ ). The output argument vn has the same size as vert and can be used to set the "VertexNormals" property of the corresponding patch.

If called with additional input arguments $x, y$, and $z$, which are 3 -dimensional arrays with the same size as val, then the volume data is taken at these points. Instead of the vertex data vert, a patch handle $h p$ can be passed to the function.

If the last input argument is the string "negate", compute the reverse vector normals of an isosurface geometry (i.e., pointed towards larger values in val).
If no output argument is given, the property "VertexNormals" of the patch associated with the patch handle $h p$ is changed directly.
See also: [isosurface], page 363, [isocolors], page 367, [smooth3], page 367.

```
fvc = isocaps (v, isoval)
fvc = isocaps (v)
fvc = isocaps (x, y, z, v, isoval)
fvc = isocaps (x, y, z, v)
fvc = isocaps (..., which_caps)
fvc = isocaps (..., which_plane)
fvc = isocaps (..., "verbose")
[faces, vertices, fvcdata] = isocaps (...)
isocaps (...)
Create end-caps for isosurfaces of 3-D data.
```

This function places caps at the open ends of isosurfaces.
The input argument $v$ is a three-dimensional array that contains data sampled over a volume.
The input isoval is a scalar that specifies the value for the isosurface. If isoval is omitted or empty, a "good" value for an isosurface is determined from v .
When called with a single output argument, isocaps returns a structure array fvc with the fields: faces, vertices, and facevertexcdata. The results are computed at the points $[x, y, z]=$ meshgrid ( $1: 1,1: m, 1: n$ ) where $[1, m, n]=\operatorname{size}(v)$. The output fve can be used directly as input to the patch function.
If called with additional input arguments $x, y$, and $z$ that are three-dimensional arrays with the same size as $v$ or vectors with lengths corresponding to the dimensions of $v$, then the volume data is taken at the specified points. If $x, y$, or $z$ are empty, the grid corresponds to the indices ( $1: \mathrm{n}$ ) in the respective direction (see [meshgrid], page 373).
The optional parameter which_caps can have one of the following string values which defines how the data will be enclosed:
"above", "a" (default)
for end-caps that enclose the data above isoval.
"below", "b"
for end-caps that enclose the data below isoval.
The optional parameter which_plane can have one of the following string values to define which end-cap should be drawn:
"all" (default)
for all of the end-caps.
"xmin" for end-caps at the lower $x$-plane of the data.
"xmax" for end-caps at the upper x-plane of the data.
"ymin" for end-caps at the lower y-plane of the data.
"ymax" for end-caps at the upper y-plane of the data.
"zmin" for end-caps at the lower z-plane of the data.
"zmax" for end-caps at the upper z-plane of the data.
The string input argument "verbose" is supported for matlab compatibility, but has no effect.
If called with two or three output arguments, the data for faces faces, vertices vertices, and the color data facevertexcdata are returned in separate arrays instead of a single structure.

If called with no output argument, the end-caps are drawn directly in the current figure with the patch command.

See also: [isosurface], page 363, [isonormals], page 365, [patch], page 425.

```
cdat = isocolors (c, v)
cdat = isocolors (x, y, z, c, v)
cdat = isocolors (x, y, z, r, g, b, v)
cdat = isocolors (r, g, b, v)
cdat = isocolors (..., p)
isocolors (...)
```

        Compute isosurface colors.
    If called with one output argument, and the first input argument \(c\) is a three-
    dimensional array that contains indexed color values, and the second input argument
    \(v\) are the vertices of an isosurface geometry, then return a matrix cdat with color data
    information for the geometry at computed points \([x, y, z]=\operatorname{meshgrid}(1: 1,1: m\),
    \(1: n\) ). The output argument \(c d a t\) can be used to manually set the "FaceVertexCData"
    property of an isosurface patch object.
    If called with additional input arguments $x, y$ and $z$ which are three-dimensional arrays of the same size as $c$ then the color data is taken at those specified points.
Instead of indexed color data $c$, isocolors can also be called with RGB values $r, g$, $b$. If input arguments $x, y, z$ are not given then meshgrid computed values are used. Optionally, a patch handle $p$ can be given as the last input argument to all function call variations and the vertex data will be extracted from the isosurface patch object. Finally, if no output argument is given then the colors of the patch given by the patch handle $p$ are changed.
See also: [isosurface], page 363, [isonormals], page 365.

```
smoothed_data = smooth3 (data)
smoothed_data = smooth3 (data, method)
smoothed_data = smooth3 (data, method, sz)
smoothed_data = smooth3 (data, method, sz, std_dev)
```

Smooth values of 3 -dimensional matrix data.
This function may be used, for example, to reduce the impact of noise in data before calculating isosurfaces.
data must be a non-singleton 3-dimensional matrix. The output smoothed_data is a matrix of the same size as data.
The option input method determines which convolution kernel is used for the smoothing process. Possible choices:
"box", "b" (default)
a convolution kernel with sharp edges.
"gaussian", "g"
a convolution kernel that is represented by a non-correlated trivariate normal distribution function.
$s z$ is either a 3 -element vector specifying the size of the convolution kernel in the $x$-, y - and z -directions, or a scalar. In the scalar case the same size is used for all three dimensions ( $[s z, s z, s z]$ ). The default value is 3 .
If method is "gaussian" then the optional input std_dev defines the standard deviation of the trivariate normal distribution function. std_dev is either a 3-element vector specifying the standard deviation of the Gaussian convolution kernel in $x-, y$ and z-directions, or a scalar. In the scalar case the same value is used for all three dimensions. The default value is 0.65 .

See also: [isosurface], page 363, [isonormals], page 365, [patch], page 425.

```
[nx, ny, nz, nv] = reducevolume (v, r)
[nx, ny, nz, nv] = reducevolume (x, y, z, v, r)
nv = reducevolume (...)
```

Reduce the volume of the dataset in $v$ according to the values in $r$.
$v$ is a matrix that is non-singleton in the first 3 dimensions.
$r$ can either be a vector of 3 elements representing the reduction factors in the x -, y -, and z -directions or a scalar, in which case the same reduction factor is used in all three dimensions.
reducevolume reduces the number of elements of $v$ by taking only every $r$-th element in the respective dimension.
Optionally, $x, y$, and $z$ can be supplied to represent the set of coordinates of $v$. They can either be matrices of the same size as $v$ or vectors with sizes according to the dimensions of $v$, in which case they are expanded to matrices (see [meshgrid], page 373).
If reducevolume is called with two arguments then $x, y$, and $z$ are assumed to match the respective indices of $v$.
The reduced matrix is returned in $n v$.
Optionally, the reduced set of coordinates are returned in $n x$, ny, and $n z$, respectively.
Examples:

```
v = reshape (1:6*8*4, [6 8 4]);
nv = reducevolume (v, [4 3 2]);
v = reshape (1:6*8*4, [6 8 4]);
x = 1:3:24; y = -14:5:11; z = linspace (16, 18, 4);
[nx, ny, nz, nv] = reducevolume (x, y, z, v, [4 3 2]);
```

See also: [isosurface], page 363, [isonormals], page 365 .

```
reduced_fv = reducepatch (fv)
reduced_fv = reducepatch (faces, vertices)
reduced_fv = reducepatch (patch_handle)
reducepatch (patch_handle)
reduced_fv = reducepatch (..., reduction_factor)
reduced_fv = reducepatch (..., "fast")
reduced_fv = reducepatch (..., "verbose")
[reduced_faces, reduces_vertices] = reducepatch (...)
```

Reduce the number of faces and vertices in a patch object while retaining the overall shape of the patch.
The input patch can be represented by a structure fv with the fields faces and vertices, by two matrices faces and vertices (see, e.g., the result of isosurface), or by a handle to a patch object patch_handle (see [patch], page 425).
The number of faces and vertices in the patch is reduced by iteratively collapsing the shortest edge of the patch to its midpoint (as discussed, e.g., here: https://libigl. github.io/libigl/tutorial/tutorial.html\#meshdecimation).
Currently, only patches consisting of triangles are supported. The resulting patch also consists only of triangles.
If reducepatch is called with a handle to a valid patch patch_handle, and without any output arguments, then the given patch is updated immediately.
If the reduction_factor is omitted, the resulting structure reduced_fv includes approximately $50 \%$ of the faces of the original patch. If reduction_factor is a fraction between 0 (excluded) and 1 (excluded), a patch with approximately the corresponding fraction of faces is determined. If reduction_factor is an integer greater than or equal to 1 , the resulting patch has approximately reduction_factor faces. Depending on the geometry of the patch, the resulting number of faces can differ from the given value of reduction_factor. This is especially true when many shared vertices are detected.
For the reduction, it is necessary that vertices of touching faces are shared. Shared vertices are detected automatically. This detection can be skipped by passing the optional string argument "fast".
With the optional string arguments "verbose", additional status messages are printed to the command window.
Any string input arguments must be passed after all other arguments.
If called with one output argument, the reduced faces and vertices are returned in a structure reduced_fv with the fields faces and vertices (see the one output option of isosurface).
If called with two output arguments, the reduced faces and vertices are returned in two separate matrices reduced_faces and reduced_vertices.
See also: [isosurface], page 363, [isonormals], page 365, [reducevolume], page 368, [patch], page 425.
shrinkfaces ( $p, s f$ )
$n f v=$ shrinkfaces ( $p, s f$ )
$n f v=$ shrinkfaces ( $f v, s f$ )
$n f v=$ shrinkfaces (f, v, sf)
[nf, nv] = shrinkfaces (...)
Reduce the size of faces in a patch by the shrink factor sf.
The patch object can be specified by a graphics handle $(p)$, a patch structure ( $f_{v}$ ) with the fields "faces" and "vertices", or as two separate matrices ( $f, v$ ) of faces and vertices.
The shrink factor $s f$ is a positive number specifying the percentage of the original area the new face will occupy. If no factor is given the default is 0.3 (a reduction to $30 \%$ of the original size). A factor greater than 1.0 will result in the expansion of faces.
Given a patch handle as the first input argument and no output parameters, perform the shrinking of the patch faces in place and redraw the patch.
If called with one output argument, return a structure with fields "faces", "vertices", and "facevertexcdata" containing the data after shrinking. This structure can be used directly as an input argument to the patch function.
Caution:: Performing the shrink operation on faces which are not convex can lead to undesirable results.
Example: a triangulated $3 / 4$ circle and the corresponding shrunken version.

```
[phi r] = meshgrid (linspace (0, 1.5*pi, 16), linspace (1, 2, 4));
tri = delaunay (phi(:), r(:));
v = [r(:).*sin(phi(:)) r(:).*cos(phi(:))];
clf ()
p = patch ("Faces", tri, "Vertices", v, "FaceColor", "none");
fv = shrinkfaces (p);
patch (fv)
axis equal
grid on
```

See also: [patch], page 425.
diffuse (sx, sy, sz, lv)
Calculate the diffuse reflection strength of a surface defined by the normal vector elements sx, sy, sz.
The light source location vector $l v$ can be given as a 2-element vector [azimuth, elevation] in degrees or as a 3-element vector $[x, y, z]$.
See also: [specular], page 370, [surfl], page 362.
specular ( $s x, s y, s z, l v, v v$ )
specular (sx, sy, sz, lv, vv, se)
Calculate the specular reflection strength of a surface defined by the normal vector elements sx, sy, sz using Phong's approximation.
The light source location and viewer location vectors are specified using parameters $l v$ and $v v$ respectively. The location vectors can given as 2-element vectors [azimuth, elevation] in degrees or as 3-element vectors [ $\mathrm{x}, \mathrm{y}, \mathrm{z}$ ].
An optional sixth argument specifies the specular exponent (spread) se. If not given, se defaults to 10 .
See also: [diffuse], page 370, [surfl], page 362.
lighting (type)
lighting (hax, type)
Set the lighting of patch or surface graphic objects.
Valid arguments for type are
"flat" Draw objects with faceted lighting effects.
"gouraud"
Draw objects with linear interpolation of the lighting effects between the vertices.
"none" Draw objects without light and shadow effects.
If the first argument hax is an axes handle, then change the lighting effects of objects in this axes, rather than the current axes returned by gca.
The lighting effects are only visible if at least one light object is present and visible in the same axes.

See also: [light], page 426, [fill], page 346, [mesh], page 358, [patch], page 425, [pcolor], page 345, [surf], page 360, [surface], page 426, [shading], page 382.

```
material shiny
material dull
material metal
material default
material ([as, ds, ss])
material ([as, ds, ss, se])
material ([as, ds, ss, se, scr])
material (hlist,...)
mtypes = material ()
refl_props = material (mtype_string)
```

Set reflectance properties for the lighting of surfaces and patches.
This function changes the ambient, diffuse, and specular strengths, as well as the specular exponent and specular color reflectance, of all patch and surface objects in the current axes. This can be used to simulate, to some extent, the reflectance properties of certain materials when used with light.
When called with a string, the aforementioned properties are set according to the values in the following table:

| mtype | ambient- <br> strength | diffuse- <br> strength | specular- <br> strength | specular- <br> exponent | specular- <br> color- <br> reflectance |
| :--- | :--- | :--- | :--- | :--- | :--- |
| "shiny" | 0.3 | 0.6 | 0.9 | 20 | 1.0 |
| "dull" | 0.3 | 0.8 | 0.0 | 10 | 1.0 |
| "metal" | 0.3 | 0.3 | 1.0 | 25 | 0.5 |
| "default" | "default" | "default" | "default" | "default" | "default" |

When called with a vector of three elements, the ambient, diffuse, and specular strengths of all patch and surface objects in the current axes are updated. An
optional fourth vector element updates the specular exponent, and an optional fifth vector element updates the specular color reflectance.
A list of graphic handles can also be passed as the first argument. In this case, the properties of these handles and all child patch and surface objects will be updated.
Additionally, material can be called with a single output argument. If called without input arguments, a column cell vector mtypes with the strings for all available materials is returned. If the one input argument mtype_string is the name of a material, a 1 x 5 cell vector refl_props with the reflectance properties of that material is returned. In both cases, no graphic properties are changed.
See also: [light], page 426, [fill], page 346, [mesh], page 358, [patch], page 425, [pcolor], page 345, [surf], page 360, [surface], page 426.

```
camlight
camlight right
camlight left
camlight headlight
camlight (az, el)
camlight (..., style)
camlight (hl,...)
camlight (hax,...)
h = camlight (...)
```

Add a light object to a figure using a simple interface.
When called with no arguments, a light object is added to the current plot and is placed slightly above and to the right of the camera's current position: this is equivalent to camlight right. The commands camlight left and camlight headlight behave similarly with the placement being either left of the camera position or centered on the camera position.
For more control, the light position can be specified by an azimuthal rotation az and an elevation angle el, both in degrees, relative to the current properties of the camera.
The optional string style specifies whether the light is a local point source ("local", the default) or placed at infinite distance ("infinite").
If the first argument hl is a handle to a light object, then act on this light object rather than creating a new object.
If the first argument hax is an axes handle, then create a new light object in this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the light object. This can be used to move or further change properties of the light object.

## Examples:

Add a light object to a plot
sphere (36);
camlight
Position the light source exactly
camlight (45, 30);

Here the light is first pitched upwards (see [camup], page 379) from the camera position (see [campos], page 376) by 30 degrees. It is then yawed by 45 degrees to the right. Both rotations are centered around the camera target (see [camtarget], page 378).
Return a handle to further manipulate the light object

```
clf
sphere (36);
hl = camlight ("left");
set (hl, "color", "r");
```

See also: [light], page 426.
lightangle (az, el)
lightangle (hax, az, el)
lightangle (hl, az, el)
hl = lightangle (...)
[az, el] = lightangle (hl)
Add a light object to the current axes using spherical coordinates.
The light position is specified by an azimuthal rotation $a z$ and an elevation angle el, both in degrees.
If the first argument hax is an axes handle, then create a new light object in this axes, rather than the current axes returned by gca.
If the first argument $h l$ is a handle to a light object, then act on this light object rather than creating a new object.
The optional return value $h l$ is a graphics handle to the light object.
Example:
Add a light object to a plot

```
clf;
sphere (36);
lightangle (45, 30);
```

See also: [light], page 426, [view], page 375, [camlight], page 372.
[xx, yy] = meshgrid (x, y)
[xx, yy, zz] = meshgrid (x, y, z)
[xx, yy] = meshgrid (x)
[xx, yy, zz] = meshgrid (x)
Given vectors of $x$ and $y$ coordinates, return matrices $x x$ and yy corresponding to a full 2-D grid.

The rows of $x x$ are copies of $x$, and the columns of $y y$ are copies of $y$. If $y$ is omitted, then it is assumed to be the same as $x$.
If the optional $z$ input is given, or $z z$ is requested, then the output will be a full 3-D grid. If $z$ is omitted and $z z$ is requested, it is assumed to be the same as $y$.
meshgrid is most frequently used to produce input for a 2 -D or 3-D function that will be plotted. The following example creates a surface plot of the "sombrero" function.

```
f = @(x,y) sin (sqrt (x.^2 + y.^2)) ./ sqrt (x.^2 + y.^2);
range = linspace (-8, 8, 41);
[X, Y] = meshgrid (range, range);
Z = f (X, Y);
surf (X, Y, Z);
```

Programming Note: meshgrid is restricted to $2-\mathrm{D}$ or $3-\mathrm{D}$ grid generation. The ndgrid function will generate 1-D through N-D grids. However, the functions are not completely equivalent. If $x$ is a vector of length $M$ and $y$ is a vector of length $N$, then meshgrid will produce an output grid which is NxM . ndgrid will produce an output which is MxN (transpose) for the same input. Some core functions expect meshgrid input and others expect ndgrid input. Check the documentation for the function in question to determine the proper input format.
See also: [ndgrid], page 374, [mesh], page 358, [contour], page 330, [surf], page 360.
[y1, y2, ..., yn] = ndgrid (x1, x2, .., xn)
[y1, y2, ..., yn] = ndgrid (x)
Given $n$ vectors $x 1, \ldots, x n$, ndgrid returns $n$ arrays of dimension $n$.
The elements of the i-th output argument contains the elements of the vector xi repeated over all dimensions different from the i-th dimension. Calling ndgrid with only one input argument $x$ is equivalent to calling ndgrid with all $n$ input arguments equal to $x$ :
$[y 1, y 2, \ldots, y n]=\operatorname{ndgrid}(x, \ldots, x)$
Programming Note: ndgrid is very similar to the function meshgrid except that the first two dimensions are transposed in comparison to meshgrid. Some core functions expect meshgrid input and others expect ndgrid input. Check the documentation for the function in question to determine the proper input format.
See also: [meshgrid], page 373.

```
plot3 (x, y, z)
plot3 (x, y, z, prop, value, ...)
plot3 (x, y, z, fmt)
plot3 (x, cplx)
plot3 (cplx)
plot3 (hax,...)
h = plot3 (...)
```

Produce 3-D plots.
Many different combinations of arguments are possible. The simplest form is

```
plot3 (x, y, z)
```

in which the arguments are taken to be the vertices of the points to be plotted in three dimensions. If all arguments are vectors of the same length, then a single continuous line is drawn. If all arguments are matrices, then each column of is treated as a separate line. No attempt is made to transpose the arguments to make the number of rows match.

If only two arguments are given, as

```
plot3 (x, cplx)
```

the real and imaginary parts of the second argument are used as the $y$ and $z$ coordinates, respectively.
If only one argument is given, as
plot3 (cplx)
the real and imaginary parts of the argument are used as the $y$ and $z$ values, and they are plotted versus their index.
Arguments may also be given in groups of three as

```
plot3 (x1, y1, z1, x2, y2, z2, ...)
```

in which each set of three arguments is treated as a separate line or set of lines in three dimensions.
To plot multiple one- or two-argument groups, separate each group with an empty format string, as

```
plot3 (x1, c1, "", c2, "", ...)
```

Multiple property-value pairs may be specified which will affect the line objects drawn by plot3. If the fmt argument is supplied it will format the line objects in the same manner as plot. The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.
Example:

```
z = [0:0.05:5];
plot3 (cos (2*pi*z), sin (2*pi*z), z, ";helix;");
plot3 (z, exp (2i*pi*z), ";complex sinusoid;");
```

See also: [ezplot3], page 384, [plot], page 316.

```
view (azimuth, elevation)
view ([azimuth elevation])
view ([x y z])
view (2)
view (3)
view (hax,...)
[azimuth, elevation] = view ()
```

Query or set the viewpoint for the current axes.
The parameters azimuth and elevation can be given as two arguments or as 2-element vector. The viewpoint can also be specified with Cartesian coordinates $x, y$, and $z$.
The call view (2) sets the viewpoint to azimuth $=0$ and elevation $=90$, which is the default for 2-D graphs.
The call view (3) sets the viewpoint to azimuth $=-37.5$ and elevation $=30$, which is the default for 3-D graphs.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
If no inputs are given, return the current azimuth and elevation.

```
camlookat ()
camlookat (h)
camlookat (handle_list)
camlookat (hax)
```

Move the camera and adjust its properties to look at objects.
When the input is a handle $h$, the camera is set to point toward the center of the bounding box of $h$. The camera's position is adjusted so the bounding box approximately fills the field of view.
This command fixes the camera's viewing direction (camtarget() - campos()), camera up vector (see [camup], page 379) and viewing angle (see [camva], page 379). The camera target (see [camtarget], page 378) and camera position (see [campos], page 376) are changed.
If the argument is a list handle_list, then a single bounding box for all the objects is computed and the camera is then adjusted as above.
If the argument is an axis object hax, then the children of the axis are used as handle_list. When called with no inputs, it uses the current axis (see [gca], page 429).
See also: [camorbit], page 377, [camzoom], page 380, [camroll], page 377.
$P=$ campos ()
campos ([x y z])
mode = campos ("mode")
campos (mode)
campos (ax, ...)
Set or get the camera position.
The default camera position is determined automatically based on the scene. For example, to get the camera position:

```
hf = figure();
peaks()
p = campos ()
p =
-27.394 -35.701 64.079
```

We can then move the camera further up the z -axis:

```
campos ( \(\left.p+\left[\begin{array}{lll}0 & 0 & 10\end{array}\right]\right)\)
campos ()
            \(\Rightarrow\) ans \(=\)
                -27.394 -35.701 74.079
```

Having made that change, the camera position mode is now manual:

```
campos ("mode")
            m manual
```

We can set it back to automatic:

```
campos ("auto")
campos ()
            \(\Rightarrow\) ans \(=\)
                    -27.394 -35.701 64.079
close (hf)
```

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.

See also: [camup], page 379, [camtarget], page 378, [camva], page 379.

```
camorbit (theta, phi)
camorbit (theta, phi, coorsys)
camorbit (theta, phi, coorsys, dir)
camorbit (theta, phi, "data")
camorbit (theta, phi, "data", "z")
camorbit (theta, phi, "data", "x")
camorbit (theta, phi, "data", "y")
camorbit (theta, phi, "data", [x y z])
camorbit (theta, phi, "camera")
camorbit (hax,...)
```

Rotate the camera up/down and left/right around its target.
Move the camera phi degrees up and theta degrees to the right, as if it were in an orbit around its target. Example:

```
sphere ()
camorbit (30, 20)
```

These rotations are centered around the camera target (see [camtarget], page 378). First the camera position is pitched up or down by rotating it phi degrees around an axis orthogonal to both the viewing direction (specifically camtarget() - campos()) and the camera "up vector" (see [camup], page 379). Example:

```
camorbit (0, 20)
```

The second rotation depends on the coordinate system coorsys and direction dir inputs. The default for coorsys is "data". In this case, the camera is yawed left or right by rotating it theta degrees around an axis specified by dir. The default for dir is " z ", corresponding to the vector $[0,0,1]$. Example:
camorbit (30, 0)
When coorsys is set to "camera", the camera is moved left or right by rotating it around an axis parallel to the camera up vector (see [camup], page 379). The input dir should not be specified in this case. Example:

```
camorbit (30, 0, "camera")
```

(Note: the rotation by phi is unaffected by "camera".)
The camorbit command modifies two camera properties: [campos], page 376, and [camup], page 379.
By default, this command affects the current axis; alternatively, an axis can be specified by the optional argument hax.
See also: [camzoom], page 380, [camroll], page 377, [camlookat], page 376.
camroll (theta)
camroll (ax, theta)
Roll the camera.

Roll the camera clockwise by theta degrees. For example, the following command will roll the camera by 30 degrees clockwise (to the right); this will cause the scene to appear to roll by 30 degrees to the left:

```
peaks ()
camroll (30)
```

Roll the camera back:

```
camroll (-30)
```

The following command restores the default camera roll:

## camup ("auto")

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.
See also: [camzoom], page 380, [camorbit], page 377, [camlookat], page 376, [camup], page 379 .
$T$ = camtarget ()
camtarget ([x y z])
mode = camtarget ("mode")
camtarget (mode)
camtarget (ax, ...)
Set or get where the camera is pointed.
The camera target is a point in space where the camera is pointing. Usually, it is determined automatically based on the scene:

```
hf = figure();
sphere (36)
v = camtarget ()
    # v =
        0 0
```

We can turn the camera to point at a new target:

```
camtarget ([\begin{array}{lll}{1}&{1}&{1}\end{array}])
camtarget ()
    m}1\begin{array}{lll}{1}&{1}&{1}
```

Having done so, the camera target mode is manual:

```
camtarget ("mode")
    => manual
```

This means, for example, adding new objects to the scene will not retarget the camera:

```
hold on;
peaks ()
camtarget ()
    #}11
```

We can reset it to be automatic:

```
camtarget ("auto")
camtarget ()
    # 0 0 0.76426
close (hf)
```

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.
See also: [campos], page 376, [camup], page 379, [camva], page 379.

```
up = camup ()
camup ([x y z])
mode = camup ("mode")
camup (mode)
camup (ax,...)
```

Set or get the camera up vector.
By default, the camera is oriented so that "up" corresponds to the positive z-axis:

```
hf = figure ();
sphere (36)
v = camup ()
    # v =
        0 0 1
```

Specifying a new "up vector" rolls the camera and sets the mode to manual:

```
camup ([llll
camup ()
    #}\begin{array}{llll}{1}&{1}&{0}
camup ("mode")
    m manual
```

Modifying the up vector does not modify the camera target (see [camtarget], page 378). Thus, the camera up vector might not be orthogonal to the direction of the camera's view:

```
camup ([\begin{array}{lll}{1}&{2}&{3}\end{array}]
dot (camup (), camtarget () - campos ())
    # 6...
```

A consequence is that "pulling back" on the up vector does not pitch the camera view (as that would require changing the target). Setting the up vector is thus typically used only to roll the camera. A more intuitive command for this purpose is [camroll], page 377 .
Finally, we can reset the up vector to automatic mode:

```
camup ("auto")
camup ()
    \(\Rightarrow \quad 0 \quad 0 \quad 1\)
close (hf)
```

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.
See also: [campos], page 376, [camtarget], page 378, [camva], page 379.
a = camva ()
camva (a)
mode = camva ("mode")
camva (mode)
camva (ax, ...)
Set or get the camera viewing angle.
The camera has a viewing angle which determines how much can be seen. By default this is:

```
hf = figure();
sphere (36)
a = camva ()
    =>a=10.340
```

To get a wider-angle view, we could double the viewing angle. This will also set the mode to manual:

```
camva (2*a)
camva ("mode")
    => manual
```

We can set it back to automatic:

```
camva ("auto")
camva ("mode")
    => auto
camva ()
    a ans = 10.340
close (hf)
```

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.
See also: [campos], page 376, [camtarget], page 378, [camup], page 379.
camzoom (zf)
camzoom (ax, zf)
Zoom the camera in or out.
A value of $z f$ larger than 1 "zooms in" such that the scene appears magnified:

```
hf = figure ();
sphere (36)
camzoom (1.2)
```

A value smaller than 1 "zooms out" so the camera can see more of the scene:
camzoom (0.5)
Technically speaking, zooming affects the "viewing angle". The following command resets to the default zoom:

```
camva ("auto")
close (hf)
```

By default, these commands affect the current axis; alternatively, an axis can be specified by the optional argument ax.
See also: [camroll], page 377, [camorbit], page 377, [camlookat], page 376, [camva], page 379 .

```
slice (x, y, z, v, sx, sy, sz)
slice (x, y, z, v, xi, yi, zi)
slice (v, sx, sy, sz)
slice (v, xi, yi, zi)
slice (..., method)
slice (hax,...)
h = slice (...)
```

Plot slices of 3-D data/scalar fields.
Each element of the 3-dimensional array $v$ represents a scalar value at a location given by the parameters $x, y$, and $z$. The parameters $x, x$, and $z$ are either 3-dimensional arrays of the same size as the array $v$ in the "meshgrid" format or vectors. The parameters xi, etc. respect a similar format to $x$, etc., and they represent the points at which the array vi is interpolated using interp3. The vectors $s x, s y$, and $s z$ contain points of orthogonal slices of the respective axes.
If $x, y, z$ are omitted, they are assumed to be $x=1$ : size $(v, 2), y=1$ : size $(v, 1)$ and $z=1$ : size $(v, 3)$.
method is one of:
"nearest"
Return the nearest neighbor.
"linear" Linear interpolation from nearest neighbors.
"cubic" Cubic interpolation from four nearest neighbors (not implemented yet).
"spline" Cubic spline interpolation-smooth first and second derivatives throughout the curve.

The default method is "linear".
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Examples:

```
[x, y, z] = meshgrid (linspace (-8, 8, 32));
v = sin (sqrt (x.^2 + y.^2 + z.^2)) ./ (sqrt (x.^2 + y.^2 + z.^2));
slice (x, y, z, v, [], 0, []);
[xi, yi] = meshgrid (linspace (-7, 7));
zi = xi + yi;
slice (x, y, z, v, xi, yi, zi);
```

See also: [interp3], page 820, [surface], page 426, [pcolor], page 345.

```
ribbon (y)
```

ribbon ( $x, y$ )
ribbon ( $x, y$, width)
ribbon (hax, ...)
$h=r i b b o n(. .$.

Draw a ribbon plot for the columns of $y$ vs. $x$.

If $x$ is omitted, a vector containing the row numbers is assumed (1:rows (Y)). Alternatively, $x$ can also be a vector with same number of elements as rows of $y$ in which case the same $x$ is used for each column of $y$.
The optional parameter width specifies the width of a single ribbon (default is 0.75). If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the surface objects representing each ribbon.
See also: [surface], page 426, [waterfall], page 383.
shading (type)
shading (hax, type)
Set the shading of patch or surface graphic objects.
Valid arguments for type are
"flat" Single colored patches with invisible edges.
"faceted"
Single colored patches with black edges.
"interp" Colors between patch vertices are interpolated and the patch edges are invisible.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

See also: [fill], page 346, [mesh], page 358, [patch], page 425, [pcolor], page 345, [surf], page 360, [surface], page 426, [hidden], page 360, [lighting], page 371.

```
scatter3 (x, y, z)
```

scatter3 (x, y, z, s)
scatter3 (x, y, z, s, c)
scatter3 (..., style)
scatter3 (. . ., "filled")
scatter3 (..., prop, val)
scatter3 (hax, ...)
$h=\operatorname{scatter} 3(. .$.

Draw a 3-D scatter plot.
A marker is plotted at each point defined by the coordinates in the vectors $x, y$, and z.

The size of the markers is determined by $s$, which can be a scalar or a vector of the same length as $x, y$, and $z$. If $s$ is not given, or is an empty matrix, then a default value of 8 points is used.
The color of the markers is determined by $c$, which can be a string defining a fixed color; a 3 -element vector giving the red, green, and blue components of the color; a vector of the same length as $x$ that gives a scaled index into the current colormap; or an Nx3 matrix defining the RGB color of each marker individually.

The marker to use can be changed with the style argument, that is a string defining a marker in the same manner as the plot command. If no marker is specified it defaults to "०" or circles. If the argument "filled" is given then the markers are filled.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the scatter object representing the points.

```
[x, y, z] = peaks (20);
scatter3 (x(:), y(:), z(:), [], z(:));
```

Programming Note: The full list of properties is documented at Section 15.3.3.9 [Scatter Properties], page 461.
See also: [scatter], page 327, [patch], page 425, [plot], page 316.

```
waterfall (x, y, z)
```

waterfall (z)
waterfall (..., c)
waterfall (..., prop, val, ...)
waterfall (hax,...)
$h=$ waterfall (...)
Plot a 3-D waterfall plot.
A waterfall plot is similar to a meshz plot except only mesh lines for the rows of $z$ (x-values) are shown.
The wireframe mesh is plotted using rectangles. The vertices of the rectangles [x, $y]$ are typically the output of meshgrid. over a 2-D rectangular region in the $x-y$ plane. $z$ determines the height above the plane of each vertex. If only a single $z$ matrix is given, then it is plotted over the meshgrid $x=1$ :columns ( $z$ ), $y=1$ : rows $(z)$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values.
The color of the mesh is computed by linearly scaling the $z$ values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally the color of the mesh can be specified independently of $z$ by supplying a color matrix, $c$.
Any property/value pairs are passed directly to the underlying surface object. The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
See also: [meshz], page 359, [mesh], page 358, [meshc], page 359, [contour], page 330, [surf], page 360, [surface], page 426, [ribbon], page 381, [meshgrid], page 373, [hidden], page 360, [shading], page 382, [colormap], page 870, [caxis], page 350.

### 15.2.2.1 Aspect Ratio

For three-dimensional plots the aspect ratio can be set for data with daspect and for the plot box with pbaspect. See Section 15.2.1.1 [Axis Configuration], page 348, for controlling the x -, y -, and z -limits for plotting.

```
data_aspect_ratio = daspect ()
daspect (data_aspect_ratio)
daspect (mode)
data_aspect_ratio_mode = daspect ("mode")
daspect (hax,...)
Query or set the data aspect ratio of the current axes.
```

The aspect ratio is a normalized 3 -element vector representing the span of the $\mathrm{x}, \mathrm{y}$, and z -axis limits.
daspect (mode)
Set the data aspect ratio mode of the current axes. mode is either "auto" or "manual".
daspect ("mode")
Return the data aspect ratio mode of the current axes.
daspect (hax, ...)
Operate on the axes in handle hax instead of the current axes.
See also: [axis], page 348, [pbaspect], page 384, [xlim], page 351, [ylim], page 351, [zlim], page 351.

```
plot_box_aspect_ratio = pbaspect ()
pbaspect (plot_box_aspect_ratio)
pbaspect (mode)
plot_box_aspect_ratio_mode = pbaspect ("mode")
pbaspect (hax,...)
```

Query or set the plot box aspect ratio of the current axes.
The aspect ratio is a normalized 3 -element vector representing the rendered lengths of the $\mathrm{x}, \mathrm{y}$, and z axes.

```
pbaspect(mode)
```

Set the plot box aspect ratio mode of the current axes. mode is either "auto" or "manual".
pbaspect ("mode")
Return the plot box aspect ratio mode of the current axes.
pbaspect (hax, ...)
Operate on the axes in handle hax instead of the current axes.
See also: [axis], page 348, [daspect], page 383, [xlim], page 351, [ylim], page 351, [zlim], page 351.

### 15.2.2.2 Three-dimensional Function Plotting

ezplot3 ( $f x, f y, f z$ )
ezplot3 (..., dom)
ezplot3 (..., n)
ezplot3 (..., "animate")
ezplot3 (hax, ...)
h = ezplot3 (...)
Plot a parametrically defined curve in three dimensions.
$f x, f y$, and $f z$ are strings, inline functions, or function handles with one argument defining the function. By default the plot is over the domain $0<=t<=2 *$ pi with 500 points.
If dom is a two element vector, it represents the minimum and maximum values of $t$.
$n$ is a scalar defining the number of points to use in plotting the function.
If the "animate" option is given then the plotting is animated in the style of comet3. If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created plot.

```
fx = @(t) cos (t);
fy = @(t) sin (t);
fz = @(t) t;
ezplot3 (fx, fy, fz, [0, 10*pi], 100);
```

See also: [plot3], page 374, [comet3], page 348, [ezplot], page 354, [ezmesh], page 385, [ezsurf], page 386.

```
ezmesh (f)
ezmesh (fx, fy,fz)
ezmesh (..., dom)
ezmesh (..., n)
ezmesh (..., "circ")
ezmesh (hax,...)
h = ezmesh (...)
```

Plot the mesh defined by a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 *$ pi $<=x \mid y<=2 *$ pi with 60 points in each dimension.
If three functions are passed, then plot the parametrically defined function [fx $(s$, $t), f y(s, t), f z(s, t)]$.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the argument "circ" is given, then the function is plotted over a disk centered on the middle of the domain dom.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Example 1: 2-argument function

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezmesh (f, [-3, 3]);
```

Example 2: parametrically defined function

```
fx = @(s,t) cos (s) .* cos (t);
fy =@(s,t) sin (s) .* cos (t);
fz = @(s,t) sin (t);
ezmesh (fx, fy, fz, [-pi, pi, -pi/2, pi/2], 20);
```

See also: [mesh], page 358, [ezmeshc], page 386, [ezplot], page 354, [ezsurf], page 386, [ezsurfc], page 387, [hidden], page 360.

```
ezmeshc (f)
ezmeshc (fx, fy,fz)
ezmeshc (..., dom)
ezmeshc (..., n)
ezmeshc (..., "circ")
ezmeshc (hax,...)
h = ezmeshc (...)
```

Plot the mesh and contour lines defined by a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 * \mathrm{pi}<=\mathrm{x} \mid \mathrm{y}<=2 * \mathrm{pi}$ with 60 points in each dimension.
If three functions are passed, then plot the parametrically defined function [fx $(s$, $t), f y(s, t), f z(s, t)]$.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the argument "circ" is given, then the function is plotted over a disk centered on the middle of the domain dom.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a 2 -element vector with a graphics handle for the created mesh plot and a second handle for the created contour plot.
Example: 2-argument function

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezmeshc (f, [-3, 3]);
```

See also: [meshc], page 359, [ezmesh], page 385, [ezplot], page 354, [ezsurf], page 386, [ezsurfc], page 387, [hidden], page 360.

```
ezsurf (f)
ezsurf (fx, fy,fz)
ezsurf (..., dom)
ezsurf (..., n)
ezsurf (..., "circ")
ezsurf (hax,...)
h = ezsurf (...)
```

Plot the surface defined by a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 * \mathrm{pi}<=\mathrm{x} \mid \mathrm{y}<=2 * \mathrm{pi}$ with 60 points in each dimension.
If three functions are passed, then plot the parametrically defined function [fx(s, $t), f y(s, t), f z(s, t)]$.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the argument "circ" is given, then the function is plotted over a disk centered on the middle of the domain dom.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Example 1: 2-argument function

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezsurf (f, [-3, 3]);
```

Example 2: parametrically defined function

```
fx = @(s,t) cos (s) .* cos (t);
fy = @(s,t) sin (s) .* cos (t);
fz = @(s,t) sin (t);
ezsurf (fx, fy, fz, [-pi, pi, -pi/2, pi/2], 20);
```

See also: [surf], page 360, [ezsurfc], page 387, [ezplot], page 354, [ezmesh], page 385, [ezmeshc], page 386, [shading], page 382.

```
ezsurfc (f)
ezsurfc (fx, fy,fz)
ezsurfc (..., dom)
ezsurfc (..., n)
ezsurfc (..., "circ")
ezsurfc (hax,...)
h = ezsurfc (...)
```

Plot the surface and contour lines defined by a function.
$f$ is a string, inline function, or function handle with two arguments defining the function. By default the plot is over the meshed domain $-2 * \mathrm{pi}<=\mathrm{x} \mid \mathrm{y}<=2 * \mathrm{pi}$ with 60 points in each dimension.
If three functions are passed, then plot the parametrically defined function [fx $(s$, $t), f y(s, t), f z(s, t)]$.
If dom is a two element vector, it represents the minimum and maximum values of both $x$ and $y$. If dom is a four element vector, then the minimum and maximum values are [xmin xmax ymin ymax].
$n$ is a scalar defining the number of points to use in each dimension.
If the argument "circ" is given, then the function is plotted over a disk centered on the middle of the domain dom.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a 2 -element vector with a graphics handle for the created surface plot and a second handle for the created contour plot.
Example:

```
f = @(x,y) sqrt (abs (x .* y)) ./ (1 + x.^2 + y.^2);
ezsurfc (f, [-3, 3]);
```

See also: [surfc], page 361, [ezsurf], page 386, [ezplot], page 354, [ezmesh], page 385, [ezmeshc], page 386, [shading], page 382.

### 15.2.2.3 Three-dimensional Geometric Shapes

cylinder
cylinder ( $r$ )
cylinder ( $r, n$ )
cylinder (hax, ...)
[x, y, z] = cylinder (...)
Plot a 3-D unit cylinder.
The optional input $r$ is a vector specifying the radius along the unit $z$-axis. The default is [11] indicating radius 1 at $\mathrm{Z}==0$ and at $\mathrm{Z}==1$.
The optional input $n$ determines the number of faces around the circumference of the cylinder. The default value is 20 .
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If outputs are requested cylinder returns three matrices in meshgrid format, such that surf ( $x, y, z$ ) generates a unit cylinder.
Example:

```
[x, y, z] = cylinder (10:-1:0, 50);
surf (x, y, z);
title ("a cone");
```

See also: [ellipsoid], page 389, [rectangle], page 356, [sphere], page 388.
sphere ()
sphere ( $n$ )
sphere (hax, ...)
[x, y, z] = sphere (...)
Plot a 3-D unit sphere.
The optional input $n$ determines the number of faces around the circumference of the sphere. The default value is 20 .
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If outputs are requested sphere returns three matrices in meshgrid format such that surf $(x, y, z)$ generates a unit sphere.

Example:

```
[x, y, z] = sphere (40);
surf (3*x, 3*y, 3*z);
axis equal;
title ("sphere of radius 3");
```

See also: [cylinder], page 388, [ellipsoid], page 389, [rectangle], page 356.
ellipsoid (xc, yc, zc, xr, yr, zr, n)
ellipsoid (..., n)
ellipsoid (hax, ...)
[x, y, z] = ellipsoid (...)
Plot a 3-D ellipsoid.
The inputs $x c, y c, z c$ specify the center of the ellipsoid. The inputs $x r, y r, z r$ specify the semi-major axis lengths.
The optional input $n$ determines the number of faces around the circumference of the cylinder. The default value is 20 .
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
If outputs are requested ellipsoid returns three matrices in meshgrid format, such that surf ( $x, y, z$ ) generates the ellipsoid.

See also: [cylinder], page 388, [rectangle], page 356, [sphere], page 388.

### 15.2.3 Plot Annotations

You can add titles, axis labels, legends, and arbitrary text to an existing plot. For example:

```
x = -10:0.1:10;
plot (x, sin (x));
title ("sin(x) for x = -10:0.1:10");
xlabel ("x");
ylabel ("sin (x)");
text (pi, 0.7, "arbitrary text");
legend ("sin (x)");
```

The functions grid and box may also be used to add grid and border lines to the plot. By default, the grid is off and the border lines are on.

Finally, arrows, text and rectangular or elliptic boxes can be added to highlight parts of a plot using the annotation function. Those objects are drawn in an invisible axes, on top of every other axes.

```
title (string)
title (string, prop, val, ...)
title (hax,...)
h = title (...)
```

Specify the string used as a title for the current axis.
An optional list of property/value pairs can be used to change the appearance of the created title text object.

If the first argument hax is an axes or legend handle, then add a title to this object, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created text object.
See also: [xlabel], page 393, [ylabel], page 393, [zlabel], page 393, [text], page 392.

```
legend ()
legend command
legend (str1, str2,...)
legend (charmat)
legend ({cellstr})
legend (..., property, value,...)
legend (hobjs,...)
legend ("command")
legend (hax,...)
legend (hleg,...)
hleg, hplt = legend (...)
```

Display a legend for the current axes using the specified strings as labels.
Legend entries may be specified as individual character string arguments, a character array, or a cell array of character strings. When label names might be confused with legend properties, or command arguments, the labels should be protected by specifying them as a cell array of strings.
If the first argument hax is an axes handle, then add a legend to this axes, rather than the current axes returned by gca.
If the first argument hleg is a legend handle, then operate on this legend rather than the legend of the current axes.
Legend labels are associated with the axes' children; The first label is assigned to the first object that was plotted in the axes, the second label to the next object plotted, etc. To label specific data objects, without labeling all objects, provide their graphic handles in the input hobjs.
The following customizations are available using command:

$$
\begin{array}{ll}
\text { "show" } & \text { Show legend on the plot } \\
\text { "hide" } & \text { Hide legend on the plot } \\
\text { "toggle" } & \text { Toggle between "hide" and "show" } \\
\text { "boxon" } & \text { Show a box around legend (default) } \\
\text { "boxoff" } & \text { Hide the box around legend } \\
\text { "right" } & \text { Place label text to the right of the keys (default) } \\
\text { "left" } & \text { Place label text to the left of the keys } \\
\text { "off" } & \text { Delete the legend object }
\end{array}
$$

The legend function creates a graphics object which has various properties that can be manipulated with get/set. Alternatively, properties can be set directly when calling legend by including property/value pairs. If using this calling form, the
labels must be specified as a cell array of strings. Graphics object properties are documented in detail at Section 15.3.3 [Graphics Object Properties], page 433.

Following is a subset of supported legend properties:

```
autoupdate: "off" | {"on"}
```

Control whether the number of legend items is updated automatically when objects are added to (or deleted from) the peer axes. For example:

```
## Create a single plot with its legend.
figure ();
plot (1:10);
legend ("Slope 1");
## Add another plot and specify its displayname so that
## the legend is correctly updated.
hold on;
plot ((1:10) * 2, "displayname", "Slope 2");
## Stop automatic updates for further plots.
legend ("autoupdate", "off");
plot ((1:10) * 3);
```

box: "off" | \{"on"\}
Control whether the legend has a surrounding box.

```
location: "best" | "bestoutside" |
    "east" | "eastoutside" | "none" | "north" | {"northeast"}
    | "northeastoutside" | "northoutside" | "northwest"|
    "northwestoutside" | "south" | "southeast" | "southeastoutside"
    | "southoutside" | "southwest" | "southwestoutside" | "west" |
    "westoutside" Control the location of the legend.
```

numcolumns: scalar interger, def. 1
Control the number of columns used in the layout of the legend items.
For example:
figure ();
plot (rand (30));
legend ("numcolumns", 3);

Setting numcolumns also forces the numcolumnsmode property to be set to "manual".
orientation: "horizontal" | \{"vertical"\}
Control whether the legend items are arranged vertically (column-wise) or horizontally (row-wise).
string: string | cell array of strings
List of labels for the legend items. For example:

```
figure ();
plot (rand (20));
## Let legend choose names automatically
hl = legend ();
## Selectively change some names
str = get (hl, "string");
str(1:5:end) = "Garbage";
set (hl, "string", str);
```

textcolor: colorspec, def. [0 0 0]
Control the color of the text strings for legend item.
The full list of supported legend specific properties can be found at Section 15.3.3.4 [Legend Properties], page 449.
A legend is implemented as an additional axes object with the tag property set to "legend". Properties of the legend object may be manipulated directly by using set.
The optional output value hleg is a handle to the legend object.
Implementation Note: The legend label text is either provided in the call to legend or is taken from the DisplayName property of the graphics objects. Only data objects, such as line, patch, and surface, have this property whereas axes, figures, etc. do not so they are never present in a legend. If no labels or DisplayName properties are available, then the label text is simply "data1", "data2", ..., "dataN".
The legend FontSize property is initially set to $90 \%$ of the axes FontSize to which it is attached. Use set to override this if necessary.

```
text (x, y, string)
text (x, y, z, string)
text (..., prop, val, ...)
text (hax,...)
h = text (...)
```

Create a text object with text string at position $x, y,(z)$ on the current axes.
Multiple locations can be specified if $x, y,(z)$ are vectors. Multiple strings can be specified with a character matrix or a cell array of strings.
Optional property/value pairs may be used to control the appearance of the text.
If the first argument hax is an axes handle, then add text to this axes, rather than the current axes returned by gca.
The optional return value $h$ is a vector of graphics handles to the created text objects.
Example 1: multi-line text via 3 different methods

```
text (0.5, 0.8, {"Line 1", "Line 2"})
text (0.5, 0.6, ["Line 1"; "Line 2"])
text (0.5, 0.4, "Line 1\nLine 2")
```

Example 2 : text at multiple locations

```
text ([0.2, 0.2], [0.8, 0.6], "Same text at two locations")
text ([0.4, 0.4], [0.8, 0.6], {"Point 1 Text", "Point 2 text"})
text ([0.6, 0.6], [0.8, 0.6], {{"Point 1 Line 1", "Point 1 Line 2},
    "Point 2 text"})
```

Example 2 : adjust appearance using text properties

```
ht = text (0.5, 0.5, "Hello World", "fontsize", 20);
set (ht, "color", "red");
```

Programming Notes: The full list of properties is documented at Section 15.3.3.6 [Text Properties], page 453.
Any numeric entries in a cell array will be converted to text using sprintf (" $\% \mathrm{~g}$ "). For more precise control of the appearance convert any numeric entries to strings using num2str, sprintf, etc., before calling text.

See also: [gtext], page 420, [title], page 389, [xlabel], page 393, [ylabel], page 393, [zlabel], page 393.
xlabel (string)
xlabel (string, property, val, ...)
xlabel (hax, ...)
$h=x l a b e l(. .$.
Specify the string used to label the x-axis of the current axis.
An optional list of property/value pairs can be used to change the properties of the created text label.
The full list of text object properties is documented at Section 15.3.3.6 [Text Properties], page 453.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created text object.
See also: [ylabel], page 393, [zlabel], page 393, [datetick], page 952, [title], page 389, [text], page 392.
clabel ( $c, h$ )
clabel ( $c, h, v$ )
clabel (c, h, "manual")
clabel (c)
clabel (..., prop, val, ...)
$h=c l a b e l(. .$.
Add labels to the contours of a contour plot.
The contour levels are specified by the contour matrix $c$ which is returned by contour, contourc, contourf, and contour3. Contour labels are rotated to match the local line orientation and centered on the line. The position of labels along the contour line is chosen randomly.
If the argument $h$ is a handle to a contour group object, then label this plot rather than the one in the current axes returned by gca.
By default, all contours are labeled. However, the contours to label can be specified by the vector $v$. If the "manual" argument is given then the contours to label can be selected with the mouse.
Additional property/value pairs that are valid properties of text objects can be given and are passed to the underlying text objects. Moreover, the contour group prop-
erty "LabelSpacing" is available which determines the spacing between labels on a contour to be specified. The default is 144 points, or 2 inches.
The optional return value $h$ is a vector of graphics handles to the text objects representing each label. The "userdata" property of the text objects contains the numerical value of the contour label.
The full list of text object properties is documented at Section 15.3.3.6 [Text Properties], page 453.

```
[c, h] = contour (peaks (), -4 : 6);
clabel (c, h, -4:2:6, "fontsize", 12);
```

See also: [contour], page 330, [contourf], page 331, [contour3], page 332, [meshc], page 359, [surfc], page 361, [text], page 392.

```
box
```

box on
box off
box (hax, ...)

Control display of the axes border.
The argument may be either "on" or "off". If it is omitted, the current box state is toggled.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.

See also: [axis], page 348, [grid], page 394.

## grid

grid on
grid off
grid minor
grid minor on
grid minor off
grid (hax, ...)
Control the display of plot grid lines.
The function state input may be either "on" or "off". If it is omitted, the current grid state is toggled.
When the first argument is "minor" all subsequent commands modify the minor grid rather than the major grid.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
To control the grid lines for an individual axes use the set function. For example:

```
set (gca, "ygrid", "on");
```

See also: [axis], page 348, [box], page 394.

```
colorbar
colorbar (..., loc)
colorbar (delete_option)
```

```
colorbar (hcb,...)
colorbar (hax,...)
colorbar (..., "peer", hax, ...)
colorbar (..., "location", loc,...)
colorbar (..., prop, val, ...)
h = colorbar (...)
```

Add a colorbar to the current axes.
A colorbar displays the current colormap along with numerical rulings so that the color scale can be interpreted.
The optional input loc determines the location of the colorbar. If present, it must be the last argument to colorbar. Valid values for loc are
"EastOutside"
Place the colorbar outside the plot to the right. This is the default.
"East" Place the colorbar inside the plot to the right.
"WestOutside"
Place the colorbar outside the plot to the left.
"West" Place the colorbar inside the plot to the left.

```
"NorthOutside"
```

Place the colorbar above the plot.
"North" Place the colorbar at the top of the plot.
"SouthOutside"
Place the colorbar under the plot.
"South" Place the colorbar at the bottom of the plot.
To remove a colorbar from a plot use any one of the following keywords for the delete_option: "off", "delete", "hide".
If the first argument hax is an axes handle, then the colorbar is added to this axes, rather than the current axes returned by gca. Alternatively, If the argument "peer" is given, then the following argument is treated as the axes handle in which to add the colorbar. The "peer" calling syntax may be removed in the future and is not recommended.

If the first argument hcb is a handle to a colorbar object, then operate on this colorbar directly.
Additional property/value pairs are passed directly to the underlying axes object. The full list of properties is documented at Section 15.3.3.3 [Axes Properties], page 441.
The optional return value $h$ is a graphics handle to the created colorbar object.
Implementation Note: A colorbar is created as an additional axes object with the "tag" property set to "colorbar". The created object has the extra property "location" which controls the positioning of the colorbar.
See also: [colormap], page 870 .

```
annotation (type)
annotation ("line", x, y)
annotation ("arrow", x, y)
annotation ("doublearrow", x, y)
annotation ("textarrow", x, y)
annotation ("textbox", pos)
annotation ("rectangle", pos)
annotation ("ellipse", pos)
annotation (..., prop, val)
annotation (hf, ...)
h = annotation (...)
```

Draw annotations to emphasize parts of a figure.
You may build a default annotation by specifying only the type of the annotation.
Otherwise you can select the type of annotation and then set its position using either $x$ and $y$ coordinates for line-based annotations or a position vector pos for others. In either case, coordinates are interpreted using the "units" property of the annotation object. The default is "normalized", which means the lower left hand corner of the figure has coordinates '[00]' and the upper right hand corner '[11]'.
If the first argument $h f$ is a figure handle, then plot into this figure, rather than the current figure returned by gcf.
Further arguments can be provided in the form of prop/val pairs to customize the annotation appearance.
The optional return value $h$ is a graphics handle to the created annotation object. This can be used with the set function to customize an existing annotation object.
All annotation objects share two properties:

- "units": the units in which coordinates are interpreted.

Its value may be one of "centimeters" | "characters" | "inches" | "\{normalized\}" | "pixels" | "points".

- "position": a four-element vector [x0 y0 width height].

The vector specifies the coordinates ( $\mathrm{x} 0, \mathrm{y} 0$ ) of the origin of the annotation object, its width, and its height. The width and height may be negative, depending on the orientation of the object.

Valid annotation types and their specific properties are described below:
"line" Constructs a line. $x$ and $y$ must be two-element vectors specifying the $x$ and $y$ coordinates of the two ends of the line.
The line can be customized using "linewidth", "linestyle", and "color" properties the same way as for line objects.
"arrow" Construct an arrow. The second point in vectors $x$ and $y$ specifies the arrowhead coordinates.
Besides line properties, the arrowhead can be customized using "headlength", "headwidth", and "headstyle" properties. Supported values for "headstyle" property are: ["diamond" | "ellipse" | "plain" | "rectangle" | "vback1" | "\{vback2\}" | "vback3"]
"doublearrow"
Construct a double arrow. Vectors $x$ and $y$ specify the arrowhead coordinates.

The line and the arrowhead can be customized as for arrow annotations, but some property names are duplicated: "head1length"/"head2length", "head1width"/"head2width", etc. The index 1 marks the properties of the arrowhead at the first point in $x$ and $y$ coordinates.
"textarrow"
Construct an arrow with a text label at the opposite end from the arrowhead.

Use the "string" property to change the text string. The line and the arrowhead can be customized as for arrow annotations, and the text can be customized using the same properties as text graphics objects. Note, however, that some text property names are prefixed with "text" to distinguish them from arrow properties: "textbackgroundcolor", "textcolor", "textedgecolor", "textlinewidth", "textmargin", "textrotation".
"textbox"
Construct a box with text inside. pos specifies the "position" property of the annotation.

Use the "string" property to change the text string. You may use "backgroundcolor", "edgecolor", "linestyle", and "linewidth" properties to customize the box background color and edge appearance. A limited set of text objects properties are also available; Besides "font..." properties, you may also use "horizontalalignment" and "verticalalignment" to position the text inside the box.
Finally, the "fitboxtotext" property controls the actual extent of the box. If "on" (the default) the box limits are fitted to the text extent.
"rectangle"
Construct a rectangle. pos specifies the "position" property of the annotation.
You may use "facecolor", "color", "linestyle", and "linewidth" properties to customize the rectangle background color and edge appearance.
"ellipse"
Construct an ellipse. pos specifies the "position" property of the annotation.

See "rectangle" annotations for customization.
See also: [xlabel], page 393, [ylabel], page 393, [zlabel], page 393, [title], page 389, [text], page 392, [gtext], page 420, [legend], page 390, [colorbar], page 394.

### 15.2.4 Multiple Plots on One Page

Octave can display more than one plot in a single figure. The simplest way to do this is to use the subplot function to divide the plot area into a series of subplot windows that are indexed by an integer. For example,

```
subplot (2, 1, 1)
fplot (@sin, [-10, 10]);
subplot (2, 1, 2)
fplot (@cos, [-10, 10]);
```

creates a figure with two separate axes, one displaying a sine wave and the other a cosine wave. The first call to subplot divides the figure into two plotting areas (two rows and one column) and makes the first plot area active. The grid of plot areas created by subplot is numbered in row-major order (left to right, top to bottom). After plotting a sine wave, the next call to subplot activates the second subplot area, but does not re-partition the figure.

```
subplot (rows, cols, index)
subplot (rows, cols, index, hax)
subplot (rcn)
subplot (hax)
subplot (..., "align")
subplot (..., "replace")
subplot ("position", pos)
subplot (..., prop, val,...)
hax = subplot (...)
```

Set up a plot grid with rows by cols subwindows and set the current axes for plotting (gca) to the location given by index.
If an axes handle hax is provided after the (rows, cols, index) arguments, the corresponding axes is turned into a subplot.
If only one numeric argument is supplied, then it must be a three digit value specifying the number of rows in digit 1 , the number of columns in digit 2, and the plot index in digit 3.
The plot index runs row-wise; First, all columns in a row are numbered and then the next row is filled.

For example, a plot with 2 x 3 grid will have plot indices running as follows:

| 1 | 2 | 3 |
| :--- | :--- | :--- |
| 4 | 5 | 6 |

index may also be a vector. In this case, the new axes will enclose the grid locations specified. The first demo illustrates this:

```
demo ("subplot", 1)
```

The index of the subplot to make active may also be specified by its axes handle, hax, returned from a previous subplot command.

If the option "align" is given then the plot boxes of the subwindows will align, but this may leave no room for axes tick marks or labels.
If the option "replace" is given then the subplot axes will be reset, rather than just switching the current axes for plotting to the requested subplot.
The "position" property can be used to exactly position the subplot axes within the current figure. The option pos is a 4 -element vector [ $\mathrm{x}, \mathrm{y}$, width, height] that determines the location and size of the axes. The values in pos are normalized in the range $[0,1]$.
Any property/value pairs are passed directly to the underlying axes object. The full list of properties is documented at Section 15.3.3.3 [Axes Properties], page 441.
Any previously existing axes that would be (partly) covered by the newly created axes are deleted.
If the output hax is requested, subplot returns the axes handle for the subplot. This is useful for modifying the properties of a subplot using set.
Under some circumstances, subplot might not be able to identify axes that it could re-use and might replace them. If subplot axes should be referenced repeatedly, consider creating and storing their axes handles beforehand instead of calling subplot repeatedly for the same position.
Example:

```
x = 1:10;
y = rand (16, 10);
for i_plot = 1:4
    hax(i_plot) = subplot (2, 2, i_plot);
    hold (hax(i_plot), "on");
    grid (hax(i_plot), "on");
endfor
for i_loop = 1:2
    for i_plot = 1:4
        iy = (i_loop - 1)*4 + i_plot;
        plotyy (hax(i_plot), x,y(iy,:), x,y(iy+1,:));
    endfor
    endfor
```

See also: [axes], page 423, [plot], page 316, [gca], page 429, [set], page 430.

### 15.2.5 Multiple Plot Windows

You can open multiple plot windows using the figure function. For example,

```
figure (1);
fplot (@sin, [-10, 10]);
figure (2);
fplot (@cos, [-10, 10]);
```

creates two figures, with the first displaying a sine wave and the second a cosine wave. Figure numbers must be positive integers.
figure
figure $n$

```
figure (n)
figure (..., "property", value,...)
h = figure (...)
```

Create a new figure window for plotting.
If no arguments are specified, a new figure with the next available number is created.
If called with an integer $n$, and no such numbered figure exists, then a new figure with the specified number is created. If the figure already exists then it is made visible and becomes the current figure for plotting.
Multiple property-value pairs may be specified for the figure object, but they must appear in pairs.
The optional return value $h$ is a graphics handle to the created figure object.
Programming Note: The full list of properties is documented at Section 15.3.3.2 [Figure Properties], page 434.
See also: [axes], page 423, [gcf], page 428, [shg], page 404, [clf], page 404, [close], page 405.

### 15.2.6 Manipulation of Plot Objects

```
pan
pan on
pan off
pan xon
pan yon
pan (hfig, option)
```

Control the interactive panning mode of a figure in the GUI.
Given the option "on" or "off", set the interactive pan mode on or off.
With no arguments, toggle the current pan mode on or off.
Given the option "xon" or "yon", enable pan mode for the x or y axis only.
If the first argument hfig is a figure, then operate on the given figure rather than the current figure as returned by gcf.
See also: [rotate3d], page 400, [zoom], page 401.

```
rotate (h, direction, alpha)
```

rotate (..., origin)

Rotate the plot object $h$ through alpha degrees around the line with direction direction and origin origin.
The default value of origin is the center of the axes object that is the parent of $h$.
If $h$ is a vector of handles, they must all have the same parent axes object.
Graphics objects that may be rotated are lines, surfaces, patches, and images.
rotate3d
rotate3d on
rotate3d off
rotate3d (hfig, option)
Control the interactive 3-D rotation mode of a figure in the GUI.

Given the option "on" or "off", set the interactive rotate mode on or off.
With no arguments, toggle the current rotate mode on or off.
If the first argument hfig is a figure, then operate on the given figure rather than the current figure as returned by gcf.
See also: [pan], page 400, [zoom], page 401.

```
zoom
zoom (factor)
zoom on
zoom off
zoom xon
zoom yon
zoom out
zoom reset
zoom (hfig, option)
```

Zoom the current axes object or control the interactive zoom mode of a figure in the GUI.

Given a numeric argument greater than zero, zoom by the given factor. If the zoom factor is greater than one, zoom in on the plot. If the factor is less than one, zoom out. If the zoom factor is a two- or three-element vector, then the elements specify the zoom factors for the $\mathrm{x}, \mathrm{y}$, and z axes respectively.
Given the option "on" or "off", set the interactive zoom mode on or off.
With no arguments, toggle the current zoom mode on or off.
Given the option "xon" or "yon", enable zoom mode for the x or y -axis only.
Given the option "out", zoom to the initial zoom setting.
Given the option "reset", store the current zoom setting so that zoom out will return to this zoom level.
If the first argument hfig is a figure, then operate on the given figure rather than the current figure as returned by gcf.
See also: [pan], page 400, [rotate3d], page 400.

### 15.2.7 Manipulation of Plot Windows

By default, Octave refreshes the plot window when a prompt is printed, or when waiting for input. The drawnow function is used to cause a plot window to be updated.
drawnow ()
drawnow ("expose")
drawnow (term, file, debug_file)
Update figure windows and their children.
The event queue is flushed and any callbacks generated are executed.
With the optional argument "expose", only graphic objects are updated and no other events or callbacks are processed.
The third calling form of drawnow is for debugging and is undocumented.
See also: [refresh], page 402.

Only figures that are modified will be updated. The refresh function can also be used to cause an update of the current figure, even if it is not modified.

```
refresh ()
refresh (h)
```

Refresh a figure, forcing it to be redrawn.
When called without an argument the current figure is redrawn. Otherwise, the figure with graphic handle $h$ is redrawn.
See also: [drawnow], page 401.
Normally, high-level plot functions like plot or mesh call newplot to determine whether the state of the target axes should be initialized (the default) or if subsequent plots should be drawn on top of previous ones. To have two plots drawn over one another, use the hold function or manually change the axes [nextplot], page 444, property. For example,

```
hold on;
x = -10:0.1:10;
plot (x, sin (x));
plot (x, cos (x));
hold off;
```

displays sine and cosine waves on the same axes. If the hold state is off, consecutive plotting commands like this will only display the last plot.

```
newplot ()
newplot (hfig)
newplot (hax)
hax = newplot (...)
```

Prepare graphics engine to produce a new plot.
This function is called at the beginning of all high-level plotting functions. It is not normally required in user programs. newplot queries the "NextPlot" field of the current figure and axes to determine what to do.

## Figure NextPlot Action

"new" Create a new figure and make it the current figure.
"add" (default) Add new graphic objects to the current figure.
"replacechildren" Delete child objects whose HandleVisibility is set to "on". Set NextPlot property to "add". This typically clears a figure, but leaves in place hidden objects such as menubars. This is equivalent to clf.
"replace" Delete all child objects of the figure and reset all figure properties to their defaults. However, the following four properties are not reset: Position, Units, PaperPosition, PaperUnits. This is equivalent to clf reset.

## Axes NextPlot Action

"add"
Add new graphic objects to the current axes. This is equivalent to hold on.
"replacechildren" Delete child objects whose HandleVisibility is set to "on", but leave axes properties unmodified. This typically clears a plot, but preserves special settings such as log scaling for axes. This is equivalent to cla.
"replace" (default) Delete all child objects of the axes and reset all axes properties to their defaults. However, the following properties are not reset: Position, Units. This is equivalent to cla reset.
If the optional input hfig or hax is given then prepare the specified figure or axes rather than the current figure and axes.
The optional return value hax is a graphics handle to the created axes object (not figure).
Caution: Calling newplot may change the current figure and current axes.
hold
hold on
hold off
hold (hax, ...)
Toggle or set the "hold" state of the plotting engine which determines whether new graphic objects are added to the plot or replace the existing objects.
hold on Retain plot data and settings so that subsequent plot commands are displayed on a single graph. Line color and line style are advanced for each new plot added.
hold all (deprecated)
Equivalent to hold on.
hold off Restore default graphics settings which clear the graph and reset axes properties before each new plot command. (default).
hold Toggle the current hold state.
When given the additional argument hax, the hold state is modified for this axes rather than the current axes returned by gca.
To query the current hold state use the ishold function.
See also: [ishold], page 403, [cla], page 404, [clf], page 404, [newplot], page 402.
ishold
ishold (hax)
ishold (hfig)
Return true if the next plot will be added to the current plot, or false if the plot device will be cleared before drawing the next plot.
If the first argument is an axes handle hax or figure handle hfig then operate on this plot rather than the current one.
See also: [hold], page 403, [newplot], page 402.

To clear the current figure, call the clf function. To clear the current axis, call the cla function. To bring the current figure to the top of the window stack, call the shg function. To delete a graphics object, call delete on its index. To close the figure window, call the close function.

## clf

clf reset
clf (hfig)
clf (hfig, "reset")
h = clf (...)
Clear the current figure window.
clf operates by deleting child graphics objects with visible handles (HandleVisibility = "on").
If the optional argument "reset" is specified, delete all child objects including those with hidden handles and reset all figure properties to their defaults. However, the following properties are not reset: Position, Units, PaperPosition, PaperUnits.
If the first argument hfig is a figure handle, then operate on this figure rather than the current figure returned by gcf.
The optional return value $h$ is the graphics handle of the figure window that was cleared.
See also: [cla], page 404, [close], page 405, [delete], page 405, [reset], page 490.
cla
cla reset
cla (hax)
cla (hax, "reset")
Clear the current or specified (hax) axes object.
cla operates by deleting child graphic objects with visible handles (HandleVisibility $=$ "on"). This typically clears the axes of any visual objects, but leaves in place axes limits, tick marks and labels, camera view, etc. In addition, the automatic coloring and styling of lines is reset by changing the axes properties ColorOrderIndex, LinestyleOrderIndex to 1.
If the optional argument "reset" is specified, delete all child objects, including those with hidden handles, and reset all axes properties to their defaults. However, the following properties are not reset: Position, Units.
If the first argument hax is an axes handle, then operate on this axes rather than the current axes returned by gca.
See also: [clf], page 404, [delete], page 405, [reset], page 490.
shg
Show the graph window.
This function makes the current figure visible, and places it on top of of all other plot windows.
Programming Note: shg is equivalent to figure (gcf) assuming that a current figure exists.
See also: [figure], page 399, [drawnow], page 401, [gcf], page 428.

```
delete (file)
delete (file1, file2,...)
delete (handle)
```

Delete the named file or graphics handle.
file may contain globbing patterns such as ' $*$ '. Multiple files to be deleted may be specified in the same function call.
handle may be a scalar or vector of graphic handles to delete.
Programming Note: Deleting graphics objects is the proper way to remove features from a plot without clearing the entire figure.

See also: [clf], page 404, [cla], page 404, [unlink], page 954, [rmdir], page 955.

```
close
```

close (h)
close figname
close all
close all hidden
close all force

Close figure window(s).
When called with no arguments, close the current figure. This is equivalent to close (gcf). If the input $h$ is a graphic handle, or vector of graphics handles, then close each figure in $h$. The figure to close may also be specified by name figname which is matched against the "Name" property of all figures.

If the argument "all" is given then all figures with visible handles (HandleVisibility = "on") are closed.
If the additional argument "hidden" is given then all figures, including hidden ones, are closed.

If the additional argument "force" is given then figures are closed even when "closerequestfcn" has been altered to prevent closing the window.
Implementation Note: close operates by making the handle $h$ the current figure, and then calling the function specified by the "closerequestfcn" property of the figure. By default, the function closereq is used. It is possible that the function invoked will delay or abort removing the figure. To remove a figure without executing any callback functions use delete. When writing a callback function to close a window do not use close to avoid recursion.

See also: [closereq], page 405, [delete], page 405.
closereq ()
Close the current figure and delete all graphics objects associated with it.
By default, the "closerequestfcn" property of a new plot figure points to this function.

See also: [close], page 405, [delete], page 405.

### 15.2.8 Use of the interpreter Property

text (such as titles, labels, legend item) and axes objects feature a [interpreter], page 454, and [ticklabelinterpreter], page 446, property respectively. It determines the manner in which special control sequences in the text are rendered.

The interpreter property can take three values: "none", "tex", "latex".

### 15.2.8.1 none interpreter

If the interpreter is set to "none" then no special rendering occurs-the displayed text is a verbatim copy of the specified text.

### 15.2.8.2 tex interpreter

The "tex" interpreter implements a subset of $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ functionality when rendering text. This allows the insertion of special glyphs such as Greek characters or mathematical symbols. Special characters are inserted by using a backslash ( $\backslash$ ) character followed by a code, as shown in Table 15.1.

Besides special glyphs, the formatting of the text can be changed within the string by using the codes

| \bf | Bold font |
| :--- | :--- |
| \it | Italic font |
| \sl | Oblique Font |
| \rm | Normal font |

These codes may be used in conjunction with the $\{$ and $\}$ characters to limit the change to a part of the string. For example,

```
xlabel ('{\bf H} = a {\bf V}')
```

where the character 'a' will not appear in bold font. Note that to avoid having Octave interpret the backslash character in the strings, the strings themselves should be in single quotes.

It is also possible to change the fontname and size within the text

$$
\begin{array}{ll}
\begin{array}{l}
\text { \fontname\{fontname\} } \\
\backslash \text { fontsize\{size }\}
\end{array} & \begin{array}{l}
\text { Specify the font to use } \\
\text { Specify the size of the font to use }
\end{array}
\end{array}
$$

The color of the text may also be changed inline using either a string (e.g., "red") or numerically with a Red-Green-Blue (RGB) specification (e.g., [100], also red).

$$
\begin{array}{ll}
\backslash \text { color }\{\text { color }\} & \text { Specify the color as a string } \\
\backslash \text { color }[\operatorname{rgb}]\{R & G B\}
\end{array} \quad \text { Specify the color numerically }
$$

Finally, superscripting and subscripting can be controlled with the '~' and ' _' characters. If the ' ${ }^{\prime}$ ' or ' $\_$' is followed by a \{ character, then all of the block surrounded by the $\}$ pair is superscripted or subscripted. Without the $\}$ pair, only the character immediately following the ' ${ }^{\prime \prime}$ ' or ' $\quad$ ' is changed.

Greek Lowercase Letters

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ alpha | $\alpha$ | $\backslash$ beta | $\beta$ | \gamma | $\gamma$ |
| $\backslash$ delta | $\delta$ | $\backslash$ epsilon | $\epsilon$ | \zeta | $\zeta$ |
| $\backslash$ eta | $\eta$ | $\backslash$ theta | $\theta$ | \vartheta | $\vartheta$ |
| $\backslash$ iota | $\iota$ | $\backslash$ kappa | $\kappa$ | $\backslash$ lambda | $\lambda$ |
| $\backslash$ mu | $\mu$ | $\backslash$ nu | $\nu$ | $\backslash$ xi | $\xi$ |
| $\backslash$ o | $o$ | $\backslash$ pi | $\pi$ | $\backslash$ varpi | $\varpi$ |
| $\backslash$ rho | $\rho$ | $\backslash$ sigma | $\sigma$ | $\backslash$ varsigma | $\varsigma$ |
| $\backslash$ tau | $\tau$ | $\backslash$ upsilon | $v$ | $\backslash$ phi | $\phi$ |
| $\backslash$ chi | $\chi$ | $\backslash$ psi | $\psi$ | $\backslash$ omega | $\omega$ |

Greek Uppercase Letters

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ Gamma | $\Gamma$ | $\backslash$ Delta | $\Delta$ | $\backslash$ Theta | $\Theta$ |
| $\backslash$ Lambda | $\Lambda$ | $\backslash$ Xi | $\Xi$ | $\backslash$ Pi | $\Pi$ |
| $\backslash$ Sigma | $\Sigma$ | $\backslash$ Upsilon | $\Upsilon$ | $\backslash$ Phi | $\Phi$ |
| $\backslash$ Psi | $\Psi$ | $\backslash$ Omega | $\Omega$ |  |  |

Misc Symbols Type Ord

| Code | Sym | Code | Sym | Code | Sym |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $\backslash$ aleph | ふ | \wp | $\wp$ | $\backslash \mathrm{Re}$ | $\Re$ |
| $\backslash \mathrm{Im}$ | $\Im$ | $\backslash$ partial | $\partial$ | $\backslash$ infty | $\infty$ |
| $\backslash$ prime | , | $\backslash$ nabla | $\nabla$ | \surd | $\sqrt{ }$ |
| $\backslash$ angle | $\angle$ | $\backslash$ forall | $\forall$ | \exists | $\exists$ |
| $\backslash \mathrm{neg}$ | $\neg$ | $\backslash$ clubsuit | $\%$ | $\backslash$ diamondsuit | $\diamond$ |
| $\backslash$ heartsuit | $\bigcirc$ | $\backslash$ spadesuit | - |  |  |

"Large" Operators

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| \int | $\int$ |  |  |  |  |

Binary operators

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ pm | $\pm$ | $\backslash$ cdot | $\cdot$ | $\backslash$ times | $\times$ |
| $\backslash$ ast | $*$ | $\backslash$ circ | $\circ$ | $\backslash$ bullet | $\bullet$ |
| $\backslash$ div | $\div$ | $\backslash$ cap | $\cap$ | $\backslash$ cup | $\cup$ |
| $\backslash$ vee | $\vee$ | $\backslash$ wedge | $\wedge$ | $\backslash$ oplus | $\oplus$ |
| $\backslash$ otimes | $\otimes$ | $\backslash$ oslash | $\oslash$ |  |  |

Table 15.1: Available special characters in $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ mode

Relations

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ leq | $\leq$ | $\backslash$ subset | $\subset$ | $\backslash$ subseteq | $\subseteq$ |
| $\backslash$ in | $\in$ | $\backslash$ geq | $\geq$ | $\backslash$ supset | $\supset$ |
| $\backslash$ supseteq | $\supseteq$ | $\backslash$ ni | $\ni$ | $\backslash$ mid | $\mid$ |
| $\backslash$ equiv | $\equiv$ | $\backslash$ sim | $\sim$ | $\backslash$ approx | $\approx$ |
| $\backslash$ cong | $\cong$ | $\backslash$ propto | $\propto$ | $\backslash$ perp | $\perp$ |

Arrows

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ leftarrow | $\leftarrow$ | $\backslash$ Leftarrow | $\Leftarrow$ | $\backslash$ rightarrow | $\rightarrow$ |
| $\backslash$ Rightarrow | $\Rightarrow$ | $\backslash$ leftrightarrow | $\leftrightarrow$ | \uparrow | $\uparrow$ |
| $\backslash$ downarrow | $\downarrow$ |  |  |  |  |

Openings and Closings

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ lfloor | $\lfloor$ | $\backslash$ langle | $\langle$ | $\backslash$ lceil | $\lceil$ |
| $\backslash$ rfloor | $\rfloor$ | $\backslash$ rangle | $\rangle$ | $\backslash$ rceil | $\rceil$ |

Alternate Names

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ neq | $\neq$ |  |  |  |  |

Other (not in Appendix F Tables)

| Code | Sym | Code | Sym | Code | Sym |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\backslash$ ldots <br> $\backslash$ deg | $\ldots$ | $\backslash 0$ | $\oslash$ | $\backslash$ copyright | C |

Table 15.1: Available special characters in $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ mode (cont.)

## Caution: Degree Symbol

Conformance to both $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ and matlab with respect to the \circ symbol is impossible. While $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ translates this symbol to Unicode 2218 ( $\mathrm{U}+2218$ ), matlab maps this to Unicode 00B0 ( $\mathrm{U}+00 \mathrm{~B} 0$ ) instead. Octave has chosen to follow the $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ specification, but has added the additional symbol $\backslash$ deg which maps to the degree symbol ( $\mathrm{U}+00 \mathrm{~B} 0$ ).

### 15.2.8.3 latex interpreter

The "latex" interpreter only works if an external $\mathrm{LT}_{\mathrm{E}} \mathrm{X}$ tool chain is present. Three binaries are needed: latex, dvipng, and dvisvgm. If those binaries are installed but not on the path, one can still provide their respective path using the following environment variables: OCTAVE_LATEX_BINARY, OCTAVE_DVIPNG_BINARY, and OCTAVE_DVISVG_BINARY.

Note that Octave does not parse or validate the text strings when in "latex" mode - it is the responsibility of the programmer to generate valid strings which may include wrapping sections that should appear in Math mode with ' $\$$ ' characters. See, for example, https: // www.latex-project.org/help/documentation/ for documentation about $\mathrm{EAT}_{\mathrm{E}} \mathrm{X}$ typesetting.

For debugging purpose, a convenience environment variable, OCTAVE_LATEX_DEBUG_FLAG, can be set to trigger more verbose output when Octave fails to have a given text compiled by an external $\mathrm{E}_{\mathrm{E}} \mathrm{X}$ engine. For example, " $\mathrm{x}{ }^{\wedge} 2$ " is not a valid $\mathrm{EA}_{\mathrm{E}} \mathrm{X}$ string and the following example should fail

```
setenv ("OCTAVE_LATEX_DEBUG_FLAG", "1")
x = 1:10;
plot (x, x.`2)
title ("x^2", "interpreter", "latex")
```

Searching the terminal output you should find some helpful info about the origin of the failure:

```
No file default.aux.
! Missing $ inserted.
<inserted text>
    $
1.6 x^
    2
! Missing $ inserted.
```

If no usable $\mathrm{EA}_{\mathrm{E}} \mathrm{X}$ tool chain is found at the first text rendering, using the "latex" interpreter is equivalent to "none".

### 15.2.9 Printing and Saving Plots

The print command allows you to send plots to you printer and to save plots in a variety of formats. For example,

```
print -dpsc
```

prints the current figure to a color PostScript printer. And,

```
print -deps foo.eps
```

saves the current figure to an encapsulated PostScript file called foo.eps.
The current graphic toolkits produce very similar graphic displays, but differ in their capability to display unusual text and in their ability to print such text. In general, the "tex" interpreter (default) is the best all-around performer for both on-screen display and printing. However, for the reproduction of complicated text formulas the "latex" interpreter is preferred. When printing with the -painters renderer, the default for all vector formats, two options may be considered:

- Use the -svgconvert option to allow for rendering $\mathrm{L}_{\mathrm{E}} \mathrm{T}_{\mathrm{E}}$ formulas. Note that the glyph are rendered as path and the original textual info are lost.
- Use one of the -d*latex* devices to produce a .tex file (plus supporting .eps or .pdf files) to be further processed by an external $\mathrm{EAT}_{\mathrm{E}} \mathrm{X}$ engine. Note that the print function
will first set the interpreter of all strings to "latex", which means all strings must be valid $\mathrm{EAT}_{\mathrm{E}} \mathrm{X}$ strings.

A complete example showing the capabilities of text printing using the -dpdflatexstandalone option is:
$\mathrm{x}=0: 0.01: 3 ;$
hf = figure ();
plot (x, erf (x));
hold on;
plot (x, x, "r");
axis ([0, 3, 0, 1]);
text (0.65, 0.6175, ...
['\$\displaystyle\leftarrow $x=\{2$ \over \sqrt\{\pi\}\}' ... '\int_\{0\}^\{x\} e^\{-t^2\} dt = 0.6175\$'],
"interpreter", "latex");
xlabel ("x");
ylabel ("erf (x)");
title ("erf (x) with text annotation");
print (hf, "plot15_7", "-dpdflatexstandalone");
system ("pdflatex plot15_7");
open plot15_7.pdf
The result of this example can be seen in Figure 15.7


Figure 15.7: Example of inclusion of text with use of -dpdflatexstandalone

```
print ()
print (options)
print (filename, options)
print (hfig,...)
```

rgb = print ("-RGBImage", ...)
Format a figure for printing and either save it to a file, send it to a printer, or return an RGB image.
filename defines the name of the output file. If the filename has no suffix then one is inferred from the specified device and appended to the filename. When neither a filename nor the "-RGBImage" option is present, the output is sent to the printer. The various options and filename arguments may be given in any order, except for the figure handle argument hfig which must be first if it is present.

Example: Print to a file using PDF and JPEG formats.

```
figure (1);
clf ();
surf (peaks);
print figure1.pdf # The extension specifies the format
print -djpg figure1 # Will produce "figure1.jpg" file
```

If the first argument is a handle hfig to a figure object then it specifies the figure to print. By default, the current figure returned by gcf is printed.

For outputs to paged formats, for example, PostScript and PDF, the page size is specified by the figure's papersize property together with the paperunits property. The location and size of the plot on the page are specified by the figure's paperposition property. The orientation of the page is specified by the figure's paperorientation property.
For non-page formats-for example, image formats like JPEG-the width and height of the output are specified by the figure's paperposition(3:4) property values.

The print command supports many options:
-fh Specify the handle, $h$, of the figure to be printed.
Example: Print figure 1.

```
figure (1);
clf ();
surf (peaks);
figure (2);
print -f1 figure1.pdf
## Equivalent functional form:
print (1, "figure1.pdf")
```

-Pprinter

Set the printer name to which the plot is sent if no filename is specified.
Example: Print to printer named PS_printer using PostScript format.

```
clf ();
surf (peaks);
print -dpswrite -PPS_printer
```

-RGBImage
Return an M-by-N-by-3 RGB image of the figure. The size of the image depends on the formatting options. This is similar to taking a screen
capture of the plot, but formatting options may be changed such as the resolution or monochrome/color.
Example: Get the pixels of a figure image.

```
clf ();
surf (peaks);
rgb = print ("-RGBImage");
```

-opengl
-painters

Specifies whether the opengl (pixel-based) or painters (vector-based) renderer is used. This is equivalent to changing the figure's "Renderer" property. When the figure "RendererMode" property is "auto" (the default) Octave will use the "opengl" renderer for raster formats (e.g., JPEG) and "painters" for vector formats (e.g., PDF). Those options are only supported for the "qt" graphics toolkit.

## -svgconvert

When using the -painters renderer, this enables a different backend toolchain with enhanced characteristics:
Font handling:
For interpreters "none" and "tex", the actual font is embedded in the output file which allows for printing arbitrary characters and fonts in all vector formats.
Strings using the "latex" interpreter, are rendered using path objects. This looks good but note that textual info (font, characters. . .) are lost.
Output Simplification:
By default, the option -painters renders patch and surface objects using assemblies of triangles. This may lead to antialiasing artifacts when viewing the file. The -svgconvert option reconstructs polygons in order to avoid those artifacts (particularly for 2-D figures).
Transparency:
Allows for printing transparent graphics objects in PDF format. For PostScript formats the presence of any transparent object will cause the output to be rasterized.
Caution: -svgconvert may lead to inaccurate rendering of image objects.

## -portrait

-landscape
Specify the orientation of the plot for printed output. For non-printed output the aspect ratio of the output corresponds to the plot area defined by the "paperposition" property in the orientation specified. This option is equivalent to changing the figure's "paperorientation" property.
-fillpage
-bestfit When using a page-based format (PDF, PostScript, printer) ignore the "paperposition" property and have the plot occupy the entire page.

The option -fillpage will stretch the plot to occupy the page with 0.25 inch margins all around. The option -bestfit will expand the plot to take up as much room as possible on the page without distorting the original aspect ratio of the plot.
-color
-mono Color or monochrome output.
-solid
-dashed Force all lines to be solid or dashed, respectively.
-noui Don't print uicontrol objects such as pushbuttons which may overlay the plot. This is the default behavior and it is not possible to include uicontrol objects in the output without using an external screen capture tool.
-rNUM Resolution of bitmaps in dots per inch (DPI). For both metafiles and SVG the default is the screen resolution; for other formats the default is 150 DPI. To specify screen resolution, use "-r0".
Example: high resolution raster output.

```
clf ();
surf (peaks (), "facelighting", "gouraud");
light ();
print ("-r600", "lit_peaks.png");
```

-Sxsize,ysize
Plot size in pixels for raster formats including PNG, JPEG, PNG, and unusually SVG. For all vector formats, including PDF, PS, and EPS, the plot size is specified in points. This option is equivalent to changing the width and height of the output by setting the figure property paperposition(3:4). When using the command form of the print function you must quote the xsize,ysize option to prevent the Octave interpreter from recognizing the embedded comma (','). For example, by writing "-S640,480".
-tight
-loose Force a tight or loose bounding box for EPS files. The default is tight.
-preview Add a preview to EPS files. Supported formats are:
-interchange
Provide an interchange preview.
-metafile
Provide a metafile preview.
-pict Provide a pict preview.
-tiff Provide a TIFF preview.
-append Append PostScript or PDF output to an existing file of the same type.
-Ffontname
-Ffontname:size
-F:size Use fontname and/or fontsize for all text. fontname is ignored for some devices: fig, etc.
-ddevice The available output format is specified by the option device, and is one of the following (devices marked with a ' $*$ ' are only available with the Gnuplot toolkit):
Vector Formats

```
svg Scalable Vector Graphics.
```

pdf
pdfcrop Portable Document Format. The pdf device formats the fig-
ure for printing on paper. The size of the surrounding page and the position of the figure inside the page are defined by the [paper* figure properties], page 437.
Use pdfcrop if you don't want the surrounding page.
By default, PDF inherits the same limitations as PostScript. For an enhanced output with complete text support and basic transparency, use the -svgconvert option.

```
eps(2)
```

epsc(2) Encapsulated PostScript (level 1 and 2, mono and color).
The OpenGL-based graphics toolkits always generate PostScript level 3.0. They have limited support for text unless using the -svgconvert option. Limitations include using only ASCII characters (e.g., no Greek letters) and support for just three base PostScript fonts: Helvetica (the default), Times, or Courier. Any other font will be replaced by Helvetica.
ps(2)
$\mathrm{psc}(2) \quad$ Same as eps except that the figure is formatted for printing on paper. The size of the surrounding page and position of the figure inside the page are defined by the [paper* figure properties], page 437 .

```
pslatex
epslatex
pdflatex
pslatexstandalone
epslatexstandalone
pdflatexstandalone
```

Generate a $\mathrm{AT}_{\mathrm{E}} \mathrm{X}$ file filename.tex for the text portions of a plot and a file filename. (ps|eps|pdf) for the remaining graphics. The graphics file suffix .pslepslpdf is determined by the specified device type. The $\mathrm{IAT}_{\mathrm{E}} \mathrm{X}$ file produced by the 'standalone' option can be processed directly by $\mathrm{LAT}_{\mathrm{E}} \mathrm{X}$. The file generated without the 'standalone' option is intended to be included from another $\mathrm{EA}_{\mathrm{E}} \mathrm{X}$ document. In either case, the ${ }^{\mathrm{LA}} \mathrm{T}_{\mathrm{E}} \mathrm{X}$ file contains an includegraphics command so that the generated graphics file is automatically included when the ${ }^{A} T_{E} \mathrm{X}$ file is processed. The text that is written to the
${ }^{\mathrm{LA}} \mathrm{T}_{\mathrm{E}} \mathrm{X}$ file contains the strings exactly as they were specified in the plot. If any special characters of the $\mathrm{T}_{\mathrm{E}} \mathrm{X}$ mode interpreter were used, the file must be edited before $\mathrm{LA}_{\mathrm{E}} \mathrm{X}$ processing. Specifically, the special characters must be enclosed with dollar signs ( $\$ \ldots \$$ ), and other characters that are recognized by $\mathrm{AT}_{\mathrm{E}} \mathrm{X}$ may also need editing (e.g., braces). The 'pdflatex' device, and any of the 'standalone' formats, are not available with the Gnuplot toolkit.

```
epscairo*
pdfcairo*
epscairolatex*
pdfcairolatex*
epscairolatexstandalone*
pdfcairolatexstandalone*
    Generate output with Cairo renderer. The devices
    epscairo and pdfcairo are synonymous with the
    epsc device. The IATEX variants generate a \ATEX file,
    filename.tex, for the text portions of a plot, and an image
    file, filename.(epslpdf), for the graph portion of the
    plot. The 'standalone' variants behave as described for
    'epslatexstandalone' above.
canvas* Javascript-based drawing on an HTML5 canvas viewable in
        a web browser.
emf
meta Microsoft Enhanced Metafile
fig XFig. For the Gnuplot graphics toolkit, the additional op-
        tions -textspecial or -textnormal (default) can be used to
        control whether the special flag should be set for the text in
        the figure.
latex*
eepic* EATEX picture environment and extended picture environ-
        ment.
tikz
tikzstandalone*
    Generate a EATEX file using PGF/TikZ format. The OpenGL-
    based toolkits create a PGF file while Gnuplot creates a TikZ
    file. The 'tikzstandalone' device produces a IATEX docu-
    ment which includes the TikZ file.
```

Raster Formats

```
png Portable Network Graphics
jpg
jpeg JPEG image
```

```
tif
tiff
tiffn TIFF image with LZW compression (tif, tiff) or
uncompressed (tiffn).
gif GIF image
pbm PBMplus
dumb* ASCII art
```

If the device is omitted, it is inferred from the file extension, or if there is no filename then it is sent to the printer as PostScript.
-dghostscript_device
Additional devices are supported by Ghostscript. Some examples are:
ljet2p HP LaserJet IIP
pcx24b 24-bit color PCX file format
ppm Portable Pixel Map file format
For a complete list of available formats and devices type system ("gs -h").
When Ghostscript output is sent to a printer the size is determined by the figure's "papersize" property. When the output is sent to a file the size is determined by the plot box defined by the figure's "paperposition" property.
-Gghostscript_command
Specify the command for calling Ghostscript. For Unix the default is "gs" and for Windows it is "gswin32c".
-TextAlphaBits=n
-GraphicsAlphaBits=n
Octave is able to produce output for various printers, bitmaps, and vector formats by using Ghostscript. For bitmap and printer output anti-aliasing is applied using Ghostscript's TextAlphaBits and GraphicsAlphaBits options. The default number of bits are 4 and 1 respectively. Allowed values for $N$ are 1,2 , or 4 .

See also: [saveas], page 416, [getframe], page 869, [savefig], page 419, [hgsave], page 417, [orient], page 417, [figure], page 399.
saveas (h, filename)
saveas (h, filename, fmt)
Save graphics object $h$ to the file filename in graphics format fmt.
If $h$ is the handle to a figure object, that figure object is saved. If $h$ is the handle to a different graphics object, the figure containing that graphics object is saved.
All device formats accepted by print may be used. Common formats are:
ofig Octave figure file format (default)

```
mfig Two files: Octave m-file filename.m containing code to open Octave
        figure file filename.ofig
ps PostScript
eps Encapsulated PostScript
pdf Portable Document Format
jpg JPEG Image
png Portable Network Graphics image
emf Enhanced MetaFile
tif TIFF Image, compressed
```

If fmt is omitted it is extracted from the extension of filename. The default format when there is no extension is "ofig".

```
clf ();
surf (peaks);
saveas (1, "figure1.png");
```

See also: [print], page 410, [savefig], page 419, [hgsave], page 417, [orient], page 417.

```
orient (orientation)
```

orient (hfig, orientation)
orientation $=$ orient ()
orientation $=$ orient (hfig)

Query or set the print orientation for figure hfig.
Valid values for orientation are "portrait", "landscape", and "tall".
The "landscape" option changes the orientation so the plot width is larger than the plot height. The "paperposition" is also modified so that the plot fills the page, while leaving a 0.25 inch border.
The "tall" option sets the orientation to "portrait" and fills the page with the plot, while leaving a 0.25 inch border.
The "portrait" option (default) changes the orientation so the plot height is larger than the plot width. It also restores the default "paperposition" property.
When called with no arguments, return the current print orientation.
If the argument hfig is omitted, then operate on the current figure returned by gcf.
See also: [print], page 410, [saveas], page 416.
print and saveas are used when work on a plot has finished and the output must be in a publication-ready format. During intermediate stages it is often better to save the graphics object and all of its associated information so that changes - to colors, axis limits, marker styles, etc.-can be made easily from within Octave. The hgsave/hgload commands can be used to save and re-create a graphics object.

```
hgsave (filename)
hgsave (h, filename)
```

hgsave (h, filename, fmt)
Save the graphics handle(s) $h$ to the file filename in the format fmt.
If unspecified, $h$ is the current figure as returned by gcf.
When filename does not have an extension the default filename extension . ofig will be appended.
If present, fmt must be one of the following:

- -binary, -float-binary
- -hdf5, -float-hdf5
- -V7, -v7, -7, -mat7-binary
- -V6, -v6, -6, -mat6-binary
- -text
- -zip, -z

The default format is -binary to minimize storage.
Programming Note: When producing graphics for final publication use print or saveas. When it is important to be able to continue to edit a figure as an Octave object, use hgsave/hgload.

See also: [hgload], page 418, [hdl2struct], page 432, [savefig], page 419, [saveas], page 416 , [print], page 410.
$h=h g l o a d ~(f i l e n a m e)$
[h, old_prop] = hgload (filename, prop_struct)
Load the graphics objects in filename into a vector of graphics handles $h$.
If filename has no extension, Octave will try to find the file with and without the default extension .ofig.
If provided, the elements of structure prop_struct will be used to override the properties of top-level objects stored in filename, and the saved values from filename will be stored in old_prop. old_prop is a cell array matching the size of $h$; each cell contains a structure of the existing property names and values before being overridden.
See also: [openfig], page 418, [hgsave], page 417, [struct2hdl], page 432.

```
openfig
openfig (filename)
openfig (..., copies)
openfig (..., visibility)
h = openfig (...)
```

Read saved figure window(s) from filename and return graphics handle(s) h.
By default, filename is "Untitled.fig". If a full path is not specified, the file opened will be the first one encountered in the load path. If filename is not found and does not have an extension, a search will take place for the first file in the load path with extension ".fig" or ".ofig", in that order.
copies is an optional input indicating whether a new figure should be created ("new") or whether an existing figure may be reused ("reuse"). An existing figure may be reused if the "FileName" property matches the specified input filename. When a
figure is reused it becomes the active figure and is shown on top of other figures. If the figure was offscreen, it is re-positioned to be onscreen. The default value for copies is "new".
visibility is an optional input indicating whether to show the figure ("visible") or not ("invisible"). When visibility is specified as an input to openfig it overrides the visibility setting stored in filename.
See also: [open], page 973, [hgload], page 418, [savefig], page 419, [struct2hdl], page 432.

```
savefig()
savefig (h)
savefig (filename)
savefig (h, filename)
savefig (h, filename, "compact")
```

Save figure windows specified by graphics handle(s) $h$ to file filename.
If unspecified, $h$ is the current figure returned by gcf.
If unspecified, filename is set to "Untitled.fig". If filename does not have an extension then the default extension ".fig" will be added.
If the optional third input "compact" is present then the data will be compressed to save more space.
See also: [hgsave], page 417, [hdl2struct], page 432, [openfig], page 418.

### 15.2.10 Interacting with Plots

The user can select points on a plot with the ginput function or select the position at which to place text on the plot with the gtext function using the mouse.
[x, y, buttons] = ginput (n)
[x, y, buttons] = ginput ()
Return the position and type of mouse button clicks and/or key strokes in the current figure window.
If $n$ is defined, then capture $n$ events before returning. When $n$ is not defined ginput will loop until the return key RET is pressed.
The return values $x, y$ are the coordinates where the mouse was clicked in the units of the current axes. The return value button is 1,2 , or 3 for the left, middle, or right button. If a key is pressed the ASCII value is returned in button.
Implementation Note: ginput is intenteded for 2-D plots. For 3-D plots see the currentpoint property of the current axes which can be transformed with knowledge of the current view into data units.
See also: [gtext], page 420, [waitforbuttonpress], page 419.

```
waitforbuttonpress ()
```

b = waitforbuttonpress ()
Wait for mouse click or key press over the current figure window.
The return value of $b$ is 0 if a mouse button was pressed or 1 if a key was pressed.
See also: [waitfor], page 936, [ginput], page 419, [kbhit], page 278.

```
gtext (s)
gtext ({s1, s2, ...})
gtext ({s1; s2; ...})
gtext (..., prop, val, ...)
h = gtext (...)
```

Place text on the current figure using the mouse.
The string argument $s$ may be a character array or a cell array of strings. If $s$ has more than one row, each row is used to create a separate text object after a mouse click. For example:
Place a single string after one mouse click

```
gtext ("I clicked here")
```

Place two strings after two mouse clicks

```
gtext ({"I clicked here"; "and there"})
```

Place two strings, each with two lines, after two mouse clicks

```
gtext ({"I clicked", "here"; "and", "there"})
```

Optional property/value pairs are passed directly to the underlying text objects.
The full list of text object properties is documented at Section 15.3.3.6 [Text Properties], page 453.
The optional return value $h$ holds the graphics handle(s) to the created text object(s).
See also: [ginput], page 419, [text], page 392.
More sophisticated user interaction mechanisms can be obtained using the ui* family of functions, see Section 35.3 [UI Elements], page 924.

### 15.2.11 Test Plotting Functions

The functions sombrero and peaks provide a way to check that plotting is working. Typing either sombrero or peaks at the Octave prompt should display a three-dimensional plot.

```
sombrero ()
sombrero (n)
z = sombrero (...)
[x, y, z] = sombrero (...)
```

Plot the familiar 3-D sombrero function.
The function plotted is

$$
z=\frac{\sin \left(\sqrt{\left(\mathrm{x}^{2}+\mathrm{y}^{2}\right)}\right)}{\sqrt{\left(\mathrm{x}^{2}+\mathrm{y}^{2}\right)}}
$$

Called without a return argument, sombrero plots the surface of the above function over the meshgrid $[-8,8]$ using surf.
If $n$ is a scalar the plot is made with $n$ grid lines. The default value for $n$ is 41 .
When called with output arguments, return the data for the function evaluated over the meshgrid. This can subsequently be plotted with $\operatorname{surf}(x, y, z)$.
See also: [peaks], page 421, [meshgrid], page 373, [mesh], page 358, [surf], page 360.

```
peaks ()
peaks (n)
peaks (x, y)
z = peaks (...)
[x, y, z] = peaks (...)
```

Plot a function with lots of local maxima and minima.
The function has the form

$$
f(x, y)=3(1-x)^{2} e^{\left(-x^{2}-(y+1)^{2}\right)}-10\left(\frac{x}{5}-x^{3}-y^{5}\right)-\frac{1}{3} e^{\left(-(x+1)^{2}-y^{2}\right)}
$$

Called without a return argument, peaks plots the surface of the above function using surf.
If $n$ is a scalar, peaks plots the value of the above function on an $n$-by- $n$ mesh over the range $[-3,3]$. The default value for $n$ is 49 .
If $n$ is a vector, then it represents the grid values over which to calculate the function. If $x$ and $y$ are specified then the function value is calculated over the specified grid of vertices.
When called with output arguments, return the data for the function evaluated over the meshgrid. This can subsequently be plotted with $\operatorname{surf}(x, y, z)$.
See also: [sombrero], page 420, [meshgrid], page 373, [mesh], page 358, [surf], page 360 .

### 15.3 Graphics Data Structures

### 15.3.1 Introduction to Graphics Structures

The graphics functions use pointers, which are of class graphics_handle, in order to address the data structures which control visual display. A graphics handle may point to any one of a number of different base object types. These objects are the graphics data structures themselves. The primitive graphic object types are: figure, axes, line, text, patch, scatter, surface, text, image, and light.

Each of these objects has a function by the same name, and each of these functions returns a graphics handle pointing to an object of the corresponding type.

In addition, there are several functions which operate on properties of the graphics objects and which also return handles. This includes but is not limited to the following functions: The functions plot and plot3 return a handle pointing to an object of type line. The function subplot returns a handle pointing to an object of type axes. The functions fill, fill3, trimesh, and trisurf return a handle pointing to an object of type patch. The function scatter3 returns a handle to an object of type scatter. The functions slice, surf, surfl, mesh, meshz, pcolor, and waterfall each return a handle of type surface. The function camlight returns a handle to an object of type light. The functions area, bar, barh, contour, contourf, contour3, surfc, meshc, errorbar, quiver, quiver3, stair, stem, stem3 each return a handle to a complex data structure as documented in [Data Sources], page 500.

The graphics objects are arranged in a hierarchy:

1. The root object is returned by groot (historically, equivalent to the handle 0 ). In other words, get (groot) returns the properties of the root object.
2. Below the root are figure objects.
3. Below the figure objects are axes or hggroup objects.
4. Below the axes or hggroup objects are line, text, patch, scatter, surface, image, and light objects.

It is possible to walk this hierarchical tree by querying the "parent" and "children" properties of the graphics objects.

Graphics handles may be distinguished from function handles (see Section 11.12.1 [Function Handles], page 233) by means of the function ishghandle. ishghandle returns true if its argument is a handle of a graphics object. In addition, a figure or axes object may be tested using isfigure or isaxes respectively. To test for a specific type of graphics handle, such as a patch or line object, use isgraphics. The more specific test functions return true only if the argument is both a graphics handle and of the correct type (figure, axes, specified object type).

The get and set commands are used to obtain and set the values of properties of graphics objects. In addition, the get command may be used to obtain property names.

For example, the property "type" of the graphics object pointed to by the graphics handle h may be displayed by:

```
get (h, "type")
```

The properties and their current values may be obtained in the form of a structure using $\mathrm{s}=\operatorname{get}(\mathrm{h})$, where h is the handle of a graphics object. If only the names of the properties and the allowed values (for radio properties only) are wanted, one may use set (h).

Thus, for example:

```
h = figure ();
get (h, "type")
=> ans = figure
set (h)
#
    alphamap:
    beingdeleted: [ {off} | on ]
    busyaction: [ cancel | {queue} ]
    buttondownfcn:
    clipping: [ off | {on} ]
    closerequestfcn:
    color:
    colormap:
    createfcn:
    currentaxes:
    deletefcn:
    dockcontrols: [ {off} | on ]
    filename:
    graphicssmoothing: [ off | {on} ]
handlevisibility: [ callback | off | {on} ]
```

The uses of get and set are further explained in [get], page 430, [set], page 430.

## res = isprop (obj, "prop")

Return true if prop is a property of the object obj.
obj may also be an array of objects in which case res will be a logical array indicating whether each handle has the property prop.
For plotting, obj is a handle to a graphics object. Otherwise, obj should be an instance of a class. isprop reports whether the class defines a property, but Access permissions or visibility restrictions (Hidden $=$ true) may prevent use by the programmer.
See also: [get], page 430, [set], page 430, [properties], page 909, [ismethod], page 891, [isobject], page 890.

### 15.3.2 Graphics Objects

The hierarchy of graphics objects was explained above. See Section 15.3.1 [Introduction to Graphics Structures], page 421. Here the specific objects are described, and the properties contained in these objects are discussed. Keep in mind that graphics objects are always referenced by handle.
root The top level of the hierarchy and the parent of all figure objects. Use groot to obtain the handle of the root graphics object.
figure A figure window.
axes A set of axes. This object is a child of a figure object and may be a parent of line, text, image, patch, surface, or light objects.
line A line in two or three dimensions.
text Text annotations.
image A bitmap image.
patch A filled polygon, currently limited to two dimensions.
surface A three-dimensional surface.
light A light object used for lighting effects on patches and surfaces.

### 15.3.2.1 Creating Graphics Objects

You can create any graphics object primitive by calling the function of the same name as the object; In other words, figure, axes, line, text, image, patch, surface, and light functions. These fundamental graphic objects automatically become children of the current axes object as if hold on was in place. Separately, axes will automatically become children of the current figure object and figures will become children of the root object.

If this auto-joining feature is not desired then it is important to call newplot first to prepare a new figure and axes for plotting. Alternatively, the easier way is to call a highlevel graphics routine which will both create the plot and then populate it with low-level graphics objects. Instead of calling line, use plot. Or use surf instead of surface. Or use fill or fill3 instead of patch.

```
axes ()
axes (property, value,...)
axes (hpar, property, value,...)
axes (hax)
h = axes (...)
```

Create a Cartesian axes object and return a handle to it, or set the current axes to hax.

Called without any arguments, or with property/value pairs, construct a new axes. The optional argument hpar is a graphics handle specifying the parent for the new axes and may be a figure, uipanel, or uitab.
Called with a single axes handle argument hax, the function makes hax the current axes (as returned by gca). It also makes the figure which contains hax the current figure (as returned by gcf). Finally, it restacks the parent object's children property so that the axes hax appears before all other axes handles in the list. This causes hax to be displayed on top of any other axes objects (Z-order stacking). In addition it restacks any legend or colorbar objects associated with hax so that they are also visible.
Programming Note: The full list of properties is documented at Section 15.3.3.3 [Axes Properties], page 441.

See also: [gca], page 429, [set], page 430, [get], page 430.
line ()
line ( $x, y$ )
line ( $x, y, z$ )
line ("xdata", x, "ydata", y)
line ("xdata", x, "ydata", y, "zdata", z)
line (..., property, value)
line (hax, ...)
$h=$ line (...)
Create a line object from $x$ and $y$ (and possibly $z$ ) and insert it in the current axes.
In the standard calling form the data $x, y$, and $z$ may be scalars, vectors, or matrices. In the case of matrix inputs, line will attempt to orient scalars and vectors so the results can be plotted. This requires that one of the dimensions of the vector match either the number of rows or the number of columns of the matrix.
In the low-level calling form ( $50 \%$ higher performance) where the data is specified by name (line ("xdata", x, ...)) the data must be vectors. If no data is specified (line ()) then $x==y=[0,1]$.
Multiple property-value pairs may be specified for the line object, but they must appear in pairs.
If called with only property/value pairs then any unspecified properties use their default values as specified on the root object.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle (or vector of handles) to the line objects created.

Programming Note: The full list of properties is documented at Section 15.3.3.5 [Line Properties], page 450.
The function line differs from plot in that line objects are inserted in to the current axes without first clearing the plot.
See also: [image], page 866, [patch], page 425, [rectangle], page 356, [surface], page 426, [text], page 392.
patch ()
patch (x, y, c)
patch (x, y, z, c)
patch ("Faces", faces, "Vertices", verts, ...)
patch (..., prop, val, ...)
patch (..., propstruct, ...)
patch (hax, ...)
$h=\operatorname{patch}(. .$.
Create patch object in the current axes with vertices at locations ( $x, y$ ) and of color c.

If the vertices are matrices of size MxN then each polygon patch has M vertices and a total of N polygons will be created. If some polygons do not have M vertices use NaN to represent "no vertex". If the $z$ input is present then 3-D patches will be created.
The color argument $c$ can take many forms. To create polygons which all share a single color use a string value (e.g., "r" for red), a scalar value which is scaled by caxis and indexed into the current colormap, or a 3 -element RGB vector with the precise TrueColor.
If $c$ is a vector of length $N$ then the ith polygon will have a color determined by scaling entry $c(\mathrm{i})$ according to caxis and then indexing into the current colormap. More complicated coloring situations require directly manipulating patch property/value pairs.
Instead of specifying polygons by matrices $x$ and $y$, it is possible to present a unique list of vertices and then a list of polygon faces created from those vertices. In this case the "Vertices" matrix will be an Nx2 (2-D patch) or Nx3 (3-D patch). The MxN "Faces" matrix describes M polygons having N vertices - each row describes a single polygon and each column entry is an index into the "Vertices" matrix to identify a vertex. The patch object can be created by directly passing the property/value pairs "Vertices"/verts, "Faces"/faces as inputs.
Instead of using property/value pairs, any property can be set by passing a structure propstruct with the respective field names.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.

The optional return value $h$ is a graphics handle to the created patch object.
Programming Note: The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457. Useful patch properties include: "cdata", "edgecolor", "facecolor", "faces", and "facevertexcdata".
See also: [fill], page 346, [get], page 430, [set], page 430.

```
surface (x, y, z,c)
surface (x, y,z)
surface (z,c)
surface (z)
surface (..., prop, val, ...)
surface (hax,...)
h = surface (...)
```

Create a surface graphic object given matrices $x$ and $y$ from meshgrid and a matrix of values $z$ corresponding to the $x$ and $y$ coordinates of the surface.
If $x$ and $y$ are vectors, then a typical vertex is $(x(j), y(i), z(i, j))$. Thus, columns of $z$ correspond to different $x$ values and rows of $z$ correspond to different $y$ values. If only a single input $z$ is given then $x$ is taken to be 1 :columns ( $z$ ) and $y$ is 1 :rows (z).

Any property/value input pairs are assigned to the surface object.
If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created surface object.
Programming Note: The full list of properties is documented at Section 15.3.3.10 [Surface Properties], page 465.

See also: [surf], page 360, [mesh], page 358, [patch], page 425, [line], page 424.
light ()
light (..., "prop", val, ...)
light (hax, ...)
$h=$ light (...)
Create a light object in the current axes or for axes hax.
When a light object is present in an axes object, and the properties "EdgeLighting" or "FaceLighting" of a patch or surface object are set to a value other than "none", these objects are drawn with lighting effects. Supported values for Lighting properties are "none" (no lighting effects), "flat" (faceted look of the objects), and "gouraud" (linear interpolation of the lighting effects between the vertices). If the lighting mode is set to "flat", the "FaceNormals" property is used for lighting. For "gouraud", the "VertexNormals" property is used.
Up to eight light objects are supported per axes. (Implementation dependent)
Lighting is only supported for OpenGL graphic toolkits (i.e., "fltk" and "qt").
A light object has the following properties which alter the appearance of the plot.
"Color": The color of the light can be passed as an RGB-vector (e.g., [100] for red) or as a string (e.g., "r" for red). The default color is white ([lll $\left.\left.1 \begin{array}{ll}1 & 1 \\ 1\end{array}\right]\right)$.
"Position": The direction from which the light emanates as a
1 x 3 -vector. The default direction is [101 101$].$
"Style": This string defines whether the light emanates from a light source at infinite distance ("infinite") or from a local point source ("local"). The default is "infinite".

If the first argument hax is an axes handle, then add the light object to this axes, rather than the current axes returned by gca.
The optional return value $h$ is a graphics handle to the created light object.
Programming Note: The full list of properties is documented at Section 15.3.3.11 [Light Properties], page 468.
See also: [lighting], page 371, [material], page 371, [patch], page 425, [surface], page 426.

### 15.3.2.2 Handle Functions

To determine whether a variable is a graphics object index, or an index to an axes or figure, use the functions ishghandle, isgraphics, isaxes, and isfigure.
ishghandle (h)
Return true if $h$ is a graphics handle and false otherwise.
$h$ may also be a matrix of handles in which case a logical array is returned that is true where the elements of $h$ are graphics handles and false where they are not.
See also: [isgraphics], page 427, [isaxes], page 427, [isfigure], page 427, [ishandle], page 427.
isgraphics (h)
isgraphics (h, type)
Return true if $h$ is a graphics handle (of type type) and false otherwise.
When no type is specified the function is equivalent to ishghandle.
See also: [ishghandle], page 427, [ishandle], page 427, [isaxes], page 427, [isfigure], page 427.
ishandle (h)
Return true if $h$ is a handle to a graphics or Java object and false otherwise.
$h$ may also be a matrix of handles in which case a logical array is returned that is true where the elements of $h$ are handles to graphics or Java objects and false where they are not.
Programming Note: It is often more useful to test for a specific object type. To determine if a handle belongs to a graphics object use ishghandle or isgraphics. To determine if a handle belongs to a Java object use isjava.
See also: [ishghandle], page 427, [isgraphics], page 427, [isjava], page 1055.
isaxes (h)
Return true if $h$ is an axes graphics handle and false otherwise.
If $h$ is a matrix then return a logical array which is true where the elements of $h$ are axes graphics handles and false where they are not.
See also: [isfigure], page 427, [ishghandle], page 427, [isgraphics], page 427.
isfigure (h)
Return true if $h$ is a figure graphics handle and false otherwise.

If $h$ is a matrix then return a logical array which is true where the elements of $h$ are figure graphics handles and false where they are not.

See also: [isaxes], page 427, [ishghandle], page 427, [isgraphics], page 427.
The function gcf returns an index to the current figure object, or creates one if none exists. Similarly, gca returns the current axes object, or creates one (and its parent figure object) if none exists.
$h=$ groot ()
Return a handle to the root graphics object.
The root graphics object is the ultimate parent of all graphics objects.
In addition, the root object contains information about the graphics system as a whole such as the ScreenSize. Use get (groot) to find out what information is available.
Defaults for the graphic system as a whole are specified by setting properties of the root graphics object that begin with "Default". For example, to set the default font for all text objects to FreeSans use

```
set (groot, "DefaultTextFontName", "FreeSans")
```

Default properties can be deleted by using set with the special property value of "remove". To undo the default font assignment above use

```
set (groot, "DefaultTextFontName", "remove")
```

Programming Note: The root graphics object is identified by the special handle value of 0 . At some point this unique value may change, but code can be made resistant to future changes by using groot which is guaranteed to always return the root graphics object.

See also: [gcf], page 428, [gca], page 429, [get], page 430, [set], page 430.
$h=\operatorname{gcf}()$
Return a handle to the current figure.
The current figure is the default target for graphics output. If multiple figures exist, gcf returns the last created figure or the last figure that was clicked on with the mouse.
If a current figure does not exist, create one and return its handle. The handle may then be used to examine or set properties of the figure. For example,

```
fplot (@sin, [-10, 10]);
fig = gcf ();
set (fig, "numbertitle", "off", "name", "sin plot")
```

plots a sine wave, finds the handle of the current figure, and then renames the figure window to describe the contents.
Note: To find the current figure without creating a new one if it does not exist, query the "CurrentFigure" property of the root graphics object.

```
get (groot, "currentfigure");
```

See also: [gca], page 429, [gco], page 429, [gcbf], page 494, [gcbo], page 494, [get], page 430, [set], page 430.
$h=\operatorname{gca}()$
Return a handle to the current axes object.
The current axes is the default target for graphics output. In the case of a figure with multiple axes, gca returns the last created axes or the last axes that was clicked on with the mouse.
If no current axes object exists, create one and return its handle. The handle may then be used to examine or set properties of the axes. For example,

```
ax = gca ();
set (ax, "position", [0.5, 0.5, 0.5, 0.5]);
```

creates an empty axes object and then changes its location and size in the figure window.
Note: To find the current axes without creating a new axes object if it does not exist, query the "CurrentAxes" property of a figure.
get (gcf, "currentaxes");
See also: [gcf], page 428, [gco], page 429, [gcbf], page 494, [gcbo], page 494, [get], page 430, [set], page 430.
$h=\operatorname{gco}()$
$h=g c o(f i g)$
Return a handle to the current object of the current figure, or a handle to the current object of the figure with handle fig.
The current object of a figure is the object that was last clicked on. It is stored in the "CurrentObject" property of the target figure.
If the last mouse click did not occur on any child object of the figure, then the current object is the figure itself.
If no mouse click occurred in the target figure, this function returns an empty matrix.
Programming Note: The value returned by this function is not necessarily the same as the one returned by gcbo during callback execution. An executing callback can be interrupted by another callback and the current object may be changed.
See also: [gcbo], page 494, [gca], page 429, [gcf], page 428, [gcbf], page 494, [get], page 430, [set], page 430.
The get and set functions may be used to examine and set properties for graphics objects. For example,

```
get (groot)
    ans =
        {
            type = root
            currentfigure = [] (0x0)
            children = [] (0x0)
            visible = on
            }
```

returns a structure containing all the properties of the root graphics object. As with all functions in Octave, the structure is returned by value, so modifying it will not modify the
internal root object. To do that, you must use the set function. Also, note that in this case, the currentfigure property is empty, which indicates that there is no current figure window.

The get function may also be used to find the value of a single property. For example,

```
get (gca (), "xlim")
    # [ 0 1 1]
```

returns the range of the x -axis for the current axes object in the current figure.
To set graphics object properties, use the set function. For example,

```
set (gca (), "xlim", [-10, 10]);
```

sets the range of the $x$-axis for the current axes object in the current figure to ' $[-10,10]$ '.
Default property values can also be queried if the set function is called without a value argument. When only one argument is given (a graphic handle) then a structure with defaults for all properties of the given object type is returned. For example,

```
set (gca ())
```

returns a structure containing the default property values for axes objects. If set is called with two arguments (a graphic handle and a property name) then only the defaults for the requested property are returned.

```
val = get (h)
val = get (h, p)
```

Return the value of the named property $p$ from the graphics handle $h$.
If $p$ is omitted, return the complete property list for $h$.
If $h$ is a vector, return a cell array including the property values or lists respectively.
See also: [set], page 430.
set (h, property, value, ...)
set (h, properties, values)
set (h, pv)
value_list = set (h, property)
all_value_list = set (h)
Set named property values for the graphics handle (or vector of graphics handles) h.
There are three ways to give the property names and values:

- as a comma separated list of property, value pairs

Here, each property is a string containing the property name, each value is a value of the appropriate type for the property.

- as a cell array of strings properties containing property names and a cell array values containing property values.
In this case, the number of columns of values must match the number of elements in properties. The first column of values contains values for the first entry in properties, etc. The number of rows of values must be 1 or match the number of elements of $h$. In the first case, each handle in $h$ will be assigned the same values. In the latter case, the first handle in $h$ will be assigned the values from the first row of values and so on.
- as a structure array $p v$

Here, the field names of $p v$ represent the property names, and the field values give the property values. In contrast to the previous case, all elements of $p v$ will be set in all handles in $h$ independent of the dimensions of $p v$.
set is also used to query the list of values a named property will take. clist $=$ set ( $h$, "property") will return the list of possible values for "property" in the cell list clist. If no output variable is used then the list is formatted and printed to the screen.
If no property is specified (slist $=$ set $(h))$ then a structure slist is returned where the fieldnames are the properties of the object $h$ and the fields are the list of possible values for each property. If no output variable is used then the list is formatted and printed to the screen.
For example,

```
hf = figure ();
set (hf, "paperorientation")
# [ landscape | {portrait} ]
```

shows the paperorientation property can take two values with the default being "portrait".

See also: [get], page 430.
parent $=$ ancestor (h, type)
parent = ancestor (h, type, "toplevel")
Return the first ancestor of handle object $h$ whose type matches type, where type is a character string.

If type is a cell array of strings, return the first parent whose type matches any of the given type strings.

If the handle object $h$ itself is of type type, return $h$.
If "toplevel" is given as a third argument, return the highest parent in the object hierarchy that matches the condition, instead of the first (nearest) one.

See also: [findobj], page 488, [findall], page 489, [allchild], page 431.
$h$ = allchild (handles)
Find all children, including hidden children, of a graphics object.
This function is similar to get (h, "children"), but also returns hidden objects (HandleVisibility $=$ "off").
If handles is a scalar, $h$ will be a vector. Otherwise, $h$ will be a cell matrix of the same size as handles and each cell will contain a vector of handles.
See also: [findall], page 489, [findobj], page 488, [get], page 430, [set], page 430.
findfigs ()
Find all visible figures that are currently off the screen and move them onto the screen.
See also: [allchild], page 431, [figure], page 399, [get], page 430, [set], page 430.

Figures can be printed or saved in many graphics formats with print and saveas. Occasionally, however, it may be useful to save the original Octave handle graphic directly so that further modifications can be made such as modifying a title or legend.

This can be accomplished with the following functions by

```
    fig_struct = hdl2struct (gcf);
save myplot.fig -struct fig_struct;
fig_struct = load ("myplot.fig");
struct2hdl (fig_struct);
```

$s=h d 12 s t r u c t(h)$

Return a structure, $s$, whose fields describe the properties of the object, and its children, associated with the handle, $h$.

The fields of the structure $s$ are "type", "handle", "properties", "children", and "special".

See also: [struct2hdl], page 432, [hgsave], page 417, [findobj], page 488.
$h=$ struct2hdl (s)
$h=s t r u c t 2 h d l(s, p)$
$h=s t r u c t 2 h d l(s, p, h i l e v)$
Construct a graphics handle object $h$ from the structure $s$.
The structure must contain the fields "handle", "type", "children", "properties", and "special".
If the handle of an existing figure or axes is specified, $p$, the new object will be created as a child of that object. If no parent handle is provided then a new figure and the necessary children will be constructed using the default values from the root object.
A third boolean argument hilev can be passed to specify whether the function should preserve listeners/callbacks, e.g., for legends or hggroups. The default is false.

See also: [hdl2struct], page 432, [hgload], page 418, [findobj], page 488.

```
hnew = copyobj (horig)
hnew = copyobj (horig, hparent)
```

Construct a copy of the graphic objects associated with the handles horig and return new handles hnew to the new objects.
If a parent handle hparent (root, figure, axes, or hggroup) is specified, the copied object will be created as a child of hparent.
If horig is a vector of handles, and hparent is a scalar, then each handle in the vector hnew has its "Parent" property set to hparent. Conversely, if horig is a scalar and hparent a vector, then each parent object will receive a copy of horig. If horig and hparent are both vectors with the same number of elements then hnew(i) will have parent hparent(i).

See also: [struct2hdl], page 432, [hdl2struct], page 432, [findobj], page 488.

### 15.3.3 Graphics Object Properties

In this section the graphics object properties are discussed in detail, starting with the root properties and continuing through the object hierarchy. The documentation about a specific graphics object can be displayed using doc function, e.g., doc ("axes properties") will show Section 15.3.3.3 [Axes Properties], page 441.

The allowed values for radio (string) properties can be retrieved programmatically or displayed using the one or two argument calling form of the set function. See [set], page 430.

In the following documentation default values are enclosed in $\}$.

### 15.3.3.1 Root Properties

Properties of the root graphics object:

```
beingdeleted: {"off"} | "on"
    beingdeleted is unused.
busyaction: "cancel" | {"queue"}
    busyaction is unused.
```

buttondownfen: string I function handle, def. [] (0x0)
buttondownfen is unused.
callbackobject (read-only): graphics handle, def. [] (0x0)

Graphics handle of the current object whose callback is executing.
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the root's children.
clipping: "off" | \{"on"\}
clipping is unused.
commandwindowsize (read-only): def. [0 0 ]
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this root object.
createfcn: string I function handle, def. [] (0x0)
createfcn is unused.
currentfigure: graphics handle, def. [] (0x0)
Graphics handle of the current figure.
deletefcn: string | function handle, def. [] (0x0)
deletefcn is unused.
fixedwidthfontname: string, def. "Courier"
handlevisibility: "callback" | "off" | \{"on"\}
handlevisibility is unused.
hittest: "off" | \{"on"\}
hittest is unused.
interruptible: "off" | \{"on"\}
interruptible is unused.
monitorpositions (read-only):
monitorpositions is unused.
parent: graphics handle, def. [] ( $0 \times 0$ )
Root object has no parent graphics object. parent is always empty.

```
pickableparts: "all" | "none" | {"visible"}
```

    pickableparts is unused.
    pointerlocation: two-element vector, def. [0 0 $]$
pointerlocation is unused.
pointerwindow (read-only): graphics handle, def. 0
pointerwindow is unused.
screendepth (read-only): double
screenpixelsperinch (read-only): double
screensize (read-only): four-element vector
selected: \{"off"\} | "on"
selected is unused.
selectionhighlight: "off" | \{"on"\}
selectionhighlight is unused.
showhiddenhandles: \{"off"\} | "on"
If showhiddenhandles is "on", all graphics objects handles are visible in their parents' children list, regardless of the value of their handlevisibility property.
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "root"
units: "centimeters" | "characters" | "inches" | "normalized" | \{"pixels"\} |
"points"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
visible is unused.

### 15.3.3.2 Figure Properties

Properties of figure objects (see [figure], page 399):
alphamap: def. 64-by-1 double
Transparency is not yet implemented for figure objects. alphamap is unused.
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible
property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfcn: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the figure's children.
clipping: "off" | \{"on"\}
clipping is unused.
closerequestfcn: string | function handle, def. "closereq"
Function that is executed when a figure is deleted. See [closereq function], page 405.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
color: colorspec, def. [111 $\left.1 \begin{array}{ll}1 & 1\end{array}\right]$
Color of the figure background. See Section 15.4.1 [colorspec], page 491.
colormap: N-by-3 matrix, def. 64-by-3 double
A matrix containing the RGB color map for the current axes.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this figure object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after figure has been created. Function is set by using default property on root object, e.g., set (groot, "defaultfigurecreatefcn", 'disp ("figure created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
currentaxes: graphics handle, def. [] ( $0 \times 0$ )
Handle to the graphics object of the current axes.
currentcharacter (read-only): def. ""
currentcharacter is unused.
currentobject (read-only): graphics handle, def. [] (0x0)
currentpoint (read-only): two-element vector, def. [0; 0]
A 1-by-2 matrix which holds the coordinates of the point over which the mouse pointer was when a mouse event occurred. The X and Y coordinates are in units defined by the figure's units property and their origin is the lower left corner of the plotting area.
Events which set currentpoint are
A mouse button was pressed
always

A mouse button was released
only if the figure's callback windowbuttonupf cn is defined
The pointer was moved while pressing the mouse button (drag) only if the figure's callback windowbuttonmotionfcn is defined
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before figure is deleted.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
dockcontrols: \{"off"\} | "on"
dockcontrols is unused.
filename: string, def.
The filename used when saving the plot figure.

```
graphicssmoothing: "off" | {"on"}
```

Use smoothing techniques to reduce the appearance of jagged lines.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the figure's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether figure processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 438.
integerhandle: "off" | \{"on"\}
Assign the next lowest unused integer as the Figure number.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
inverthardcopy: "off" | \{"on"\}
Replace the figure and axes background color with white when printing.
keypressfcn: string I function handle, def. [] (0x0)
Callback function executed when a keystroke event happens while the figure has focus. The actual key that was pressed can be retrieved using the second argument 'evt' of the function. For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
keyreleasefcn: string I function handle, def. [] (0x0)
With keypressfcn, the keyboard callback functions. These callback functions are called when a key is pressed/released respectively. The functions are called
with two input arguments. The first argument holds the handle of the calling figure. The second argument holds an event structure which has the following members:

Character:
The ASCII value of the key
Key: Lowercase value of the key
Modifier:
A cell array containing strings representing the modifiers pressed with the key.

For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
menubar: \{"figure"\} | "none"
Control the display of the figure menu bar at the top of the figure.
name: string, def. ""
Name to be displayed in the figure title bar. The name is displayed to the right of any title determined by the numbertitle property.
nextplot: \{"add"\} | "new" | "replace" | "replacechildren"
nextplot is used by high level plotting functions to decide what to do with axes already present in the figure. See [newplot function], page 402.
number (read-only):
numbertitle: "off" | \{"on"\}
Display "Figure" followed by the numerical figure handle value in the figure title bar.
outerposition: four-element vector, def. [-1-1-1-1]
Specify the position and size of the figure including the top menubar and the bottom status bar. The four elements of the vector are the coordinates of the lower left corner and width and height of the figure. See [units property], page 439.
paperorientation: "landscape" | \{"portrait"\}
The value for the papersize, and paperposition properties depends upon paperorientation. The horizontal and vertical values for papersize and paperposition reverse order when paperorientation is switched between "portrait" and "landscape".
paperposition: four-element vector, def. [1.3422 3.3191 5.8156 4.3617]
Vector [left bottom width height] defining the position and size of the figure (in paperunits units) on the printed page. The position [left bottom] defines the lower left corner of the figure on the page, and the size is defined by [width height]. For output formats not implicitly rendered on paper, width and height define the size of the image and the position information is ignored. Setting paperposition also forces the paperpositionmode property to be set to "manual".
paperpositionmode: \{"auto"\} | "manual"
If paperpositionmode is set to "auto", the paperposition property is automatically computed: the printed figure will have the same size as the on-screen figure and will be centered on the output page. Setting the paperpositionmode to "auto" does not modify the value of the paperposition property.
papersize: two-element vector, def. [8.5000 11.0000]
Vector [width height] defining the size of the paper for printing. Setting the papersize property to a value, not associated with one of the defined papertypes and consistent with the setting for paperorientation, forces the papertype property to the value "<custom>". If papersize is set to a value associated with a supported papertype and consistent with the paperorientation, the papertype value is modified to the associated value.

```
papertype: "<custom>" | "a" | "a0" | "a1" | "a2" | "a3" | "a4" | "a5" | "arch-a" |
"arch-b" | "arch-c" | "arch-d" | "arch-e" | "b" | "b0" | "b1" | "b2" | "b3" | "b4"
| "b5" | "c" | "d" | "e" | "tabloid" | "uslegal" | {"usletter"}
```

Name of the paper used for printed output. Setting papertype also changes papersize, while maintaining consistency with the paperorientation property.
paperunits: "centimeters" | \{"inches"\} | "normalized" | "points"
The unit used to compute the paperposition property. The conversion from physical units (e.g., "inches") is dependent on the screenpixelsperinch property of the root object.
parent: graphics handle, def. 0
Handle of the parent graphics object.
pickableparts (read-only): "all" | "none" | \{"visible"\}
pickableparts is unused.
pointer: \{"arrow"\} | "botl" | "botr" | "bottom" | "circle" | "cross" |
"crosshair" | "custom" | "fleur" | "hand" | "ibeam" | "left" | "right" | "top" |
"topl" | "topr" | "watch"
Name of the mouse pointer shape associated with the canvas of the figure. When pointer is "custom", the shape is determined by the pointershapecdata property.
pointer has no effect when the figure is in zoom, pan, or rotate mode. In this case, Octave automatically uses a pointer shape appropriate to the mode.
pointershapecdata: 16-by-16 or 32 -by- 32 Matrix, def. 16 -by-16 double
m-by-m matrix defining a custom pointer. Each element defines a pixel with the element $(1,1)$ representing the top-left pixel. A value of 1 is colored black, a value of 2 is colored white, and all other values are rendered as transparent.
pointershapehotspot: two-element vector, def. [11]
For custom pointers only pointershapehotspot defines the row and column of the pixel in pointershapecdata that is used as the pointer location.
position: four-element vector, def. [300 200560 420]
Specify the position and size of the figure canvas. The four elements of the vector are the coordinates of the lower left corner and width and height of the figure. See [units property], page 439.
renderer: \{"opengl"\} | "painters" Rendering engine used for printing when renderermode is "manual". Setting renderer also forces the renderermode property to be set to "manual".
renderermode: \{"auto"\} | "manual"
Control whether the rendering engine used for printing is chosen automatically or specified by the renderer property. See [print function], page 410.
resize: "off" | \{"on"\}
Control whether the figure can be resized by dragging the window borders and corners using a mouse. When resize is "off" mouse interactions are disabled but the figure can still be resized by changing its "position" property.
resizefcn: string | function handle, def. [] (0x0)
resizefcn is deprecated. Use sizechangedfcn instead.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
selectiontype: "alt" | "extend" | \{"normal"\} | "open"
Selection type of the latest mouse click.
selectiontype may take different values depending on the combination of mouse button and keyboard modifier that were used:
normal: Left-click.
alt: Right-click or Ctrl+Left-click.
extend: Shitf+Left-click, Middle click, or combined Left-click and Rightclick.
open: Double Left-click.
sizechangedfcn: string | function handle, def. [] ( $0 \times 0$ )
Callback triggered when the figure window size is changed.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
tag: string, def. ""
A user-defined string to label the graphics object.
toolbar: \{"auto"\} | "figure" | "none"
Control the display of the toolbar (along the bottom of the menubar) and the status bar. When set to "auto", the display is based on the value of the menubar property.
type (read-only): string
Class name of the graphics object. type is always "figure"

```
units: "centimeters" | "characters" | "inches" | "normalized" | {"pixels"} |
```

"points"

The unit used to compute the position and outerposition properties.
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the figure is not rendered on screen.
windowbuttondownfcn: string | function handle, def. [] (0x0)
See [windowbuttonupfen property], page 440.
windowbuttonmotionfcn: string | function handle, def. [] ( 0 x 0 )
See [windowbuttonupfen property], page 440.
windowbuttonupf cn: string | function handle, def. [] (0x0)
With windowbuttondownfcn and windowbuttonmotionf $c n$, the mouse callback functions. These callback functions are called when a mouse button is pressed, dragged, or released respectively. When these callback functions are executed, the currentpoint property holds the current coordinates of the cursor.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
windowkeypressfen: string | function handle, def. [] (0x0)
Function that is executed when a key is pressed and the figure has focus.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
windowkeyreleasefcn: string | function handle, def. [] (0x0)
Function that is executed when a key is released and the figure has focus.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
windowscrollwheelfcn: string | function handle, def. [] (0x0)
Function that is executed when a user manipulates the mouse wheel over this figure. The function is called with two input arguments. The first argument holds the handle of the calling figure. The second argument holds an event structure which has the following members:

VerticalScrollCount:
The number of wheel steps, typically 1 when scrolling down and -1 when scrolling up.

VerticalScrollAmount:
The number of lines a wheel step should scroll. This value is always 3.

## EventName:

The event name which is "WindowScrollWheel".
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
windowstyle: "docked" | "modal" | \{"normal"\}
The window style of a figure. One of the following values:
normal Set the window style as non modal.
modal Set the window as modal so that it will stay on top of all normal figures.
docked Setting the window style as docked currently does not dock the window.

Changing modes of a visible figure may cause the figure to close and reopen.

### 15.3.3.3 Axes Properties

Properties of axes objects (see [axes], page 423):
alim: def. [0 1]
Transparency is not yet implemented for axes objects. alim is unused.
alimmode: \{"auto"\} | "manual"
alphamap: def. [] (0x0)
alphascale: \{"linear"\} | "log"
ambientlightcolor: def. [1 1 1]
ambientlightcolor is unused.
beingdeleted: \{"off"\} | "on"
box: \{"off"\} | "on"
Control whether the axes has a surrounding box.
boxstyle: \{"back"\} | "full"
For 3-D axes, control whether the "full" box is drawn or only the 3 "back" axes
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
cameraposition: three-element vector, def. [0.5000 0.5000 9.1603]
camerapositionmode: \{"auto"\} | "manual"
cameratarget: three-element vector, def. [0.5000 0.5000 0.5000]
cameratargetmode: \{"auto"\} | "manual"
cameraupvector: three-element vector, def. [0 10 0]
cameraupvectormode: \{"auto"\} | "manual"
cameraviewangle: scalar, def. 6.6086
cameraviewanglemode: \{"auto"\} | "manual"
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the axes's children.
clim: two-element vector, def. [0 1]
Define the limits for the color axis of image children. Setting clim also forces the climmode property to be set to "manual". See [pcolor function], page 345.
climmode: \{"auto"\} | "manual"
clipping: "off" | \{"on"\}
clipping is unused.
clippingstyle: \{"3dbox"\} | "rectangle"
clippingstyle is unused.
color: colorspec, def. [llll $\left.\begin{array}{lll}1 & 1 & 1\end{array}\right]$
Color of the axes background. See Section 15.4.1 [colorspec], page 491.
colormap: def. 64-by-3 double
colororder: N-by-3 RGB matrix, def. 7-by-3 double
RGB values used by plot function for automatic line coloring.
colororderindex: def. 1
colororderindex is unused.
colorscale: \{"linear"\} | "log"
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this axes object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after axes has been created. Function is set by using default property on root object, e.g., set (groot, "defaultaxescreatefcn", 'disp ("axes created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
currentpoint: 2-by-3 matrix, def. 2-by-3 double
Matrix [xf, $\mathrm{yf}, \mathrm{zf} ; \mathrm{xb}, \mathrm{yb}, \mathrm{zb}$ ] which holds the coordinates (in axes data units) of the point over which the mouse pointer was when the mouse button was pressed. If a mouse callback function is defined, currentpoint holds the pointer coordinates at the time the mouse button was pressed. For 3-D plots, the first row of the returned matrix specifies the point nearest to the current camera position and the second row the furthest point. The two points forms a line which is perpendicular to the screen.
dataaspectratio: three-element vector, def. [1111]
Specify the relative height and width of the data displayed in the axes. Setting dataaspectratio to $[1,2]$ causes the length of one unit as displayed on the x axis to be the same as the length of 2 units on the y-axis. See [daspect function], page 383. Setting dataaspectratio also forces the dataaspectratiomode property to be set to "manual".
dataaspectratiomode: \{"auto"\} | "manual"
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before axes is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 443.
fontsizemode: \{"auto"\} | "manual"
fontsmoothing: "off" | \{"on"\}
Control whether any text associated with axes is anti-aliased.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
gridalpha: def. 0.1500
Transparency is not yet implemented for axes objects. gridalpha is unused.
gridalphamode: \{"auto"\} | "manual"
gridalphamode is unused.
gridcolor: def. [0. 15000.15000 .1500 ]
gridcolor is unused.
gridcolormode: \{"auto"\} | "manual"
gridcolormode is unused.
gridlinestyle: \{"-"\} | "--" | "-." | ":" | "none"
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the axes's handle is not visible in its parent's
"children" property.
hittest: "off" | \{"on"\}
Specify whether axes processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 445.
innerposition: four-element vector, def. [0.1300 0.1100 0.7750 0.8150]
The "innerposition" property is the same as the ["position" property], page 445.
interactions (read-only): def. [] (0x0)
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
labelfontsizemultiplier: def. 1.1000
Ratio between the $\mathrm{x} / \mathrm{y} /$ zlabel fontsize and the tick label fontsize
layer: \{"bottom"\} | "top"
Control whether the axes is drawn below child graphics objects (ticks, labels, etc. covered by plotted objects) or above.
layout (read-only): def. [] (0x0)
legend (read-only): def. [] ( $0 x 0$ )
linestyleorder: def. "-"
linestyleorder is unused.
linestyleorderindex: def. 1
linestyleorderindex is unused.
linewidth: def. 0.5000
Width of the main axes lines
minorgridalpha: def. 0.2500
Transparency is not yet implemented for axes objects. minorgridalpha is unused.
minorgridalphamode: \{"auto"\} | "manual"
minorgridalphamode is unused.
minorgridcolor: def. [0.1000 0.1000 0.1000]
minorgridcolor is unused.
minorgridcolormode: \{"auto"\} | "manual"
minorgridcolormode is unused.
minorgridlinestyle: "-" | "--" | "-." | \{":"\} | "none"
mousewheelzoom: scalar in the range $(0,1)$, def. 0.5000
Fraction of axes limits to zoom for each wheel movement.

```
nextplot: "add" | {"replace"} | "replacechildren"
    nextplot is used by high level plotting functions to decide what to do with
    graphics objects already present in the axes. See [newplot function], page 402.
    The state of nextplot is typically controlled using the hold function. See
    [hold function], page 403.
nextseriesindex (read-only): def. 1
outerposition: four-element vector, def. [0 0 1 1]
Specify the position of the plot including titles, axes, and legend. The four elements of the vector are the coordinates of the lower left corner and width and height of the plot, in units normalized to the width and height of the plot window. For example, \([0.2,0.3,0.4,0.5]\) sets the lower left corner of the axes at \((0.2,0.3)\) and the width and height to be 0.4 and 0.5 respectively. See [position property], page 445.
```

parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether axes will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the axes or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 443.
plotboxaspectratio: def. [1 111$]$
See [pbaspect function], page 384. Setting plotboxaspectratio also forces the plotboxaspectratiomode property to be set to "manual".
plotboxaspectratiomode: \{"auto"\} | "manual"
position: four-element vector, def. [0.1300 0.11000 .77500 .8150 ]
Specify the position of the plot excluding titles, axes, and legend. The four elements of the vector are the coordinates of the lower left corner and width and height of the plot, in units normalized to the width and height of the plot window. For example, $[0.2,0.3,0.4,0.5]$ sets the lower left corner of the axes at $(0.2,0.3)$ and the width and height to be 0.4 and 0.5 respectively. See [outerposition property], page 445.
positionconstraint: "innerposition" | \{"outerposition"\}
Specify which of "innerposition" or "outerposition" properties takes precedence when axes annotations extent changes. See ["innerposition" property], page 444, and ["outerposition" property], page 445.
projection: \{"orthographic"\} | "perspective"
projection is unused.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
sortmethod: "childorder" | \{"depth"\}
sortmethod is unused.
tag: string, def. ""
A user-defined string to label the graphics object.
tickdir: \{"in"\} | "out"
Control whether axes tick marks project "in" to the plot box or "out". Setting tickdir also forces the tickdirmode property to be set to "manual".
tickdirmode: \{"auto"\} | "manual"
ticklabelinterpreter: "latex" | "none" | \{"tex"\}
Control the way $x / y / z t i c k l a b e l$ properties are interpreted. See [Use of the interpreter property], page 406.
ticklength: two-element vector, def. [0.010000 0.025000]
Two-element vector [2Dlen 3Dlen] specifying the length of the tickmarks relative to the longest visible axis.
tightinset (read-only): four-element vector
Size of the [left bottom right top] margins around the axes that enclose labels and title annotations.
title: graphics handle
Graphics handle of the title text object.
titlefontsizemultiplier: positive scalar, def. 1.1000
Ratio between the title fontsize and the tick label fontsize
titlefontweight: \{"bold"\} | "normal"
Control variant of base font used for the axes title.
toolbar (read-only): def. [] (0x0)
type (read-only): string
Class name of the graphics object. type is always "axes"
units: "centimeters" | "characters" | "inches" | \{"normalized"\} | "pixels" | "points"

Units used to interpret the "position", "outerposition", and "tightinset" properties.
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
view: two-element vector, def. [0 90]
Two-element vector [azimuth elevation] specifying the viewpoint for threedimensional plots
visible: "off" | \{"on"\}
If visible is "off", the axes is not rendered on screen.
xaxis (read-only): def. [] (0x0)
xaxislocation: \{"bottom"\} | "origin" | "top"
Control the x axis location.
xcolor: $\{$ colorspec $\}$ | "none", def. [0. 1500 0.1500 0.1500]
Color of the x-axis. See Section 15.4 .1 [colorspec], page 491. Setting xcolor also forces the xcolormode property to be set to "manual".
xcolormode: \{"auto"\} | "manual"
xdir: \{"normal"\} | "reverse"
Direction of the x axis: "normal" is left to right.
xgrid: \{"off"\} | "on"
Control whether major x grid lines are displayed.
xlabel: graphics handle
Graphics handle of the x label text object.
xlim: two-element vector, def. [0 1]
Two-element vector [xmin xmax] specifying the limits for the x-axis. Setting xlim also forces the xlimmode property to be set to "manual". See [xlim function], page 351.
xlimmode: \{"auto"\} | "manual"
xminorgrid: \{"off"\} | "on"
Control whether minor x grid lines are displayed.
xminortick: \{"off"\} | "on"
xscale: \{"linear"\} | "log"
xtick: vector
Position of x tick marks. Setting xtick also forces the xtickmode property to be set to "manual".
xticklabel: string | cell array of strings, def. 1-by-6 cell
Labels of x tick marks. Setting xticklabel also forces the xticklabelmode property to be set to "manual".
xticklabelmode: \{"auto"\} | "manual"
xticklabelrotation: def. 0
xticklabelrotation is unused.
xtickmode: \{"auto"\} | "manual"
yaxis (read-only): def. [] (0x0)
yaxislocation: \{"left"\} | "origin" | "right"
Control the y-axis location.
ycolor: \{colorspec\} | "none", def. [0. 1500 0.1500 0.1500]
Color of the y-axis. See Section 15.4.1 [colorspec], page 491.
ycolormode: \{"auto"\} | "manual"
ydir: \{"normal"\} | "reverse"
Direction of the y-axis: "normal" is bottom to top.
ygrid: \{"off"\} | "on"
Control whether major y grid lines are displayed.
ylabel: graphics handle
Graphics handle of the y label text object.
ylim: two-element vector, def. [0 1]
Two-element vector [ymin ymax] specifying the limits for the y-axis. Setting ylim also forces the ylimmode property to be set to "manual". See [ylim function], page 351.

```
ylimmode: {"auto"} | "manual"
yminorgrid: {"off"} | "on"
```

Control whether minor y grid lines are displayed.
yminortick: \{"off"\} | "on"
yscale: \{"linear"\} | "log"
ytick: vector
Position of y tick marks. Setting ytick also forces the ytickmode property to be set to "manual".
yticklabel: string | cell array of strings, def. 1-by-6 cell
Labels of y tick marks. Setting yticklabel also forces the yticklabelmode property to be set to "manual".
yticklabelmode: \{"auto"\} | "manual"
yticklabelrotation: def. 0
yticklabelrotation is unused.
ytickmode: \{"auto"\} | "manual"
zaxis (read-only): def. [] (0x0)
zcolor: \{colorspec\} | "none", def. [0.1500 0.1500 0.1500]
Color of the z-axis. See Section 15.4.1 [colorspec], page 491.
zcolormode: \{"auto"\} | "manual"
zdir: \{"normal"\} | "reverse"
zgrid: \{"off"\} | "on"
Control whether major z grid lines are displayed.
zlabel: graphics handle
Graphics handle of the z label text object.
zlim: two-element vector, def. [0 1]
Two-element vector [zmin zmaz] specifying the limits for the z-axis. Setting zlim also forces the zlimmode property to be set to "manual". See [zlim function], page 351.
zlimmode: \{"auto"\} | "manual"
zminorgrid: \{"off"\} | "on"
Control whether minor z grid lines are displayed.
zminortick: \{"off"\} | "on"
zscale: \{"linear"\} | "log"
ztick: vector
Position of z tick marks. Setting ztick also forces the ztickmode property to be set to "manual".
zticklabel: string | cell array of strings, def. 1-by-6 cell
Labels of z tick marks. Setting zticklabel also forces the zticklabelmode property to be set to "manual".
zticklabelmode: \{"auto"\} | "manual"
zticklabelrotation: def. 0
zticklabelrotation is unused.
ztickmode: \{"auto"\} | "manual"

### 15.3.3.4 Legend Properties

Properties of legend objects (see [legend], page 390):
autoupdate: "off" | \{"on"\}
Control whether the number of legend items is updated automatically when objects are added to (or deleted from) the peer axes. For example:

```
## Create a single plot with its legend.
figure ();
plot (1:10);
legend ("Slope 1");
## Add another plot and specify its displayname so that
## the legend is correctly updated.
hold on;
plot ((1:10) * 2, "displayname", "Slope 2");
## Stop automatic updates for further plots.
legend ("autoupdate", "off");
plot ((1:10) * 3);
```

box: "off" | \{"on"\}

Control whether the legend has a surrounding box.
color: colorspec, def. [111 $\left.1 \begin{array}{ll}1 & 1\end{array}\right]$
Color of the legend background. See Section 15.4.1 [colorspec], page 491.
edgecolor: colorspec, def. [0.1500 0.1500 0.1500]
Control the color of the legend outline.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 9
Size of the font used for text rendering. See [fontunits property], page 449.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
location: "best" | "bestoutside" | "east" | "eastoutside" | "none" | "north" |
\{"northeast"\} | "northeastoutside" | "northoutside" | "northwest" |
"northwestoutside" | "south" | "southeast" | "southeastoutside" |
"southoutside" | "southwest" | "southwestoutside" | "west" | "westoutside"
Control the location of the legend.
orientation: "horizontal" | \{"vertical"\}
Control whether the legend items are arranged vertically (column-wise) or horizontally (row-wise).
position: four-element vector
Specify the position of the legend excluding its title. The four elements of the vector are the coordinates of the lower left corner and width and height of the legend. Changing this property also switches the "location" to "none".
string: string | cell array of strings
List of labels for the legend items. For example:

```
figure ();
plot (rand (20));
## Let legend choose names automatically
hl = legend ();
## Selectively change some names
str = get (hl, "string");
str(1:5:end) = "Garbage";
set (hl, "string", str);
Control the color of the text strings for legend items.
```

textcolor: colorspec, def. [0 0 0]
title: graphics handle
Graphics handle of the title text object.
units: "centimeters" | "characters" | "inches" | \{"normalized"\} | "pixels" |
"points"
Units used to interpret the "position", property.

### 15.3.3.5 Line Properties

Properties of line objects (see [line], page 424):
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( $0 x 0$ )
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the line is clipped in its parent axes limits.
color: colorspec, def. $\left[\begin{array}{lll}0 & 0 & 0\end{array}\right]$
Color of the line object. See Section 15.4.1 [colorspec], page 491.
contextmenu: graphics handle, def. [] ( $0 \times 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this line object.
createfcn: string I function handle, def. [] (0x0)
Callback function executed immediately after line has been created. Function is set by using default property on root object, e.g., set (groot, "defaultlinecreatefcn", 'disp ("line created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before line is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
displayname: string | cell array of strings, def. ""
Text for the legend entry corresponding to this line.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the line's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether line processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 452.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
linejoin: "chamfer" | "miter" | \{"round"\}
Control the shape of the junction of line segments. This property currently only affects the printed output.
linestyle: \{"-"\} | "--" | "-." | ":" | "none"
See Section 15.4.2 [Line Styles], page 492.
linewidth: def. 0.5000
Width of the line object measured in points.
marker: "*" | "+" | "." | "<" | ">" | "~" | "_" | "d" | "diamond" | "h" | "hexagram"
| \{"none"\} | "o" | "p" | "pentagram" | "s" | "square" | "v" | "x" | "|"
Shape of the marker for each data point. See Section 15.4.3 [Marker Styles], page 492.
markeredgecolor: \{"auto"\} | "none"
Color of the edge of the markers. When set to "auto", the marker edges have the same color as the line. If set to "none", no marker edges are displayed. This property can also be set to any color. See Section 15.4.1 [colorspec], page 491.
markerfacecolor: "auto" | \{"none"\}
Color of the face of the markers. When set to "auto", the marker faces have the same color as the line. If set to "none", the marker faces are not displayed. This property can also be set to any color. See Section 15.4.1 [colorspec], page 491.
markersize: scalar, def. 6
Size of the markers measured in points.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether line will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the line or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one.
When an object is configured to accept mouse clicks the "hittest" property
will determine how they are processed. See [hittest property], page 451.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "line"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the line is not rendered on screen.
xdata: vector, def. [0 1]
Vector of x data to be plotted.
xdatasource: string, def. ""
Name of a vector in the current base workspace to use as x data.
ydata: vector, def. [01]
Vector of y data to be plotted.
ydatasource: string, def. ""
Name of a vector in the current base workspace to use as y data.
zdata: vector, def. [] ( 0 x 0 )
Vector of z data to be plotted.
zdatasource: string, def. ""
Name of a vector in the current base workspace to use as z data.

### 15.3.3.6 Text Properties

Properties of text objects (see [text], page 392):
backgroundcolor: colorspec, def. "none"
Color of the background area. See Section 15.4.1 [colorspec], page 491.
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfcn: string I function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the text is clipped in its parent axes limits.
color: colorspec, def. [1000]
Color of the text. See Section 15.4.1 [colorspec], page 491.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this text object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after text has been created. Function is set by using default property on root object, e.g., set (groot, "defaulttextcreatefcn", 'disp ("text created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before text is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
edgecolor: colorspec, def. "none"
Color of the outline of the background area. See Section 15.4.1 [colorspec], page 491.
editing: \{"off"\} | "on"
editing is unused.
extent (read-only): four-element vector
Vector [x0 y0 width height] indicating the size and location of the text string.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 454.
fontsmoothing: "off" | \{"on"\}
Control whether anti-aliasing is used when rendering text.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the text's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether text processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 455.
horizontalalignment: "center" | \{"left"\} | "right"
interpreter: "latex" | "none" | \{"tex"\}
Control the way the "string" property is interpreted. See [Use of the interpreter property], page 406.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
linestyle: \{"-"\} | "--" | "-." | ":" | "none"
Style of the outline. See Section 15.4.2 [Line Styles], page 492.
linewidth: scalar, def. 0.5000
Width of the outline.
margin: scalar, def. 3
Margins between the borders of the background area and the texts. The value is currently interpreted as pixels, regardless of the "fontunits" property.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether text will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the text or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 454.
position: three-element vector, def. [000]
Vector [X0 Y0 Z0] where X0, Y0, and Z0 indicate the position of the text anchor as defined by verticalalignment and horizontalalignment.
rotation: scalar, def. 0
The angle of rotation for the displayed text, measured in degrees.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
string: string, def. ""
The text object string content.
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "text"
units: "centimeters" | \{"data"\} | "inches" | "normalized" | "pixels" | "points" userdata: Any Octave data, def. [] (0x0)

User-defined data to associate with the graphics object.

```
verticalalignment: "baseline" | "bottom" | "cap" | {"middle"} | "top"
visible: "off" | {"on"}
    If visible is "off", the text is not rendered on screen.
```


### 15.3.3.7 Image Properties

Properties of image objects (see [image], page 866):
alphadata: scalar | matrix, def. 1
Transparency is not yet implemented for image objects. alphadata is unused.
alphadatamapping: "direct" | \{"none"\} | "scaled"
Transparency is not yet implemented for image objects. alphadatamapping is unused.
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( $0 \times 0$ )
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
cdata: matrix, def. 64-by-64 double
cdatamapping: \{"direct"\} | "scaled"
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the image is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this image object.
createfcn: string | function handle, def. [] ( $0 \times \mathrm{x} 0$ )
Callback function executed immediately after image has been created. Function is set by using default property on root object, e.g., set (groot, "defaultimagecreatefcn", 'disp ("image created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before image is deleted.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the image's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether image processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 457.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of
drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether image will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the image or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 456.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "image"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the image is not rendered on screen.
xdata: two-element vector, def. [1 64]
Two-element vector [xfirst xlast] specifying the x coordinates of the centers of the first and last columns of the image.

Setting xdata to the empty matrix ([]) will restore the default value of [1 columns(image)].
ydata: two-element vector, def. [1 64]
Two-element vector [yfirst ylast] specifying the y coordinates of the centers of the first and last rows of the image.
Setting ydata to the empty matrix ([]) will restore the default value of [1 rows (image)].

### 15.3.3.8 Patch Properties

Properties of patch objects (see [patch], page 425):
alphadatamapping: "direct" | "none" | \{"scaled"\}
Transparency is not yet implemented for patch objects. alphadatamapping is unused.
ambientstrength: scalar, def. 0.3000
Strength of the ambient light. Value between 0.0 and 1.0
backfacelighting: "lit" | \{"reverselit"\} | "unlit"
"lit": The normals are used as is for lighting. "reverselit": The normals are always oriented towards the point of view. "unlit": Faces with normals pointing away from the point of view are unlit.
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
cdata: scalar | matrix, def. [] (0x0)
Data defining the patch object color. Patch color can be defined for faces or for vertices.
If cdata is a scalar index into the current colormap or a RGB triplet, it defines the color of all faces.

If cdata is an N-by-1 vector of indices or an N-by-3 (RGB) matrix, it defines the color of each one of the N faces.
If cdata is an N-by-M or an N-by-M-by-3 (RGB) matrix, it defines the color at each vertex.
cdatamapping: "direct" | \{"scaled"\}
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the patch is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] ( $0 \times 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this patch object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after patch has been created. Function is set by using default property on root object, e.g., set (groot, "defaultpatchcreatefcn", 'disp ("patch created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] ( $0 \times \mathrm{x} 0$ )
Callback function executed immediately before patch is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
diffusestrength: scalar, def. 0.6000
Strength of the diffuse reflex. Value between 0.0 (no diffuse reflex) and 1.0 (full diffuse reflex).
displayname: def. ""
Text of the legend entry corresponding to this patch.
edgealpha: scalar | matrix, def. 1
Transparency is not yet implemented for patch objects. edgealpha is unused.
edgecolor: def. [0 00 0]
edgelighting: "flat" | "gouraud" | \{"none"\} | "phong"
When set to a value other than "none", the edges of the object are drawn with light and shadow effects. Supported values are "none" (no lighting effects), "flat" (facetted look) and "gouraud" (linear interpolation of the lighting effects between the vertices). "phong" is deprecated and has the same effect as "gouraud".
facealpha: scalar | "flat" | "interp", def. 1
Transparency level of the faces of the patch object. Only double values are supported at present where a value of 0 means complete transparency and a value of 1 means solid faces without transparency. Setting the property to "flat" or "interp" causes the faces to not being rendered. Additionally, the faces are not sorted from back to front which might lead to unexpected results when rendering layered transparent faces.
facecolor: \{colorspec\} | "none" | "flat" | "interp", def. [0 00 0 $]$
facelighting: \{"flat"\} | "gouraud" | "none" | "phong"
When set to a value other than "none", the faces of the object are drawn with light and shadow effects. Supported values are "none" (no lighting effects), "flat" (facetted look) and "gouraud" (linear interpolation of the lighting effects between the vertices). "phong" is deprecated and has the same effect as "gouraud".
facenormals: def. [] (0x0)
Face normals are used for lighting the edges or faces if the edgelighting or facelighting properties are set to "flat". Setting facenormals also forces the facenormalsmode property to be set to "manual"
facenormalsmode: \{"auto"\} | "manual"
If this property is set to "auto", facenormals are automatically calculated if the edgelighting or facelighting property are set to "flat" and at least one light object is present and visible in the same axes.
faces: vector | matrix, def. [1 2 3]
facevertexalphadata: scalar | matrix, def. [] (0x0)
Transparency is not yet implemented for patch objects. facevertexalphadata is unused.
facevertexcdata: def. [] (0x0)
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the patch's handle is not visible in its parent's "children" property.

```
hittest: "off" | \{"on"\}
```

Specify whether patch processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 460.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
linestyle: \{"-"\} | "--" | "-." | ":" | "none"
linewidth: def. 0.5000
marker: "*" | "+" | "." | "<" | ">" | "~" | "_" | "d" | "diamond" | "h" | "hexagram"
| \{"none"\} | "o" | "p" | "pentagram" | "s" | "square" | "v" | "x" | " |"
See [line marker property], page 451.
markeredgecolor: \{"auto"\} | "flat" | "none"
See [line markeredgecolor property], page 451.
markerfacecolor: "auto" | "flat" | \{"none"\}
See [line markerfacecolor property], page 452.
markersize: scalar, def. 6
See [line markersize property], page 452.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether patch will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the patch or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 459.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
specularcolorreflectance: scalar, def. 1
Reflectance for specular color. Value between 0.0 (color of underlying face) and 1.0 (color of light source).
specularexponent: scalar, def. 10
Exponent for the specular reflex. The lower the value, the more the reflex is spread out.
specularstrength: scalar, def. 0.9000
Strength of the specular reflex. Value between 0.0 (no specular reflex) and 1.0 (full specular reflex).
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "patch"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
vertexnormals: def. [] (0x0)
Vertex normals are used for lighting the edges or faces if the edgelighting or facelighting properties are set to "gouraud". Setting vertexnormals also forces the vertexnormalsmode property to be set to "manual"
vertexnormalsmode: \{"auto"\} | "manual"
If this property is set to "auto", vertexnormals are automatically calculated if the edgelighting or facelighting property are set to "gouraud" and at least one light object is present and visible in the same axes.
vertices: vector | matrix, def. 3-by-2 double
visible: "off" | \{"on"\}
If visible is "off", the patch is not rendered on screen.
xdata: vector | matrix, def. [0; 1; 0]
ydata: vector I matrix, def. [1; 1; 0]
zdata: vector I matrix, def. [] (0x0)

### 15.3.3.9 Scatter Properties

Properties of scatter objects (see [scatter], page 327):
annotation: def. [] ( $0 x 0$ )
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
cdata: scalar | matrix, def. [0 0.44700 .7410 ]
Data defining the scatter object color.
If cdata is a scalar index into the current colormap or a RGB triplet, it defines the color of all scatter markers.

If cdata is an N -by- 1 vector of indices or an N -by- 3 (RGB) matrix, it defines the color of each one of the N scatter markers.
cdatamode: \{"auto"\} | "manual"
If cdatamode is "auto", cdata is set to the color from the colororder of the ancestor axes corresponding to the seriesindex.
cdatasource: def. ""
Data from workspace variables is not yet implemented for scatter objects. cdatasource is unused.
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the scatter is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this scatter object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after scatter has been created. Function is set by using default property on root object, e.g., set (groot, "defaultscattercreatefcn", 'disp ("scatter created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
datatiptemplate: def. [] (0x0)
deletefcn: string | function handle, def. [] ( $0 x 0$ )
Callback function executed immediately before scatter is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
displayname: def. ""
Text of the legend entry corresponding to this scatter object.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the scatter's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether scatter processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 463.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.

```
latitudedata: def. [] (0x0)
latitudedatasource: def. ""
linewidth: def. 0.5000
```

Line width of the edge of the markers.

```
longitudedata: def. [] (0x0)
```

longitudedatasource: def. ""
marker: "*" | "+" | "." | "<" | ">" | "~" | "_" | "d" | "diamond" | "h" | "hexagram"
| "none" | \{"o"\} | "p" | "pentagram" | "s" | "square" | "v" | "x" | "|"
See [line marker property], page 451.
markeredgealpha: scalar, def. 1
Transparency level of the faces of the markers where a value of 0 means complete transparency and a value of 1 means solid faces without transparency. Note that the markers are not sorted from back to front which might lead to unexpected results when rendering layered transparent markers or in combination with other transparent objects.
markeredgecolor: \{"none"\} | \{"flat"\} | colorspec
Color of the edge of the markers. "none" means that the edges are transparent and "flat" means that the value from cdata is used. See [line markeredgecolor property], page 451.
markerfacealpha: scalar, def. 1
Transparency level of the faces of the markers where a value of 0 means complete transparency and a value of 1 means solid faces without transparency. Note that the markers are not sorted from back to front which might lead to unexpected results when rendering layered transparent markers or in combination with other transparent objects.
markerfacecolor: \{\{"none"\}\} | "flat" | "auto" | colorspec
Color of the face of the markers. "none" means that the faces are transparent, "flat" means that the value from cdata is used, and "auto" uses the color property of the ancestor axes. See [line markerfacecolor property], page 452.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether scatter will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the scatter or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 462.
rdata: def. [] (0x0)
rdatasource: def. ""
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
seriesindex: def. 1
Each scatter object in the same axes is asigned an incrementing integer. This corresponds to the index into the colororder of the ancestor axes that is used if cdatamode is set to "auto".
sizedata: [] | scalar | vector, def. [] (0x0)
Size of the area of the marker. A scalar value applies to all markers. If cdata is an N-by- 1 vector, it defines the color of each one of the N scatter markers.
sizedatasource: def. ""
Data from workspace variables is not yet implemented for scatter objects. sizedatasource is unused.
tag: string, def. ""
A user-defined string to label the graphics object.
thetadata: def. [] ( $0 x 0$ )
thetadatasource: def. ""
type (read-only): string
Class name of the graphics object. type is always "scatter"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the scatter is not rendered on screen.
xdata: vector, def. [] (0x0)
Vector with the x coordinates of the scatter object.
xdatasource: def. ""
Data from workspace variables is not yet implemented for scatter objects. xdatasource is unused.
ydata: vector, def. [] (0x0)
Vector with the y coordinates of the scatter object.
ydatasource: def. ""
Data from workspace variables is not yet implemented for scatter objects. ydatasource is unused.
zdata: [] | vector, def. [] (0x0)
For 3D data, vector with the y coordinates of the scatter object.
zdatasource: def. ""
Data from workspace variables is not yet implemented for scatter objects. zdatasource is unused.

### 15.3.3.10 Surface Properties

Properties of surface objects (see [surface], page 426):
alphadata: scalar | matrix, def. 1
Transparency is not yet implemented for surface objects. alphadata is unused.
alphadatamapping: "direct" | "none" | \{"scaled"\}
Transparency is not yet implemented for surface objects. alphadatamapping is unused.
ambientstrength: scalar, def. 0.3000
Strength of the ambient light. Value between 0.0 and 1.0
backfacelighting: "lit" | \{"reverselit"\} | "unlit"
"lit": The normals are used as is for lighting. "reverselit": The normals are always oriented towards the point of view. "unlit": Faces with normals pointing away from the point of view are unlit.
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfcn: string I function handle, def. [] ( 0 x 0 )
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
cdata: matrix, def. 3-by-3 double
cdatamapping: "direct" | \{"scaled"\}
cdatasource: def. ""
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the surface is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] ( $0 \times 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this surface object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after surface has been created. Function is set by using default property on root object, e.g., set (groot, "defaultsurfacecreatefcn", 'disp ("surface created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before surface is deleted.

For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
diffusestrength: scalar, def. 0.6000
Strength of the diffuse reflex. Value between 0.0 (no diffuse reflex) and 1.0 (full diffuse reflex).
displayname: def. ""
Text for the legend entry corresponding to this surface.
edgealpha: scalar, def. 1
Transparency is not yet implemented for surface objects. edgealpha is unused.
edgecolor: def. [ 0000$]$
edgelighting: "flat" | "gouraud" | \{"none"\} | "phong"
When set to a value other than "none", the edges of the object are drawn with light and shadow effects. Supported values are "none" (no lighting effects), "flat" (facetted look) and "gouraud" (linear interpolation of the lighting effects between the vertices). "phong" is deprecated and has the same effect as "gouraud".
facealpha: scalar | "flat" | "interp" | "texturemap", def. 1
Transparency level of the faces of the surface object. Only double values are supported at present where a value of 0 means complete transparency and a value of 1 means solid faces without transparency. Setting the property to "flat", "interp" or "texturemap" causes the faces to not being rendered. Additionally, the faces are not sorted from back to front which might lead to unexpected results when rendering layered transparent faces.
facecolor: \{"flat"\} | "interp" | "none" | "texturemap"
facelighting: \{"flat"\} | "gouraud" | "none" | "phong"
When set to a value other than "none", the faces of the object are drawn with light and shadow effects. Supported values are "none" (no lighting effects), "flat" (facetted look) and "gouraud" (linear interpolation of the lighting effects between the vertices). "phong" is deprecated and has the same effect as "gouraud".
facenormals: def. [] (0x0)
Face normals are used for lighting the edges or faces if the edgelighting or facelighting properties are set to "flat". Setting facenormals also forces the facenormalsmode property to be set to "manual"
facenormalsmode: \{"auto"\} | "manual"
If this property is set to "auto", facenormals are automatically calculated if the edgelighting or facelighting property are set to "flat" and at least one light object is present and visible in the same axes.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the surface's handle is not visible in its parent's "children" property.

```
hittest: "off" | {"on"}
```

Specify whether surface processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 467.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
linestyle: \{"-"\} | "--" | "-." | ":" | "none"
See Section 15.4.2 [Line Styles], page 492.
linewidth: def. 0.5000
See [line linewidth property], page 451.
marker: "*" | "+" | "." | "<" | ">" | "~" | "_" | "d" | "diamond" | "h" | "hexagram"
| \{"none"\} | "o" | "p" | "pentagram" | "s" | "square" | "v" | "x" | " |"
See Section 15.4.3 [Marker Styles], page 492.
markeredgecolor: \{"auto"\} | "flat" | "none"
See [line markeredgecolor property], page 451.
markerfacecolor: "auto" | "flat" | \{"none"\}
See [line markerfacecolor property], page 452.
markersize: scalar, def. 6
See [line markersize property], page 452.
meshstyle: \{"both"\} | "column" | "row"
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether surface will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the surface or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 466.

```
selected: {"off"} | "on"
selectionhighlight: "off" | {"on"}
specularcolorreflectance: scalar, def. 1
```

Reflectance for specular color. Value between 0.0 (color of underlying face) and 1.0 (color of light source).
specularexponent: scalar, def. 10
Exponent for the specular reflex. The lower the value, the more the reflex is spread out.
specularstrength: scalar, def. 0.9000
Strength of the specular reflex. Value between 0.0 (no specular reflex) and 1.0 (full specular reflex).
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "surface"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
vertexnormals: def. [] (0x0)
Vertex normals are used for lighting the edges or faces if the edgelighting or facelighting properties are set to "gouraud". Setting vertexnormals also forces the vertexnormalsmode property to be set to "manual"
vertexnormalsmode: \{"auto"\} | "manual"
If this property is set to "auto", vertexnormals are automatically calculated if the edgelighting or facelighting property are set to "gouraud" and at least one light object is present and visible in the same axes.
visible: "off" | \{"on"\}
If visible is "off", the surface is not rendered on screen.
xdata: matrix, def. [1 2 3]
xdatasource: def. ""
ydata: matrix, def. [1; 2; 3]
ydatasource: def. ""
zdata: matrix, def. 3-by-3 double
zdatasource: def. ""

### 15.3.3.11 Light Properties

Properties of light objects (see [light], page 426):

```
beingdeleted: {"off"} | "on"
```

busyaction: "cancel" | \{"queue"\}

Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
children is unused.
clipping: "off" | \{"on"\}
If clipping is "on", the light is clipped in its parent axes limits.
color: colorspec, def. [11ll $\left.1 \begin{array}{ll}1 & 1\end{array}\right]$
Color of the light source. See Section 15.4.1 [colorspec], page 491.
contextmenu: graphics handle, def. [] ( $0 \times 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this light object.
createfcn: string I function handle, def. [] ( $0 \times 0$ )
Callback function executed immediately after light has been created. Function is set by using default property on root object, e.g., set (groot, "defaultlightcreatefcn", 'disp ("light created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before light is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the light's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether light processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 469.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether light will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the light or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one.

When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 469.

## position: def. [lll 101$]$

Position of the light source.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
style: \{"infinite"\} | "local"
This string defines whether the light emanates from a light source at infinite distance ("infinite") or from a local point source ("local").
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "light"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.

```
visible: "off" | {"on"}
```

If visible is "off", the light is not rendered on screen.

### 15.3.3.12 Uimenu Properties

Properties of uimenu objects (see [uimenu], page 930):
accelerator: def. ""
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
buttondownfen is unused.
callback: def. [] ( 0 x0)
checked: \{"off"\} | "on"
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uimenu's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uimenu is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] ( 0 x 0 )
Graphics handle of the uicontextmenu object that is currently associated to this uimenu object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uimenu has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuimenucreatefcn", 'disp ("uimenu created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before uimenu is deleted.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
enable: "off" | \{"on"\}
foregroundcolor: def. [000]
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uimenu's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether uimenu processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 471.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
label: def. ""
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uimenu will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uimenu or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 471.

```
position: def. 0
selected: {"off"} | "on"
selectionhighlight: "off" | {"on"}
separator:{"off"} | "on"
tag: string, def. ""
```

A user-defined string to label the graphics object.
text: def. ""
type (read-only): string
Class name of the graphics object. type is always "uimenu"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uimenu is not rendered on screen.

### 15.3.3.13 Uibuttongroup Properties

Properties of uibuttongroup objects (see [uibuttongroup], page 924):
backgroundcolor: def. [0.9400 0.9400 0.9400]
beingdeleted: \{"off"\} | "on"
bordertype: "beveledin" | "beveledout" | \{"etchedin"\} | "etchedout" | "line" |
"none"
borderwidth: def. 1
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( $0 \times \mathrm{x} 0$ )
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uibuttongroup's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uibuttongroup is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] ( $0 \times x 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this uibuttongroup object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uibuttongroup has been
created. Function is set by using default property on root object, e.g., set
(groot, "defaultuibuttongroupcreatefcn", 'disp ("uibuttongroup
created!")').
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uibuttongroup is deleted.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 473.

```
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | {"points"}
```

Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
foregroundcolor: def. [0 000 ]
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uibuttongroup's handle is not visible in its parent's "children" property.
highlightcolor: def. [1111]
hittest: "off" | \{"on"\}
Specify whether uibuttongroup processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 473.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uibuttongroup will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uibuttongroup or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured
to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 473.
position: def. [0 $\left.0 \begin{array}{lll}0 & 1 & 1\end{array}\right]$
resizefcn: string | function handle, def. [] (0x0)
resizefcn is deprecated. Use sizechangedfen instead.
selected: \{"off"\} | "on"
selectedobject: def. [] (0x0)
selectionchangedfcn: def. [] (0x0)
selectionhighlight: "off" | \{"on"\}
shadowcolor: def. [0.7000 0.7000 0.7000]
sizechangedfcn: string | function handle, def. [] (0x0)
Callback triggered when the uibuttongroup size is changed.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
tag: string, def. ""
A user-defined string to label the graphics object.
title: def. ""
titleposition: "centerbottom" | "centertop" | "leftbottom" | \{"lefttop"\} |
"rightbottom" | "righttop"
type (read-only): string
Class name of the graphics object. type is always "uibuttongroup"
units: "centimeters" | "characters" | "inches" | \{"normalized"\} | "pixels" |
"points"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uibuttongroup is not rendered on screen.

### 15.3.3.14 Uicontextmenu Properties

Properties of uicontextmenu objects (see [uicontextmenu], page 931):
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( 0 x 0 )
buttondownfen is unused.
callback: def. [] ( $0 x 0$ )
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uicontextmenu's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uicontextmenu is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] ( 0 x 0 )
Graphics handle of the uicontextmenu object that is currently associated to this uicontextmenu object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uicontextmenu has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuicontextmenucreatefcn", 'disp ("uicontextmenu created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uicontextmenu is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uicontextmenu's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether uicontextmenu processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 475.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uicontextmenu will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uicontextmenu or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 475.
position: def. [0 0]
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
tag: string, def. ""
A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "uicontextmenu"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uicontextmenu is not rendered on screen.

### 15.3.3.15 Uipanel Properties

Properties of uipanel objects (see [uipanel], page 924):
backgroundcolor: def. [0.9400 0.94000 .9400 ]
beingdeleted: \{"off"\} | "on"
bordertype: "beveledin" | "beveledout" | \{"etchedin"\} | "etchedout" | "line" |
"none"
borderwidth: def. 1
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( $0 \times \mathrm{x} 0$ )
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uipanel's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uipanel is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this uipanel object.
createfcn: string | function handle, def. [] ( $0 x 0$ )
Callback function executed immediately after uipanel has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuipanelcreatefcn", 'disp ("uipanel created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uipanel is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 477.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
foregroundcolor: def. [0 00 0 $]$
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uipanel's handle is not visible in its parent's "children" property.
highlightcolor: def. [1111]
hittest: "off" | \{"on"\}
Specify whether uipanel processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 477.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uipanel will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uipanel or its children may react to
mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 477.
position: def. [0 0011 1]
resizefcn: string | function handle, def. [] (0x0)
resizefcn is deprecated. Use sizechangedfen instead.
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
shadowcolor: def. [0.7000 0.7000 0.7000]
sizechangedfcn: string | function handle, def. [] (0x0)
Callback triggered when the uipanel size is changed.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
tag: string, def. ""
A user-defined string to label the graphics object.
title: def. ""
titleposition: "centerbottom" | "centertop" | "leftbottom" | \{"lefttop"\} |
"rightbottom" | "righttop"
type (read-only): string
Class name of the graphics object. type is always "uipanel"
units: "centimeters" | "characters" | "inches" | \{"normalized"\} | "pixels" |
"points"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uipanel is not rendered on screen.

### 15.3.3.16 Uicontrol Properties

Properties of uicontrol objects (see [uicontrol], page 925):
backgroundcolor: def. [0.9400 0.94000 .9400 ]
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
callback: def. [] (0x0)
cdata: def. [] (0x0)
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uicontrol's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uicontrol is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this uicontrol object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uicontrol has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuicontrolcreatefcn", 'disp ("uicontrol created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uicontrol is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
enable: "inactive" | "off" | \{"on"\}
extent (read-only): four-element vector
Size of the text string associated to the uicontrol returned in the form [00 width height] (the two first elements are always zero).
For multi-line strings the returned width and height indicate the size of the rectangle enclosing all lines.
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 479.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
foregroundcolor: def. [0 000 ]
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uicontrol's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether uicontrol processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 480.
horizontalalignment: \{"center"\} | "left" | "right"
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
keypressfcn: def. [] (0x0)
listboxtop: def. 1
max: def. 1
min: def. 0
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uicontrol will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uicontrol or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 480.
position: def. [0 080 30]
selected: \{"off"\} | "on"
selectionhighlight: "off" | \{"on"\}
sliderstep: def. [0.010000 0.100000]
string: def. ""
style: "checkbox" | "edit" | "frame" | "listbox" | "popupmenu" | \{"pushbutton"\}
| "radiobutton" | "slider" | "text" | "togglebutton"
tag: string, def. ""
A user-defined string to label the graphics object.
tooltipstring: def. ""
type (read-only): string
Class name of the graphics object. type is always "uicontrol"

```
units: "centimeters" | "characters" | "inches" | "normalized" | {"pixels"} |
"points"
userdata: Any Octave data, def. [] (0x0)
    User-defined data to associate with the graphics object.
value: def. 0
verticalalignment: "bottom" | {"middle"} | "top"
visible: "off" | {"on"}
    If visible is "off", the uicontrol is not rendered on screen.
```


### 15.3.3.17 Uitable Properties

Properties of uitable objects (see [uitable], page 926):
backgroundcolor: def. 2-by-3 double
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] (0x0)
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
celleditcallback: def. [] (0x0)
cellselectioncallback: def. [] ( $0 x 0$ )
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uitable's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uitable is clipped in its parent axes limits.
columneditable: def. [] (0x0)
columnformat: def. \{\} (0x0)
columnname: def. "numbered"
columnwidth: def. "auto"
contextmenu: graphics handle, def. [] ( $0 \times 0$ )
Graphics handle of the uicontextmenu object that is currently associated to this uitable object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uitable has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuitablecreatefcn", 'disp ("uitable created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
data: def. [] (0x0)
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uitable is deleted.
For information on how to write graphics listener functions see Section 15.4.4
[Callbacks section], page 492.
enable: "off" | \{"on"\}
extent (read-only): four-element vector
fontangle: "italic" | \{"normal"\}
Control whether the font is italic or normal.
fontname: string, def. "*"
Name of font used for text rendering. When setting this property, the text rendering engine will search for a matching font in your system. If none is found then text is rendered using a default sans serif font (same as the default "*" value).
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.
fontsize: scalar, def. 10
Size of the font used for text rendering. See [fontunits property], page 482.
fontunits: "centimeters" | "inches" | "normalized" | "pixels" | \{"points"\}
Units used to interpret the "fontsize" property.
fontweight: "bold" | \{"normal"\}
Control the variant of the base font used for text rendering.
foregroundcolor: def. [000]
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uitable's handle is not visible in its parent's
"children" property.
hittest: "off" | \{"on"\}
Specify whether uitable processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 482.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
keypressfcn: def. [] (0x0)
keyreleasef cn: def. [] (0x0)
parent: graphics handle
Handle of the parent graphics object.

```
pickableparts: "all" | "none" | {"visible"}
Specify whether uitable will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uitable or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 482.
```

```
position: def. [20 20 300 300]
rearrangeablecolumns:{"off"} | "on"
rowname: def. "numbered"
rowstriping: "off" | {"on"}
selected: {"off"} | "on"
selectionhighlight: "off" | {"on"}
tag: string, def. ""
```

A user-defined string to label the graphics object.

```
tooltipstring: def. ""
```

type (read-only): string

Class name of the graphics object. type is always "uitable"

```
units: "centimeters" | "characters" | "inches" | "normalized" | {"pixels"} |
"points"
userdata: Any Octave data, def. [] (0x0)
User-defined data to associate with the graphics object.
```

```
visible: "off" | {"on"}
```

If visible is "off", the uitable is not rendered on screen.

### 15.3.3.18 Uitoolbar Properties

Properties of uitoolbar objects (see [uitoolbar], page 931):

```
beingdeleted: {"off"} | "on"
```

busyaction: "cancel" | \{"queue"\}

Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string I function handle, def. [] (0x0) buttondownfen is unused.
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uitoolbar's children.
clipping: "off" | \{"on"\}
If clipping is "on", the uitoolbar is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this uitoolbar object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uitoolbar has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuitoolbarcreatefcn", 'disp ("uitoolbar created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uitoolbar is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uitoolbar's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether uitoolbar processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 484.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uitoolbar will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uitoolbar or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 484.

```
selected: {"off"} | "on"
selectionhighlight: "off" | {"on"}
tag: string, def. ""
```

A user-defined string to label the graphics object.
type (read-only): string
Class name of the graphics object. type is always "uitoolbar"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uitoolbar is not rendered on screen.

### 15.3.3.19 Uipushtool Properties

Properties of uipushtool objects (see [uipushtool], page 932):
__named_icon__: def. ""
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfcn: string I function handle, def. [] (0x0)
buttondownfen is unused.
cdata: def. [] (0x0)
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uipushtool's children.
clickedcallback: def. [] (0x0)
clipping: "off" | \{"on"\}
If clipping is "on", the uipushtool is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this uipushtool object.
createfcn: string I function handle, def. [] (0x0)
Callback function executed immediately after uipushtool has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuipushtoolcreatefcn", 'disp ("uipushtool created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string | function handle, def. [] (0x0)
Callback function executed immediately before uipushtool is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
enable: "off" | \{"on"\}
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uipushtool's handle is not visible in its parent's "children" property.

```
hittest: "off" | {"on"}
```

Specify whether uipushtool processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 486.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uipushtool will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uipushtool or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 485.

```
selected: {"off"} | "on"
selectionhighlight: "off" | {"on"}
separator:{"off"} | "on"
tag: string, def. ""
```

A user-defined string to label the graphics object.
tooltipstring: def. ""
type (read-only): string
Class name of the graphics object. type is always "uipushtool"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uipushtool is not rendered on screen.

### 15.3.3.20 Uitoggletool Properties

Properties of uitoggletool objects (see [uitoggletool], page 933):
__named_icon__: def. ""
beingdeleted: \{"off"\} | "on"
busyaction: "cancel" | \{"queue"\}
Define how Octave handles the execution of this object's callback properties when it is unable to interrupt another object's executing callback. This is only
relevant when the currently executing callback object has its interruptible property set to "off". The busyaction property of the interrupting callback object indicates whether the interrupting callback is queued ("queue" (default)) or discarded ("cancel"). See Section 15.4.4 [Callbacks section], page 492.
buttondownfen: string | function handle, def. [] ( 0 x 0 ) buttondownfen is unused.
cdata: def. [] (0x0)
children (read-only): vector of graphics handles, def. [] (0x1)
Graphics handles of the uitoggletool's children.
clickedcallback: def. [] (0x0)
clipping: "off" | \{"on"\}
If clipping is "on", the uitoggletool is clipped in its parent axes limits.
contextmenu: graphics handle, def. [] (0x0)
Graphics handle of the uicontextmenu object that is currently associated to this uitoggletool object.
createfcn: string | function handle, def. [] (0x0)
Callback function executed immediately after uitoggletool has been created. Function is set by using default property on root object, e.g., set (groot, "defaultuitoggletoolcreatefcn", 'disp ("uitoggletool created!")').
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
deletefcn: string I function handle, def. [] (0x0)
Callback function executed immediately before uitoggletool is deleted.
For information on how to write graphics listener functions see Section 15.4.4 [Callbacks section], page 492.
enable: "off" | \{"on"\}
handlevisibility: "callback" | "off" | \{"on"\}
If handlevisibility is "off", the uitoggletool's handle is not visible in its parent's "children" property.
hittest: "off" | \{"on"\}
Specify whether uitoggletool processes mouse events or passes them to ancestors of the object. When enabled, the object may respond to mouse clicks by evaluating the "buttondownfcn", showing the uicontextmenu, and eventually becoming the root "currentobject". This property is only relevant when the object can accept mouse clicks which is determined by the "pickableparts" property. See [pickableparts property], page 488.
interruptible: "off" | \{"on"\}
Specify whether this object's callback functions may be interrupted by other callbacks. By default interruptible is "on" and callbacks that make use of drawnow, figure, waitfor, getframe or pause functions are eventually interrupted. See Section 15.4.4 [Callbacks section], page 492.
offcallback: def. [] (0x0)
oncallback: def. [] (0x0)
parent: graphics handle
Handle of the parent graphics object.
pickableparts: "all" | "none" | \{"visible"\}
Specify whether uitoggletool will accept mouse clicks. By default, pickableparts is "visible" and only visible parts of the uitoggletool or its children may react to mouse clicks. When pickableparts is "all" both visible and invisible parts (or children) may react to mouse clicks. When pickableparts is "none" mouse clicks on the object are ignored and transmitted to any objects underneath this one. When an object is configured to accept mouse clicks the "hittest" property will determine how they are processed. See [hittest property], page 487.

```
selected: {"off"} | "on"
```

selectionhighlight: "off" | \{"on"\}
separator: \{"off"\} | "on"
state: \{"off"\} | "on"
tag: string, def. ""

A user-defined string to label the graphics object.
tooltipstring: def. ""
type (read-only): string
Class name of the graphics object. type is always "uitoggletool"
userdata: Any Octave data, def. [] ( $0 x 0$ )
User-defined data to associate with the graphics object.
visible: "off" | \{"on"\}
If visible is "off", the uitoggletool is not rendered on screen.

### 15.3.4 Searching Properties

h = findobj ()
h = findobj (prop_name, prop_value, ...)
h = findobj (prop_name, prop_value, "-logical_op", prop_name, prop_value)
h = findobj ("-property", prop_name)
h = findobj ("-regexp", prop_name, pattern)
h = findobj (hlist,...)
h = findobj (hlist, "flat", ...)
h = findobj (hlist, "-depth", d, ...)
Find graphics objects with specified properties.
When called without arguments, return all graphic objects beginning with the root object (0) and including all of its descendants.

The simplest form for narrowing the results is
findobj (prop_name, prop_value)
which returns the handles of all objects which have a property named prop_name that has the value prop_value. If multiple property/value pairs are specified then only objects meeting all of the conditions (equivalent to -and) are returned.
The search can be limited to a particular set of objects and their descendants, by passing a handle or set of handles hlist as the first argument.
The depth of the object hierarchy to search can be limited with the "-depth" argument. An example of searching through only three generations of children is:

```
findobj (hlist, "-depth", 3, prop_name, prop_value)
```

Specifying a depth $d$ of 0 limits the search to the set of objects passed in hlist. A depth of 0 is also equivalent to the "flat" argument. The default depth value is $\operatorname{Inf}$ which includes all descendants.
A specified logical operator may be used between prop_name, prop_value pairs. The supported logical operators are: "-and", "-or", "-xor", "-not". Example code to locate all figure and axes objects is
findobj ("type", "figure", "-or", "type", "axes")
Objects may also be matched by comparing a regular expression to the property values, where property values that match regexp (prop_value, pattern) are returned.
Finally, objects which have a property name can be found with the "-property" option. For example, code to locate objects with a "meshstyle" property is
findobj ("-property", "meshstyle")
Implementation Note: The search only includes objects with visible handles (HandleVisibility $=$ "on"). See [findall], page 489, to search for all objects including hidden ones.
See also: [findall], page 489, [allchild], page 431, [get], page 430, [set], page 430.

```
h = findall ()
h = findall (prop_name, prop_value, ...)
h = findall (prop_name, prop_value, "-logical_op", prop_name,
    prop_value)
h = findall ("-property", prop_name)
h = findall ("-regexp", prop_name, pattern)
h = findall (hlist,...)
h = findall (hlist, "flat", ...)
h = findall (hlist, "-depth", d,...)
```

Find graphics object, including hidden ones, with specified properties.
The return value $h$ is a list of handles to the found graphic objects.
findall performs the same search as findobj, but it includes hidden objects (HandleVisibility $=$ "off" $"$. For full documentation, see [findobj], page 488.
See also: [findobj], page 488, [allchild], page 431, [get], page 430, [set], page 430.

### 15.3.5 Managing Default Properties

Object properties have two classes of default values, factory defaults (the initial values) and user-defined defaults, which may override the factory defaults.

Although default values may be set for any object, they are set in parent objects and apply to child objects, of the specified object type. For example, setting the default color property of line objects to "green", for the root object, will result in all line objects inheriting the color "green" as the default value.

```
set (groot, "defaultlinecolor", "green");
```

sets the default line color for all objects. The rule for constructing the property name to set a default value is

```
default + object-type + property-name
```

This rule can lead to some strange looking names, for example defaultlinelinewidth" specifies the default linewidth property for line objects.

The example above used the root object so the default property value will apply to all line objects. However, default values are hierarchical, so defaults set in a figure objects override those set in the root object. Likewise, defaults set in an axes object override those set in figure or root objects. For example,

```
subplot (2, 1, 1);
set (groot, "defaultlinecolor", "red");
set (1, "defaultlinecolor", "green");
set (gca (), "defaultlinecolor", "blue");
line (1:10, rand (1, 10));
subplot (2, 1, 2);
line (1:10, rand (1, 10));
figure (2)
line (1:10, rand (1, 10));
```

produces two figures. The line in first subplot window of the first figure is blue because it inherits its color from its parent axes object. The line in the second subplot window of the first figure is green because it inherits its color from its parent figure object. The line in the second figure window is red because it inherits its color from the global root object.

To remove a user-defined default setting, set the default property to the value "remove". For example,

```
set (gca (), "defaultlinecolor", "remove");
```

removes the user-defined default line color setting from the current axes object. To quickly remove all user-defined defaults use the reset function.

By default, high level plotting functions such as plot reset and redefine axes properties independently from the defaults. An example of such property is the axes box property: it is set on by high level 2-D graphics functions regardless of the property "defaultaxesbox". Use the hold function to prevent this behavior:

```
set (groot, "defaultaxesbox", "off");
subplot (2, 1, 1);
plot (1:10)
title ("Box is on anyway")
subplot (2, 1, 2);
hold on
plot (1:10)
title ("Box is off")
```

reset (h)
Reset the properties of the graphic object $h$ to their default values.
For figures, the properties "position", "units", "windowstyle", and "paperunits" are not affected. For axes, the properties "position" and "units" are not affected.
The input $h$ may also be a vector of graphic handles in which case each individual object will be reset.
See also: [cla], page 404, [clf], page 404, [newplot], page 402.
Getting the "default" property of an object returns a list of user-defined defaults set for the object. For example,
get (gca (), "default");
returns a list of user-defined default values for the current axes object.
Factory default values are stored in the root object. The command

```
get (groot, "factory");
```

returns a list of factory defaults.

### 15.4 Advanced Plotting

### 15.4.1 Colors

Colors may be specified in three ways: 1) RGB triplets, 2) by name, or 3) by HTML notation.
RGB triplet
An RGB triplet is a 1 x 3 vector where each value is between 0 and 1 inclusive.
The first value represents the percentage of Red, the second value the percentage of Green, and the third value the percentage of Blue. For example, [1, 0, 1] represents full Red and Blue channels resulting in the color magenta.
short or long name
Eight colors can be specified directly by name or by a single character short name.

| Name | Color |
| :--- | :--- |
| ' $k$ ', "black" | blacK |
| 'r', "red" | Red |
| 'g', "green" | Green |
| 'b', "blue" | Blue |
| 'y', "yellow" | Yellow |
| 'm', "magenta" | Magenta |
| 'c', "cyan" | Cyan |
| 'w', "white" | White |

HTML notation
HTML notation is a string that begins with the character '\#' and is followed by either 3 or 6 hexadecimal digits. As with RGB triplets, each hexadecimal number represents the fraction of the Red, Green, and Blue channels present in the specified color. For example, "\#FF00FF" represents the color magenta.

### 15.4.2 Line Styles

Line styles are specified by the following properties:
linestyle
May be one of
"-" Solid line. [default]
"--" Dashed line.
":" Dotted line.
"-." A dash-dot line.
"none" No line. Points will still be marked using the current Marker Style.
linewidth
A number specifying the width of the line. The default is 1 . A value of 2 is twice as wide as the default, etc.

### 15.4.3 Marker Styles

Marker styles are specified by the following properties:
marker A character indicating a plot marker to be place at each data point, or "none", meaning no markers should be displayed.
markeredgecolor
The color of the edge around the marker, or "auto", meaning that the edge color is the same as the face color. See Section 15.4.1 [Colors], page 491.
markerfacecolor
The color of the marker, or "none" to indicate that the marker should not be filled. See Section 15.4.1 [Colors], page 491.
markersize
A number specifying the size of the marker. The default is 1 . A value of 2 is twice as large as the default, etc.

The colstyle function will parse a plot-style specification and will return the color, line, and marker values that would result.
[style, color, marker, msg] = colstyle (style)
Parse the line specification style and return the line style, color, and markers given.
In the case of an error, the string msg will return the text of the error.

### 15.4.4 Callbacks

Callback functions can be associated with graphics objects and triggered after certain events occur. The basic structure of all callback function is

```
function mycallback (hsrc, evt)
endfunction
```

where hsrc is a handle to the source of the callback, and evt gives some event specific data.

The function can be provided as a function handle to a plain Octave function, as an anonymous function, or as a string representing an Octave command. The latter syntax is not recommended since syntax errors will only occur when the string is evaluated. See Section 11.12 [Function Handles section], page 233.

This can then be associated with an object either at the object's creation, or later with the set function. For example,

```
plot (x, "DeleteFcn", @(h, e) disp ("Window Deleted"))
```

where at the moment that the plot is deleted, the message "Window Deleted" will be displayed.

Additional user arguments can be passed to callback functions, and will be passed after the two default arguments. For example:

```
plot (x, "DeleteFcn", {@mycallback, "1"})
function mycallback (h, evt, arg1)
    fprintf ("Closing plot %d\n", arg1);
endfunction
```

Caution: The second argument in callback functions-evt-is only partially implemented in the Qt graphics toolkit:

- Mouse click events: evt is a class double value: 1 for left, 2 for middle, and 3 for right click.
- Key press events: evt is a structure with fields Key (string), Character (string), and Modifier (cell array of strings).
- Other events: evt is a class double empty matrix.

The basic callback functions that are available for all graphics objects are

- CreateFcn: called at the moment of the objects creation. It is not called if the object is altered in any way, and so it only makes sense to define this callback in the function call that defines the object. Callbacks that are added to CreateFcn later with the set function will never be executed.
- DeleteFcn: called at the moment an object is deleted.
- ButtonDownFcn: called if a mouse button is pressed while the pointer is over this object. Note, that the gnuplot interface does not implement this callback.

By default callback functions are queued (they are executed one after the other in the event queue) unless the drawnow, figure, waitfor, getframe, or pause functions are used. If an executing callback invokes one of those functions, it causes Octave to flush the event queue, which results in the executing callback being interrupted.

It is possible to specify that an object's callbacks should not be interrupted by setting the object's interruptible property to "off". In this case, Octave decides what to do based on the busyaction property of the interrupting callback object:
queue (the default)
The interrupting callback is executed after the executing callback has returned.
cancel The interrupting callback is discarded.

The interruptible property has no effect when the interrupting callback is a deletefcn, or a figure resizefcn or closerequestfcn. Those callbacks always interrupt the executing callback.

The handle to the object that holds the callback being executed can be obtained with the gcbo function. The handle to the ancestor figure of this object may be obtained using the gcbf function.

```
h = gcbo ()
[h, fig] = gcbo ()
```

Return a handle to the object whose callback is currently executing.
If no callback is executing, this function returns the empty matrix. This handle is obtained from the root object property "CallbackObject".
When called with a second output argument, return the handle of the figure containing the object whose callback is currently executing. If no callback is executing the second output is also set to the empty matrix.
See also: [gcbf], page 494, [gco], page 429, [gca], page 429, [gcf], page 428, [get], page 430, [set], page 430.
fig $=\operatorname{gcbf}()$
Return a handle to the figure containing the object whose callback is currently executing.
If no callback is executing, this function returns the empty matrix. The handle returned by this function is the same as the second output argument of gcbo.
See also: [gcbo], page 494, [gcf], page 428, [gco], page 429, [gca], page 429, [get], page 430, [set], page 430.

Callbacks can equally be added to properties with the addlistener function described below.

### 15.4.5 Application-defined Data

Octave has a provision for attaching application-defined data to a graphics handle. The data can be anything which is meaningful to the application, and will be completely ignored by Octave.

```
setappdata (h, name, value)
setappdata (h, name1, value1, name2, value3,...)
setappdata (h, {name1, name2,...}, {value1, value2, ...})
```

Set the application data name to value for the graphics object with handle $h$.
$h$ may also be a vector of graphics handles. If the application data with the specified name does not exist, it is created.
Multiple name/value pairs can be specified. Alternatively, a cell array of names and a corresponding cell array of values can be specified. For details on obtaining a list of valid application data properties, see [getappdata], page 495.
See also: [getappdata], page 495, [isappdata], page 495, [rmappdata], page 495, [guidata], page 933, [get], page 430, [set], page 430, [getpref], page 937, [setpref], page 937.

```
value = getappdata (h, name)
```

appdata $=$ getappdata $(h)$
Return the value of the application data name for the graphics object with handle $h$. $h$ may also be a vector of graphics handles. If no second argument name is given then getappdata returns a structure, appdata, whose fields correspond to the appdata properties.

See also: [setappdata], page 494, [isappdata], page 495, [rmappdata], page 495, [guidata], page 933, [get], page 430, [set], page 430, [getpref], page 937, [setpref], page 937.
rmappdata (h, name)
rmappdata (h, name1, name2, ...)
Delete the application data name from the graphics object with handle $h$.
$h$ may also be a vector of graphics handles. Multiple application data names may be supplied to delete several properties at once.

See also: [setappdata], page 494, [getappdata], page 495, [isappdata], page 495.

```
valid = isappdata (h, name)
```

Return true if the named application data, name, exists for the graphics object with handle $h$.
$h$ may also be a vector of graphics handles.
See also: [getappdata], page 495, [setappdata], page 494, [rmappdata], page 495, [guidata], page 933, [get], page 430, [set], page 430, [getpref], page 937, [setpref], page 937.

### 15.4.6 Object Groups

A number of Octave high level plot functions return groups of other graphics objects or they return graphics objects that have their properties linked in such a way that changes to one of the properties results in changes in the others. A graphic object that groups other objects is an hggroup
hggroup ()
hggroup (hax)
hggroup (..., property, value, ...)
$h=$ hggroup (...)
Create handle graphics group object with axes parent hax.
If no parent is specified, the group is created in the current axes.
Multiple property/value pairs may be specified for the hggroup, but they must appear in pairs. The full list of properties is documented at Section 15.3.3.3 [Axes Properties], page 441.

The optional return value $h$ is a graphics handle to the created hggroup object.
Programming Note: An hggroup is a way to group base graphics objects such as line objects or patch objects into a single unit which can react appropriately. For example, the individual lines of a contour plot are collected into a single hggroup so that they
can be made visible/invisible with a single command, set (hg_handle, "visible", "off").

See also: [addproperty], page 496, [addlistener], page 497.
For example a simple use of a hggroup might be

```
x = 0:0.1:10;
hg = hggroup ();
plot (x, sin (x), "color", [1, 0, 0], "parent", hg);
hold on
plot (x, cos (x), "color", [0, 1, 0], "parent", hg);
set (hg, "visible", "off");
```

which groups the two plots into a single object and controls their visibility directly. The default properties of an hggroup are the same as the set of common properties for the other graphics objects. Additional properties can be added with the addproperty function.
addproperty (name, h, type)
addproperty (name, h, type, arg, ...)
Create a new property named name in graphics object $h$.
type determines the type of the property to create. args usually contains the default value of the property, but additional arguments might be given, depending on the type of the property.
The supported property types are:
string A string property. arg contains the default string value.
any An un-typed property. This kind of property can hold any octave value. args contains the default value.
radio A string property with a limited set of accepted values. The first argument must be a string with all accepted values separated by a vertical bar ('I'). The default value can be marked by enclosing it with a ' $\{$ ' ' $\}$ ' pair. The default value may also be given as an optional second string argument.
boolean A boolean property. This property type is equivalent to a radio property with "onloff" as accepted values. arg contains the default property value.
double A scalar double property. arg contains the default value.
handle A handle property. This kind of property holds the handle of a graphics object. arg contains the default handle value. When no default value is given, the property is initialized to the empty matrix.
data A data (matrix) property. arg contains the default data value. When no default value is given, the data is initialized to the empty matrix.
color A color property. arg contains the default color value. When no default color is given, the property is set to black. An optional second string argument may be given to specify an additional set of accepted string values (like a radio property).
type may also be the concatenation of a core object type and a valid property name for that object type. The property created then has the same characteristics as the referenced property (type, possible values, hidden state. . .). This allows one to clone an existing property into the graphics object $h$.

Examples:

```
addproperty ("my_property", gcf, "string", "a string value");
addproperty ("my_radio", gcf, "radio", "val_1|val_2|{val_3}");
addproperty ("my_style", gcf, "linelinestyle", "--");
```

See also: [addlistener], page 497, [hggroup], page 495.

Once a property in added to an hggroup, it is not linked to any other property of either the children of the group, or any other graphics object. Add so to control the way in which this newly added property is used, the addlistener function is used to define a callback function that is executed when the property is altered.

## addlistener (h, prop, fcn)

Register fcn as listener for the property prop of the graphics object $h$.
Property listeners are executed (in order of registration) when the property is set. The new value is already available when the listeners are executed.
prop must be a string naming a valid property in $h$.
fcn can be a function handle, a string or a cell array whose first element is a function handle. If fcn is a function handle, the corresponding function should accept at least 2 arguments, that will be set to the object handle and the empty matrix respectively. If $f c n$ is a string, it must be any valid octave expression. If $f c n$ is a cell array, the first element must be a function handle with the same signature as described above. The next elements of the cell array are passed as additional arguments to the function.

Example:

```
function my_listener (h, dummy, p1)
    fprintf ("my_listener called with p1=%s\n", p1);
endfunction
addlistener (gcf, "position", {@my_listener, "my string"})
```

See also: [dellistener], page 497, [addproperty], page 496, [hggroup], page 495.
dellistener (h, prop, fcn)
Remove the registration of $f c n$ as a listener for the property prop of the graphics object $h$.

The function fcn must be the same variable (not just the same value), as was passed to the original call to addlistener.

If fcn is not defined then all listener functions of prop are removed.
Example:

```
function my_listener (h, dummy, p1)
    fprintf ("my_listener called with p1=%s\n", p1);
endfunction
c = {@my_listener, "my string"};
addlistener (gcf, "position", c);
dellistener (gcf, "position", c);
```

See also: [addlistener], page 497.
An example of the use of these two functions might be

```
x = 0:0.1:10;
hg = hggroup ();
h = plot (x, sin (x), "color", [1, 0, 0], "parent", hg);
addproperty ("linestyle", hg, "linelinestyle", get (h, "linestyle"));
addlistener (hg, "linestyle", @update_props);
hold on
plot (x, cos (x), "color", [0, 1, 0], "parent", hg);
function update_props (h, d)
    set (get (h, "children"), "linestyle", get (h, "linestyle"));
endfunction
```

that adds a linestyle property to the hggroup and propagating any changes its value to the children of the group. The linkprop function can be used to simplify the above to be

```
x = 0:0.1:10;
hg = hggroup ();
h1 = plot (x, sin (x), "color", [1, 0, 0], "parent", hg);
addproperty ("linestyle", hg, "linelinestyle", get (h, "linestyle"));
hold on
h2 = plot (x, cos (x), "color", [0, 1, 0], "parent", hg);
hlink = linkprop ([hg, h1, h2], "color");
hlink = linkprop (h, "prop")
hlink = linkprop (h, {"prop1", "prop2", ...})
```

Link graphic object properties, such that a change in one is propagated to the others.
The input $h$ is a vector of graphic handles to link.
prop may be a string when linking a single property, or a cell array of strings for multiple properties. During the linking process all properties in prop will initially be set to the values that exist on the first object in the list $h$.
The function returns hlink which is a special object describing the link. As long as the reference hlink exists, the link between graphic objects will be active. This means that hlink must be preserved in a workspace variable, a global variable, or otherwise stored using a function such as setappdata or guidata. To unlink properties, execute clear hlink.
An example of the use of linkprop is

```
x = 0:0.1:10;
subplot (1,2,1);
h1 = plot (x, sin (x));
subplot (1,2,2);
h2 = plot (x, cos (x));
hlink = linkprop ([h1, h2], {"color","linestyle"});
set (h1, "color", "green");
set (h2, "linestyle", "--");
```

See also: [linkaxes], page 499, [addlistener], page 497.
linkaxes (hax)
linkaxes (hax, optstr)
Link the axis limits of 2-D plots such that a change in one is propagated to the others.
The axes handles to be linked are passed as the first argument hax.
The optional second argument is a string which defines which axis limits will be linked. The possible values for optstr are:
"x" Link x-axes
"y" Link y-axes
"xy" (default)
Link both axes
"off" Turn off linking
If unspecified the default is to link both X and Y axes.
When linking, the limits from the first axes in hax are applied to the other axes in the list. Subsequent changes to any one of the axes will be propagated to the others.
See also: [linkprop], page 498, [addproperty], page 496.
These capabilities are used in a number of basic graphics objects. The hggroup objects created by the functions of Octave contain one or more graphics object and are used to:

- group together multiple graphics objects,
- create linked properties between different graphics objects, and
- to hide the nominal user data, from the actual data of the objects.

For example the stem function creates a stem series where each hggroup of the stem series contains two line objects representing the body and head of the stem. The ydata property of the hggroup of the stem series represents the head of the stem, whereas the body of the stem is between the baseline and this value. For example

```
h = stem (1:4)
get (h, "xdata")
# [ [\begin{array}{llll}{1}&{2}&{3}&{4]'}\end{array}]
get (get (h, "children")(1), "xdata")
# [\begin{array}{ccccccll}{1}&{1 NaN 2 NaN 3 3 NaN 4 4 NaN]'}\end{array}]
```

shows the difference between the xdata of the hggroup of a stem series object and the underlying line.

The basic properties of such group objects is that they consist of one or more linked hggroup, and that changes in certain properties of these groups are propagated to other members of the group. Whereas, certain properties of the members of the group only apply to the current member.

In addition the members of the group can also be linked to other graphics objects through callback functions. For example the baseline of the bar or stem functions is a line object, whose length and position are automatically adjusted, based on changes to the corresponding hggroup elements.

### 15.4.6.1 Data Sources in Object Groups

All of the group objects contain data source parameters. There are string parameters that contain an expression that is evaluated to update the relevant data property of the group when the refreshdata function is called.

```
refreshdata ()
refreshdata (h)
refreshdata (h, workspace)
```

Evaluate any 'datasource' properties of the current figure and update the plot if the corresponding data has changed.
If the first argument $h$ is a list of graphic handles, then operate on these objects rather than the current figure returned by gcf.
The optional second argument workspace can take the following values:
"base" Evaluate the datasource properties in the base workspace. (default).
"caller" Evaluate the datasource properties in the workspace of the function that called refreshdata.

An example of the use of refreshdata is:

```
\(\mathrm{x}=0: 0.1: 10 ;\)
\(y=\sin (x) ;\)
plot (x, y, "ydatasource", "y");
for \(i=1\) : 100
    pause (0.1);
    \(\mathrm{y}=\sin (\mathrm{x}+0.1 * \mathrm{i})\);
    refreshdata ();
endfor
```


### 15.4.6.2 Area Series

Area series objects are created by the area function. Each of the hggroup elements contains a single patch object. The properties of the area series are
basevalue
The value where the base of the area plot is drawn.
linewidth
linestyle
The line width and style of the edge of the patch objects making up the areas. See Section 15.4.2 [Line Styles], page 492.

```
edgecolor
facecolor
The line and fill color of the patch objects making up the areas. See Section 15.4.1 [Colors], page 491.
xdata
ydata The x and y coordinates of the original columns of the data passed to area
    prior to the cumulative summation used in the area function.
xdatasource
ydatasource
Data source variables.
```


### 15.4.6.3 Bar Series

Bar series objects are created by the bar or barh functions. Each hggroup element contains a single patch object. The properties of the bar series are

```
showbaseline
baseline
basevalue
```

The property showbaseline flags whether the baseline of the bar series is displayed (default is "on"). The handle of the graphics object representing the baseline is given by the baseline property and the y -value of the baseline by the basevalue property.
Changes to any of these properties are propagated to the other members of the bar series and to the baseline itself. Equally, changes in the properties of the base line itself are propagated to the members of the corresponding bar series.

```
barwidth
barlayout
horizontal
```

The property barwidth is the width of the bar corresponding to the width variable passed to bar or barh. Whether the bar series is "grouped" or "stacked" is determined by the barlayout property and whether the bars are horizontal or vertical by the horizontal property.
Changes to any of these property are propagated to the other members of the bar series.
linewidth
linestyle
The line width and style of the edge of the patch objects making up the bars. See Section 15.4.2 [Line Styles], page 492.
edgecolor
facecolor
The line and fill color of the patch objects making up the bars. See Section 15.4.1 [Colors], page 491.
xdata The nominal x positions of the bars. Changes in this property and propagated to the other members of the bar series.
ydata The y value of the bars in the hggroup.
xdatasource
ydatasource
Data source variables.

### 15.4.6.4 Contour Groups

Contour group objects are created by the contour, contourf, and contour3 functions. They are also one of the handles returned by the surfc and meshc functions. The properties of the contour group are
contourmatrix
A read only property that contains the data return by contourc used to create the contours of the plot.
fill A radio property that can have the values "on" or "off" that flags whether the contours to plot are to be filled.
zlevelmode
zlevel The radio property zlevelmode can have the values "none", "auto", or "manual". When its value is "none" there is no z component to the plotted contours. When its value is "auto" the $z$ value of the plotted contours is at the same value as the contour itself. If the value is "manual", then the $z$ value at which to plot the contour is determined by the zlevel property.
levellistmode
levellist
levelstepmode
levelstep
If levellistmode is "manual", then the levels at which to plot the contours is determined by levellist. If levellistmode is set to "auto", then the distance between contours is determined by levelstep. If both levellistmode and levelstepmode are set to "auto", then there are assumed to be 10 equal spaced contours.

```
textlistmode
```

textlist
textstepmode
textstep If textlistmode is "manual", then the labeled contours is determined by textlist. If textlistmode is set to "auto", then the distance between labeled contours is determined by textstep. If both textlistmode and textstepmode are set to "auto", then there are assumed to be 10 equal spaced labeled contours.
showtext Flag whether the contour labels are shown or not.
labelspacing
The distance between labels on a single contour in points.

```
linewidth
linestyle
linecolor
```

The properties of the contour lines. The properties linewidth and linestyle are similar to the corresponding properties for lines. The property linecolor is a color property (see Section 15.4.1 [Colors], page 491), that can also have the values of "none" or "auto". If linecolor is "none", then no contour line is drawn. If linecolor is "auto" then the line color is determined by the colormap.

```
xdata
ydata
zdata The original x, y, and z data of the contour lines.
xdatasource
ydatasource
zdatasource
```

Data source variables.

### 15.4.6.5 Error Bar Series

Error bar series are created by the errorbar function. Each hggroup element contains two line objects representing the data and the errorbars separately. The properties of the error bar series are
color The RGB color or color name of the line objects of the error bars. See Section 15.4.1 [Colors], page 491.

```
linewidth
linestyle
```

The line width and style of the line objects of the error bars. See Section 15.4.2 [Line Styles], page 492.

```
marker
markeredgecolor
markerfacecolor
markersize
```

The line and fill color of the markers on the error bars. See Section 15.4.1 [Colors], page 491.

```
xdata
ydata
ldata
udata
xldata
xudata The original x, y, l, u, xl, xu data of the error bars.
```

```
xdatasource
ydatasource
ldatasource
udatasource
xldatasource
xudatasource
Data source variables.
```


### 15.4.6.6 Line Series

Line series objects are created by the plot and plot3 functions and are of the type line. The properties of the line series with the ability to add data sources.
color The RGB color or color name of the line objects. See Section 15.4.1 [Colors], page 491.

```
linewidth
```

linestyle

The line width and style of the line objects. See Section 15.4.2 [Line Styles], page 492.
marker
markeredgecolor
markerfacecolor
markersize
The line and fill color of the markers. See Section 15.4.1 [Colors], page 491.
xdata
ydata
zdata The original $\mathrm{x}, \mathrm{y}$ and z data.
xdatasource
ydatasource
zdatasource
Data source variables.

### 15.4.6.7 Quiver Group

Quiver series objects are created by the quiver or quiver3 functions. Each hggroup element of the series contains three line objects as children representing the body and head of the arrow, together with a marker as the point of origin of the arrows. The properties of the quiver series are
autoscale
autoscalefactor
Flag whether the length of the arrows is scaled or defined directly from the $u, v$ and $w$ data. If the arrow length is flagged as being scaled by the autoscale property, then the length of the autoscaled arrow is controlled by the autoscalefactor.
maxheadsize
This property controls the size of the head of the arrows in the quiver series. The default value is 0.2 .

```
showarrowhead
    Flag whether the arrow heads are displayed in the quiver plot.
color The RGB color or color name of the line objects of the quiver. See Section 15.4.1
    [Colors], page 491.
linewidth
linestyle
The line width and style of the line objects of the quiver. See Section 15.4.2 [Line Styles], page 492.
marker
markerfacecolor
markersize
The line and fill color of the marker objects at the original of the arrows. See Section 15.4.1 [Colors], page 491.
```

```
xdata
```

xdata
ydata
zdata The origins of the values of the vector field.
udata
vdata
wdata The values of the vector field to plot.
xdatasource
ydatasource
zdatasource
udatasource
vdatasource
wdatasource
Data source variables.

```

\subsection*{15.4.6.8 Stair Group}

Stair series objects are created by the stair function. Each hggroup element of the series contains a single line object as a child representing the stair. The properties of the stair series are
color \(\quad\) The RGB color or color name of the line objects of the stairs. See Section 15.4.1 [Colors], page 491.
linewidth
linestyle
The line width and style of the line objects of the stairs. See Section 15.4.2 [Line Styles], page 492.
marker
markeredgecolor
markerfacecolor
markersize
The line and fill color of the markers on the stairs. See Section 15.4.1 [Colors], page 491.
xdata
ydata The original x and y data of the stairs.
```

xdatasource

```
ydatasource
Data source variables.

\subsection*{15.4.6.9 Stem Series}

Stem series objects are created by the stem or stem3 functions. Each hggroup element contains a single line object as a child representing the stems. The properties of the stem series are
showbaseline
baseline
basevalue
The property showbaseline flags whether the baseline of the stem series is displayed (default is "on"). The handle of the graphics object representing the baseline is given by the baseline property and the \(y\)-value (or \(z\)-value for stem3) of the baseline by the basevalue property.
Changes to any of these property are propagated to the other members of the stem series and to the baseline itself. Equally changes in the properties of the base line itself are propagated to the members of the corresponding stem series.
color \(\quad\) The RGB color or color name of the line objects of the stems. See Section 15.4.1 [Colors], page 491.
linewidth
linestyle
The line width and style of the line objects of the stems. See Section 15.4.2 [Line Styles], page 492.
```

marker
markeredgecolor
markerfacecolor
markersize

```

The line and fill color of the markers on the stems. See Section 15.4.1 [Colors], page 491.
```

xdata

```
ydata
zdata The original \(x, y\) and \(z\) data of the stems.
xdatasource
ydatasource
zdatasource

Data source variables.

\subsection*{15.4.6.10 Surface Group}

Surface group objects are created by the surf or mesh functions, but are equally one of the handles returned by the surfc or meshc functions. The surface group is of the type surface.

The properties of the surface group are
```

edgecolor
facecolor

```

The RGB color or color name of the edges or faces of the surface. See Section 15.4.1 [Colors], page 491.
```

linewidth
linestyle

```

The line width and style of the lines on the surface. See Section 15.4.2 [Line Styles], page 492.
```

marker
markeredgecolor
markerfacecolor
markersize

```

The line and fill color of the markers on the surface. See Section 15.4.1 [Colors], page 491.
```

xdata
ydata
zdata
xdatasource
ydatasource
zdatasource
cdatasource

```
cdata The original \(\mathrm{x}, \mathrm{y}, \mathrm{z}\) and c data.

Data source variables.

\subsection*{15.4.7 Transform Groups}
```

h = hgtransform ()
h = hgtransform (property, value,...)
h = hgtransform (hax,...)

```
    Create a graphics transform object.

FIXME: Need to write documentation. FIXME: Add <makehgtform> to seealso list when it is implemented.

See also: [hggroup], page 495.

\subsection*{15.4.8 Graphics Toolkits}
```

name = graphics_toolkit ()
name = graphics_toolkit (hlist)
graphics_toolkit (name)
graphics_toolkit (hlist, name)

```

Query or set the default graphics toolkit which is assigned to new figures.
With no inputs, return the current default graphics toolkit. If the input is a list of figure graphic handles, hlist, then return the name of the graphics toolkit in use for each figure.

When called with a single input name set the default graphics toolkit to name. If the toolkit is not already loaded, it is initialized by calling the function __init_name__. If the first input is a list of figure handles, hlist, then the graphics toolkit is set to name for these figures only.

See also: [available_graphics_toolkits], page 508.
available_graphics_toolkits ()
Return a cell array of registered graphics toolkits.
See also: [graphics_toolkit], page 507, [register_graphics_toolkit], page 508.
loaded_graphics_toolkits ()
Return a cell array of the currently loaded graphics toolkits.
See also: [available_graphics_toolkits], page 508.
register_graphics_toolkit (toolkit)
List toolkit as an available graphics toolkit.
See also: [available_graphics_toolkits], page 508.

\subsection*{15.4.8.1 Customizing Toolkit Behavior}

The specific behavior of the backend toolkit may be modified using the following utility functions. Note: Not all functions apply to every graphics toolkit.
```

[prog, args] = gnuplot_binary ()
[old_prog, old_args] = gnuplot_binary (new_prog)
[old_prog, old_args] = gnuplot_binary (new_prog, arg1,...)

```

Query or set the name of the program invoked by the plot command when the graphics toolkit is set to "gnuplot".
Additional arguments to pass to the external plotting program may also be given. The default value is "gnuplot" with no additional arguments. See Appendix E [Installation], page 1085.
See also: [graphics_toolkit], page 507.
In addition, the gnuplot program usually provides a number of different interfaces, known as terminals. Octave normally chooses a default terminal, but you can override this with the environment variable GNUTERM. This variable may be set in the shell before starting Octave or from within Octave before plotting for the first time. For example:
```

setenv ("GNUTERM", "wxt")
graphics_toolkit ("gnuplot")
plot (1:10)

```

\subsection*{15.4.8.2 Hardware vs. Software Rendering}

On Windows platforms, Octave uses software rendering for the OpenGL graphics toolkits ("qt" and "fltk") by default. This is done to avoid rendering and printing issues due to imperfect OpenGL driver implementations for diverse graphic cards from different vendors. As a down-side, software rendering might be considerably slower than hardware accelerated
rendering. To permanently switch back to hardware accelerated rendering with your graphic card drivers, rename the following file while Octave is closed:
octave-home \(\backslash\) bin \(\backslash o p e n g 132\).dll
where octave-home is the directory in which Octave is installed (the default is C: \Octave\Octave-version).

\subsection*{15.4.8.3 Precision issues}

The OpenGL graphics toolkits ("qt" and "fltk") use single precision for rendering. This limitation in particular applies to plots of time series against serial dates as used by the datenum, datestr, datestruct, and datetick functions.

Serial dates encode timestamps as days elapsed since the year zero with hours, minutes, seconds as the fractional part. On December 31st 1999, the serial date representation was 730485. A double precision variable with this integer part allows for a resolution in its fractional part of \(1.2 \mathrm{e}-10\), representing about 5 microseconds. But with single precision, the resolution is reduced to about 0.06 , representing 45 minutes. Any attempt to plot timestamped data with finer granularity will result in a distorted graph.

As a workaround, it is possible to use the "gnuplot" graphics toolkit or subtract 2000 years-i.e., datenum ( \(2000,0,0\) ) or 730485 -from the time values. Due to the fact that the calendar structure repeats every 2000 years, the relation between year, month, day of month and day of week will stay unchanged and the ticks and ticklabels produced by the datetick function will still be correct. Only years will lack the millennium digit. Thus, "2020" will be printed as "20". For example:
```


# timestamps of 24 hours in one minute steps

t = datenum (2020, 1, 1):(1/1440):datenum (2020, 1, 2);

# some example time series data

x = -cos (2*pi*t) + rand (size (t)) / 10;
subplot (1, 2, 1);
plot (t, x);
datetick ("x");
xlabel ("serial date");
title ("problem");
subplot (1, 2, 2);
plot (t - 730485, x);
datetick ("x");
xlabel ("2000 years off");
title ("workaround");

```

The result of which can be seen in Figure 15.8.


Figure 15.8: Single precision issues with OpenGL graphics toolkits
Similarly, other data can be translated or re-scaled to work around this issue.

\section*{16 Matrix Manipulation}

There are a number of functions available for checking to see if the elements of a matrix meet some condition, and for rearranging the elements of a matrix. For example, Octave can easily tell you if all the elements of a matrix are finite, or are less than some specified value. Octave can also rotate the elements, extract the upper- or lower-triangular parts, or sort the columns of a matrix.

\subsection*{16.1 Finding Elements and Checking Conditions}

The functions any and all are useful for determining whether any or all of the elements of a matrix satisfy some condition. The find function is also useful in determining which elements of a matrix meet a specified condition.
any ( \(x\) )
any (x, dim)
For a vector argument, return true (logical 1) if any element of the vector is nonzero.
For a matrix argument, return a row vector of logical ones and zeros with each element indicating whether any of the elements of the corresponding column of the matrix are nonzero. For example:
```

any (eye (2, 4))
\# [ 1, 1, 0, 0 ]

```

If the optional argument dim is supplied, work along dimension dim. For example:
```

any (eye (2, 4), 2)
\# [ 1; 1]

```

See also: [all], page 511.
all (x)
all (x, dim)
For a vector argument, return true (logical 1) if all elements of the vector are nonzero.
For a matrix argument, return a row vector of logical ones and zeros with each element indicating whether all of the elements of the corresponding column of the matrix are nonzero. For example:
\[
\begin{aligned}
& \text { all }([2,3 ; 1,0]) \\
& \quad \Rightarrow[1,0]
\end{aligned}
\]

If the optional argument dim is supplied, work along dimension dim.
See also: [any], page 511.
Since the comparison operators (see Section 8.4 [Comparison Ops], page 164) return matrices of ones and zeros, it is easy to test a matrix for many things, not just whether the elements are nonzero. For example,
```

all (all (rand (5) < 0.9))
=> 0

```
tests a random 5 by 5 matrix to see if all of its elements are less than 0.9 .
Note that in conditional contexts (like the test clause of if and while statements) Octave treats the test as if you had typed all (all (condition)).
\(z=\operatorname{xor}(x, y)\)
\(z=\) xor (x1, x2, ..)
Return the exclusive or of \(x\) and \(y\).
For boolean expressions \(x\) and \(y\), xor \((x, y)\) is true if and only if one of \(x\) or \(y\) is true. Otherwise, if \(x\) and \(y\) are both true or both false, xor returns false.
The truth table for the xor operation is
\[
\begin{array}{ccc}
x & y & z \\
- & - & - \\
0 & 0 & 0 \\
1 & 0 & 1 \\
0 & 1 & 1 \\
1 & 1 & 0
\end{array}
\]

If more than two arguments are given the xor operation is applied cumulatively from left to right:
(...((x1 XOR x2) XOR x3) XOR ...)

See also: [and], page 166, [or], page 167, [not], page 167 .
```

diff (x)

```
diff ( \(x, k\) )
diff (x, k, dim)

If x is a vector of length \(n\), diff \((\mathrm{x})\) is the vector of first differences \(x_{2}-x_{1}, \ldots, x_{n}-\) \(x_{n-1}\).
If \(x\) is a matrix, diff ( \(x\) ) is the matrix of column differences along the first nonsingleton dimension.
The second argument is optional. If supplied, \(\operatorname{diff}(x, k)\), where \(k\) is a non-negative integer, returns the \(k\)-th differences. It is possible that \(k\) is larger than the first nonsingleton dimension of the matrix. In this case, diff continues to take the differences along the next non-singleton dimension.
The dimension along which to take the difference can be explicitly stated with the optional variable dim. In this case the \(k\)-th order differences are calculated along this dimension. In the case where \(k\) exceeds size ( \(x\), dim) an empty matrix is returned.
See also: [sort], page 521, [merge], page 168.
isinf (x)
Return a logical array which is true where the elements of \(x\) are infinite and false where they are not.
For example:
```

isinf ([13, Inf, NA, NaN])
[ [0, 1, 0, 0 ]

```

See also: [isfinite], page 513, [isnan], page 512, [isna], page 45.
isnan (x)
Return a logical array which is true where the elements of \(x\) are NaN values and false where they are not.

NA values are also considered NaN values. For example:
```

isnan ([13, Inf, NA, NaN])
[ [0, 0, 1, 1]

```

See also: [isna], page 45, [isinf], page 512, [isfinite], page 513.
isfinite (x)
Return a logical array which is true where the elements of \(x\) are finite values and false where they are not.
For example:
```

isfinite ([13, Inf, NA, NaN])
| [1, 0, 0, 0 ]

```

See also: [isinf], page 512, [isnan], page 512, [isna], page 45.
[err, yi, ...] = common_size (xi,...)
Determine if all input arguments are either scalar or of common size.
If true, err is zero, and yi is a matrix of the common size with all entries equal to xi if this is a scalar or xi otherwise. If the inputs cannot be brought to a common size, err is 1 , and yi is xi. For example:
```

[err, a, b] = common_size ([1 2; 3 4], 5)
merr = 0
a a = [1, 2; 3, 4 ]
=>b=[5,5;5,5 ]

```

This is useful for implementing functions where arguments can either be scalars or of common size.

See also: [size], page 47, [size_equal], page 49, [numel], page 46, [ndims], page 46.
```

idx = find (x)
idx = find (x, n)
idx = find (x, n, direction)
[i, j] = find (...)
[i, j, v] = find (...)

```

Return a vector of indices of nonzero elements of a matrix, as a row if \(x\) is a row vector or as a column otherwise.
To obtain a single index for each matrix element, Octave pretends that the columns of a matrix form one long vector (like Fortran arrays are stored). For example:
```

find (eye (2))
$\Rightarrow[1 ; 4]$

```

If two inputs are given, \(n\) indicates the maximum number of elements to find from the beginning of the matrix or vector.
If three inputs are given, direction should be one of "first" or "last", requesting only the first or last \(n\) indices, respectively. However, the indices are always returned in ascending order.

If two outputs are requested, find returns the row and column indices of nonzero elements of a matrix. For example:
\[
\begin{aligned}
{[i,} & j] \\
\Rightarrow & \text { find }(2 * \text { eye }(2)) \\
\Rightarrow & =[1 ; 2] \\
\Rightarrow j & =[1 ; 2]
\end{aligned}
\]

If three outputs are requested, find also returns a vector containing the nonzero values. For example:
```

[i, j, v] = find (3 * eye (2))
\# i = [ 1; 2 ]
\# j = [ 1; 2 ]
| v = [ 3; 3 ]

```

If x is a multi-dimensional array of size \(\mathrm{m} \times \mathrm{n} \times \mathrm{p} \times \ldots, j\) contains the column locations as if \(x\) was flattened into a two-dimensional matrix of size \(m x(n+p+\) ...).
Note that this function is particularly useful for sparse matrices, as it extracts the nonzero elements as vectors, which can then be used to create the original matrix.
For example:
```

sz = size (a);
[i, j, v] = find (a);
b = sparse (i, j, v, sz(1), sz(2));

```

See also: [nonzeros], page 666.
```

idx = lookup (table, y)
idx = lookup (table, y, opt)

```

Lookup values in a sorted table.
This function is usually used as a prelude to interpolation.
If table is increasing, of length N and \(i d x=\) lookup (table, y ), then table(idx(i)) \(<=y(i)<t a b l e(i d x(i+1))\) for all \(y(i)\) within the table. If \(y(i)<t a b l e(1)\) then \(i d x(i)\) is 0 . If \(y(i)>=t a b l e(e n d)\) or isnan ( \(y(i))\) then \(i d x(i)\) is \(N\).
If the table is decreasing, then the tests are reversed. For non-strictly monotonic tables, empty intervals are always skipped. The result is undefined if table is not monotonic, or if table contains a NaN .
The complexity of the lookup is \(\mathrm{O}\left(\mathrm{M}^{*} \log (\mathrm{~N})\right)\) where M is the size of \(y\). In the special case when \(y\) is also sorted, the complexity is \(\mathrm{O}\left(\min \left(\mathrm{M}^{*} \log (\mathrm{~N}), \mathrm{M}+\mathrm{N}\right)\right)\).
table and \(y\) can also be cell arrays of strings (or \(y\) can be a single string). In this case, string lookup is performed using lexicographical comparison.
If opts is specified, it must be a string with letters indicating additional options.
\(m \quad\) Match. table(idx(i)) == \(y\) (i) if \(y(i)\) occurs in table; otherwise, \(i d x(i)\) is zero.
b
Boolean. \(\mathrm{idx}(\mathrm{i})\) is a logical 1 or 0 , indicating whether \(\mathrm{y}(\mathrm{i})\) is contained in table or not.

1
Left. For numeric lookups the leftmost subinterval shall be extended to minus infinity (i.e., all indices at least 1).
r
Right. For numeric lookups the rightmost subinterval shall be extended to infinity (i.e., all indices at most \(\mathrm{N}-1\) ).

Note: If table is not sorted the results from lookup will be unpredictable.
If you wish to check if a variable exists at all, instead of properties its elements may have, consult Section 7.3 [Status of Variables], page 142.

\subsection*{16.2 Rearranging Matrices}
fliplr (x)
Flip array left to right.
Return a copy of \(x\) with the order of the columns reversed. In other words, \(x\) is flipped left-to-right about a vertical axis. For example:
```

fliplr ([1, 2; 3, 4])
\# 2 1
4 3

```

See also: [flipud], page 515, [flip], page 515, [rot90], page 516, [rotdim], page 516.
flipud (x)
Flip array upside down.
Return a copy of \(x\) with the order of the rows reversed. In other words, \(x\) is flipped upside-down about a horizontal axis. For example:
```

flipud ([1, 2; 3, 4])
=> 3}\begin{array}{l}{3}<br>{1}

```

See also: [fliplr], page 515, [flip], page 515, [rot90], page 516, [rotdim], page 516.
flip (x)
flip (x, dim)
Return a copy of array x flipped across dimension dim.
If dim is unspecified it defaults to the first non-singleton dimension.
Examples:
```


## row vector

flip ([llllll
\#

## column vector

flip ([1; 2; 3; 4])
\# 4
3
2
1

## 2-D matrix along dimension 1

flip ([1 2; 3 4])

```
```

    m 3 4
        1 2
    
## 2-D matrix along dimension 2

flip ([1 2; 3 4], 2)
\# 2 1
4 3

```

See also: [fliplr], page 515, [flipud], page 515, [rot90], page 516, [rotdim], page 516, [permute], page 518, [transpose], page 164.
rot90 (A)
rot90 (A, k)
Rotate array by 90 degree increments.
Return a copy of \(A\) with the elements rotated counterclockwise in 90 -degree increments.
The second argument is optional, and specifies how many 90 -degree rotations are to be applied (the default value is 1 ). Negative values of \(k\) rotate the matrix in a clockwise direction. For example,
```

rot90 ([1, 2; 3, 4], -1)
\# 3 1
4 2

```
rotates the given matrix clockwise by 90 degrees. The following are all equivalent statements:
```

rot90 ([1, 2; 3, 4], -1)
rot90 ([1, 2; 3, 4], 3)
rot90 ([1, 2; 3, 4], 7)

```

The rotation is always performed on the plane of the first two dimensions, i.e., rows and columns. To perform a rotation on any other plane, use rotdim.
See also: [rotdim], page 516, [fliplr], page 515, [flipud], page 515, [flip], page 515.
rotdim ( \(x\) )
rotdim ( \(x, n\) )
rotdim (x, n, plane)
Return a copy of \(x\) with the elements rotated counterclockwise in 90 -degree increments.
The second argument \(n\) is optional, and specifies how many 90 -degree rotations are to be applied (the default value is 1 ). Negative values of \(n\) rotate the matrix in a clockwise direction.
The third argument is also optional and defines the plane of the rotation. If present, plane is a two element vector containing two different valid dimensions of the matrix. When plane is not given the first two non-singleton dimensions are used.
For example,
```

rotdim ([1, 2; 3, 4], -1, [1, 2])
m 3 1
4 2

```
rotates the given matrix clockwise by 90 degrees. The following are all equivalent statements:
```

rotdim ([1, 2; 3, 4], -1, [1, 2])
rotdim ([1, 2; 3, 4], 3, [1, 2])
rotdim ([1, 2; 3, 4], 7, [1, 2])

```

See also: [rot90], page 516, [fliplr], page 515, [flipud], page 515, [flip], page 515.
cat (dim, array1, array2, ..., arrayN)
Return the concatenation of N-D array objects, array1, array \(2, \ldots\), arrayN along dimension dim.
```

A = ones (2, 2);
B = zeros (2, 2);
cat (2, A, B)
\# 1 1 0 0
1100

```

Alternatively, we can concatenate \(A\) and \(B\) along the second dimension in the following way:
\[
[\mathrm{A}, \mathrm{~B}]
\]
dim can be larger than the dimensions of the N-D array objects and the result will thus have dim dimensions as the following example shows:
```

cat (4, ones (2, 2), zeros (2, 2))
$\Rightarrow \operatorname{ans}(:,:, 1,1)=$
11
11
ans(:, : , 1, 2) =
00
00

```

See also: [horzcat], page 517, [vertcat], page 517.
horzcat (array1, array2, ..., arrayN)
Return the horizontal concatenation of N-D array objects, array1, array \(2, \ldots\), array \(N\) along dimension 2 .
Arrays may also be concatenated horizontally using the syntax for creating new matrices. For example:
\[
\text { hcat }=[\text { array1, array2, ... ] }
\]

See also: [cat], page 517, [vertcat], page 517.
vertcat (array1, array2, ..., arrayN)
Return the vertical concatenation of N-D array objects, array1, array \(2, \ldots\), array \(N\) along dimension 1 .

Arrays may also be concatenated vertically using the syntax for creating new matrices. For example:
```

vcat = [ array1; array2; ... ]

```

See also: [cat], page 517, [horzcat], page 517.
permute (A, perm)
Return the generalized transpose for an N-D array object \(A\).
The permutation vector perm must contain the elements 1:ndims (A) (in any order, but each element must appear only once). The \(N\) th dimension of \(A\) gets remapped to dimension \(\operatorname{PERM}(N)\). For example:
```

x = zeros ([2, 3, 5, 7]);
size (x)
\#
size (permute (x, [2, 1, 3, 4]))
\#
size (permute (x, [1, 3, 4, 2]))
\# 2 5 5 7 3

## The identity permutation

size (permute (x, [1, 2, 3, 4]))
m}\begin{array}{lllll}{2}\&{3}\&{5}\&{7}

```

See also: [ipermute], page 518.
ipermute ( \(A\), iperm)
The inverse of the permute function.
The expression
ipermute (permute (A, perm), perm)
returns the original array \(A\).
See also: [permute], page 518.
reshape ( \(A, m, n, \ldots\) )
reshape ( \(A,[m \cap \ldots]\) )
reshape ( \(A, \ldots,[], \ldots\) )
reshape (A, size)
Return a matrix with the specified dimensions ( \(m, n, \ldots\) ) whose elements are taken from the matrix \(A\).
The elements of the matrix are accessed in column-major order (like Fortran arrays are stored).
The following code demonstrates reshaping a 1 x 4 row vector into a 2 x 2 square matrix.
```

reshape ([1, 2, 3, 4], 2, 2)
\# 1 3
24

```

Note that the total number of elements in the original matrix (prod (size (A))) must match the total number of elements in the new matrix ( \(\operatorname{prod}([m n \ldots])\) ).
A single dimension of the return matrix may be left unspecified and Octave will determine its size automatically. An empty matrix ([]) is used to flag the unspecified dimension.
See also: [resize], page 519, [vec], page 524, [postpad], page 525, [cat], page 517, [squeeze], page 49.
resize ( \(x, m\) )
resize ( \(x, m, n, \ldots\) )
resize ( \(x,[m n \ldots]\) )
Resize \(x\) cutting off elements as necessary.
In the result, element with certain indices is equal to the corresponding element of \(x\) if the indices are within the bounds of \(x\); otherwise, the element is set to zero.
In other words, the statement
\[
y=\operatorname{resize}(x, d v)
\]
is equivalent to the following code:
```

y = zeros (dv, class (x));
sz = min (dv, size (x));
for i = 1:length (sz)
idx{i} = 1:sz(i);
endfor
y(idx{:}) = x(idx{:});

```
but is performed more efficiently.
If only \(m\) is supplied, and it is a scalar, the dimension of the result is \(m\)-by- \(m\). If \(m\), \(n, \ldots\) are all scalars, then the dimensions of the result are \(m\)-by- \(n\)-by-. ... If given a vector as input, then the dimensions of the result are given by the elements of that vector.
An object can be resized to more dimensions than it has; in such case the missing dimensions are assumed to be 1 . Resizing an object to fewer dimensions is not possible.
See also: [reshape], page 518, [postpad], page 525, [prepad], page 524, [cat], page 517.
\(y=\operatorname{circshift}(x, n)\)
\(y=\operatorname{circshift~(x,~n,~dim)~}\)
Circularly shift the values of the array \(x\).
\(n\) must be a vector of integers no longer than the number of dimensions in \(x\). The values of \(n\) can be either positive or negative, which determines the direction in which the values of \(x\) are shifted. If an element of \(n\) is zero, then the corresponding dimension of \(x\) will not be shifted. If \(n\) is a scalar and no dim is specified then the shift is applied to the first non-singular dimension.
If a scalar \(d i m\) is given then operate along the specified dimension. In this case \(n\) must be a scalar as well.
Examples:
\[
x=[1,2,3
\]
```

        4, 5, 6;
        7, 8, 9];
    
## positive shift on rows (1st non-singular dim)

circshift (x, 1)
\#
7 8 9
1 2 3
4 5

## negative shift on rows (1st non-singular dim)

circshift (x, -2)
\#
7 8 9
1 2 3
4 5

## no shift of rows, shift columns by 1 (2nd dimension)

circshift (x, [0,1])
\#
3 1 2
6 4 5
9 7 8

## shift columns (2nd dimension)

circshift (x, 1, 2)
\#
3 1 2
6 4 5
9 7 8

```

See also: [permute], page 518, [ipermute], page 518, [shiftdim], page 520.
shift ( \(x, b\) )
shift ( \(x, b, d i m\) )
If \(x\) is a vector, perform a circular shift of length \(b\) of the elements of \(x\).
If \(x\) is a matrix, do the same for each column of \(x\).
If the optional dim argument is given, operate along this dimension.
\(y=\operatorname{shiftdim}(x, n)\)
[y, ns] = shiftdim (x)
Shift the dimensions of \(x\) by \(n\), where \(n\) must be an integer scalar.
When \(n\) is positive, the dimensions of \(x\) are shifted to the left, with the leading dimensions circulated to the end. If \(n\) is negative, then the dimensions of \(x\) are shifted to the right, with \(n\) leading singleton dimensions added.

Called with a single argument, shiftdim, removes the leading singleton dimensions, returning the number of dimensions removed in the second output argument ns.

For example:
```

x = ones (1, 2, 3);
size (shiftdim (x, -1))
\#}1\begin{array}{llll}{1}\&{1}\&{2}\&{3}
size (shiftdim (x, 1))
\# 2 3
[b, ns] = shiftdim (x)
b b =
1 1 1
1 1 1
\# ns = 1

```

See also: [reshape], page 518, [permute], page 518, [ipermute], page 518, [circshift], page 519, [squeeze], page 49.
\([s, i]=\operatorname{sort}(x)\)
[s, i] \(=\operatorname{sort}(x, \operatorname{dim})\)
[s, i] \(=\operatorname{sort}(x\), mode \()\)
[s, i] = sort (x, dim, mode)
Return a copy of \(x\) with the elements arranged in increasing order.
For matrices, sort orders the elements within columns
For example:
```

sort ([1, 2; 2, 3; 3, 1])
\# 1 1
2 2
3 3

```

If the optional argument dim is given, then the matrix is sorted along the dimension defined by dim. The optional argument mode defines the order in which the values will be sorted. Valid values of mode are "ascend" or "descend".

The sort function may also be used to produce a matrix containing the original row indices of the elements in the sorted matrix. For example:
```

[s,i] = sort ([1, 2; 2, 3; 3, 1])
ms=1 1
2 2
3 3
\# i = 1 3
2 1
3 2

```

For equal elements, the indices are such that equal elements are listed in the order in which they appeared in the original list.

Sorting of complex entries is done first by magnitude (abs (z)) and for any ties by phase angle (angle (z)). For example:
```

sort ([1+i; 1; 1-i])
=> 1 + 0i
1 - 1i
1 + 1i

```

NaN values are treated as being greater than any other value and are sorted to the end of the list.
The sort function may also be used to sort strings and cell arrays of strings, in which case ASCII dictionary order (uppercase 'A' precedes lowercase 'a') of the strings is used.
The algorithm used in sort is optimized for the sorting of partially ordered lists.
See also: [sortrows], page 522, [issorted], page 522.
[s, i] = sortrows (A)
[s, i] = sortrows (A, c)
Sort the rows of the matrix \(A\) according to the order of the columns specified in \(c\).
By default ( \(c\) omitted, or a particular column unspecified in \(c\) ) an ascending sort order is used. However, if elements of \(c\) are negative then the corresponding column is sorted in descending order. If the elements of \(A\) are strings then a lexicographical sort is used.
Example: sort by column 2 in descending order, then 3 in ascending order
```

x = [ 7, 1, 4;
8, 3, 5;
9, 3, 6 ];
sortrows (x, [-2, 3])
\# 8 3 5
9 3 6
7 1 4

```

See also: [sort], page 521.
issorted (a)
issorted (a, mode)
issorted (a, "rows", mode)
Return true if the array is sorted according to mode, which may be either "ascend", "descend", or "either".
By default, mode is "ascend". NaNs are treated in the same manner as sort.
If the optional argument "rows" is supplied, check whether the array is sorted by rows as output by the function sortrows (with no options).
This function does not support sparse matrices.
See also: [sort], page 521, [sortrows], page 522.
nth_element ( \(\mathrm{x}, \mathrm{n}\) )
nth_element ( \(x, n\), dim)
Select the n -th smallest element of a vector, using the ordering defined by sort.
The result is equivalent to sort \((x)(n)\).
\(n\) can also be a contiguous range, either ascending \(1: u\) or descending \(u:-1: 1\), in which case a range of elements is returned.
If \(x\) is an array, nth_element operates along the dimension defined by dim, or the first non-singleton dimension if dim is not given.

Programming Note: nth_element encapsulates the C++ standard library algorithms nth_element and partial_sort. On average, the complexity of the operation is \(\mathrm{O}\left(\mathrm{M}^{*} \log (\mathrm{~K})\right)\), where \(\mathrm{M}=\operatorname{size}(\mathrm{x}, \operatorname{dim})\) and \(\mathrm{K}=\) length ( n\()\). This function is intended for cases where the ratio \(K / M\) is small; otherwise, it may be better to use sort.

See also: [sort], page 521, [min], page 559, [max], page 558.
\(A_{-} L O=\operatorname{tril}(A)\)
\(A_{-} L O=\operatorname{tril}(A, k)\)
\(A_{-} L O=\operatorname{tril}(A, k, p a c k)\)
Return a new matrix formed by extracting the lower triangular part of the matrix \(A\), and setting all other elements to zero.
The optional second argument specifies how many diagonals above or below the main diagonal should also be set to zero. The default value of \(k\) is zero which includes the main diagonal as part of the result. If the value of \(k\) is a nonzero integer then the selection of elements starts at an offset of \(k\) diagonals above the main diagonal for positive \(k\) or below the main diagonal for negative \(k\). The absolute value of \(k\) may not be greater than the number of subdiagonals or superdiagonals.
Example 1: exclude main diagonal
```

tril (ones (3), -1)
=> 0 0 0
1 0 0
1 1 0

```

Example 2 : include first superdiagonal
```

tril (ones (3), 1)
\# 1 1 0
1 1 1
1 1 1

```

If the optional third argument "pack" is given then the extracted elements are not inserted into a matrix, but instead stacked column-wise one above another, and returned as a column vector.

See also: [triu], page 523, [istril], page 70, [diag], page 525.
\(A_{-} U P=\operatorname{triu}(A)\)
\(A_{-} U P=\operatorname{triu}(A, k)\)
\(A_{-} U P=\operatorname{triu}(A, k, p a c k)\)
Return a new matrix formed by extracting the upper triangular part of the matrix \(A\), and setting all other elements to zero.
The optional second argument specifies how many diagonals above or below the main diagonal should also be set to zero. The default value of \(k\) is zero which includes the main diagonal as part of the result. If the value of \(k\) is a nonzero integer then the selection of elements starts at an offset of \(k\) diagonals above the main diagonal for positive \(k\) or below the main diagonal for negative \(k\). The absolute value of \(k\) may not be greater than the number of subdiagonals or superdiagonals.

Example 1 : exclude main diagonal
```

triu (ones (3), 1)
$\Rightarrow \quad 0 \quad 1 \quad 1$
$0 \quad 0 \quad 1$
000

```

Example 2 : include first subdiagonal
```

triu (ones (3), -1)
m 1 1 1
1 1 1
0 1 1

```

If the optional third argument "pack" is given then the extracted elements are not inserted into a matrix, but instead stacked column-wise one above another, and returned as a column vector.

See also: [tril], page 523, [istriu], page 70, [diag], page 525.
\(\mathrm{v}=\operatorname{vec}(\mathrm{x})\)
\(v=\operatorname{vec}(x, d i m)\)
Return the vector obtained by stacking the columns of the matrix \(x\) one above the other.

Without dim this is equivalent to \(x(:)\).
If dim is supplied, the dimensions of \(v\) are set to dim with all elements along the last dimension. This is equivalent to shiftdim ( \(x(:), 1-\operatorname{dim}\) ).

See also: [vech], page 524, [resize], page 519, [cat], page 517.
vech (x)
Return the vector obtained by eliminating all superdiagonal elements of the square matrix \(x\) and stacking the result one column above the other.
This has uses in matrix calculus where the underlying matrix is symmetric and it would be pointless to keep values above the main diagonal.

See also: [vec], page 524.
prepad ( \(x, 1\) )
prepad ( \(x, 1, c\) )
prepad (x, l, c, dim)
Prepend the scalar value \(c\) to the vector \(x\) until it is of length \(l\). If \(c\) is not given, a value of 0 is used.

If length ( \(x\) ) > l, elements from the beginning of \(x\) are removed until a vector of length \(l\) is obtained.
If \(x\) is a matrix, elements are prepended or removed from each row.
If the optional argument dim is given, operate along this dimension.
If dim is larger than the dimensions of \(x\), the result will have dim dimensions.
See also: [postpad], page 525, [cat], page 517, [resize], page 519.
postpad (x, 1)
postpad ( \(x, 1, c\) )
postpad (x, l, c, dim)
Append the scalar value \(c\) to the vector \(x\) until it is of length \(l\). If \(c\) is not given, a value of 0 is used.
If length \((x)>1\), elements from the end of \(x\) are removed until a vector of length \(l\) is obtained.
If \(x\) is a matrix, elements are appended or removed from each row.
If the optional argument dim is given, operate along this dimension.
If dim is larger than the dimensions of \(x\), the result will have dim dimensions.
See also: [prepad], page 524, [cat], page 517, [resize], page 519.
\(M=\operatorname{diag}(v)\)
\(M=\operatorname{diag}(v, k)\)
\(M=\operatorname{diag}(v, m, n)\)
\(v=\operatorname{diag}(M)\)
\(v=\operatorname{diag}(M, k)\)
Return a diagonal matrix with vector \(v\) on diagonal \(k\).
The second argument is optional. If it is positive, the vector is placed on the \(k\)-th superdiagonal. If it is negative, it is placed on the \(-k\)-th subdiagonal. The default value of \(k\) is 0 , and the vector is placed on the main diagonal. For example:
\[
\begin{aligned}
& \text { diag ([1, 2, 3], 1) } \\
& \Rightarrow \quad 0 \quad 1 \quad 0 \quad 0 \\
& 0 \quad 0 \quad 2 \quad 0 \\
& 0 \quad 0 \quad 0 \quad 3 \\
& 0 \quad 0 \quad 0 \quad 0
\end{aligned}
\]

The 3 -input form returns a diagonal matrix with vector \(v\) on the main diagonal and the resulting matrix being of size \(m\) rows \(x n\) columns.
Given a matrix argument, instead of a vector, diag extracts the \(k\)-th diagonal of the matrix.
blkdiag ( \(A, B, C, \ldots\) )
Build a block diagonal matrix from \(A, B, C, \ldots\)
All arguments must be numeric and either two-dimensional matrices or scalars. If any argument is of type sparse, the output will also be sparse.
See also: [diag], page 525, [horzcat], page 517, [vertcat], page 517, [sparse], page 664.

\subsection*{16.3 Special Utility Matrices}
eye ( \(n\) )
eye ( \(m, n\) )
eye ( \([m n]\) )
eye (..., class)
Return an identity matrix.

If invoked with a single scalar argument \(n\), return a square NxN identity matrix.
If supplied two scalar arguments \((m, n)\), eye takes them to be the number of rows and columns. If given a vector with two elements, eye uses the values of the elements as the number of rows and columns, respectively. For example:
```

eye (3)
\# 1 0 0
0}1
0 0 1

```

The following expressions all produce the same result:
```

eye (2)
\equiv
eye (2, 2)
\equiv
eye (size ([1, 2; 3, 4]))

```

The optional argument class, allows eye to return an array of the specified type, like
```

val = zeros (n,m, "uint8")

```

Calling eye with no arguments is equivalent to calling it with an argument of 1. Any negative dimensions are treated as zero. These odd definitions are for compatibility with matlab.
See also: [speye], page 662, [ones], page 526, [zeros], page 526.
```

val = ones (n)
val = ones (m, n)
val = ones (m, n, k, ...)
val = ones ([mn...])
val = ones (..., "like", var))
val = ones (..., class)

```

Return a matrix or N-dimensional array whose elements are all 1.
If invoked with a single scalar integer argument \(n\), return a square NxN matrix.
If invoked with two or more scalar integer arguments, or a vector of integer values, return an array with the given dimensions.
To create a constant matrix whose values are all the same use an expression such as
```

val_matrix = val * ones (m, n)

```

If a variable var is specified after "like", the output val will have the same data type, complexity, and sparsity as var.
The optional argument class specifies the class of the return array and defaults to double. For example:
```

val = ones (m,n, "uint8")

```

See also: [zeros], page 526.
```

val = zeros (n)
val = zeros (m, n)
val = zeros (m, n, k,···)
val = zeros ([mn...])

```
val = zeros (..., "like", var))
val \(=\) zeros (..., class)
Return a matrix or N-dimensional array whose elements are all 0.
If invoked with a single scalar integer argument, return a square \(\mathrm{NxN}^{\mathrm{N}}\) matrix.
If invoked with two or more scalar integer arguments, or a vector of integer values, return an array with the given dimensions.
If a variable var is specified after "like", the output val will have the same data type, complexity, and sparsity as var.
The optional argument class specifies the class of the return array and defaults to double. For example:
```

val = zeros (m,n, "uint8")

```

See also: [ones], page 526 .
repmat ( \(A, m\) )
repmat ( \(A, m, n\) )
repmat ( \(A, m, n, p \ldots\) )
repmat (A, [mn])
repmat ( \(A,[m \mathrm{n} p \ldots]\) )
Repeat matrix or N-D array.
Form a block matrix of size \(m\) by \(n\), with a copy of matrix \(A\) as each element.
If \(n\) is not specified, form an \(m\) by \(m\) block matrix. For copying along more than two dimensions, specify the number of times to copy across each dimension \(m, n, p, \ldots\), in a vector in the second argument.

See also: [bsxfun], page 628, [kron], page 610, [repelems], page 527.
repelems ( \(x, r\) )
Construct a vector of repeated elements from \(x\).
\(r\) is a \(2 \times N\) integer matrix specifying which elements to repeat and how often to repeat each element. Entries in the first row, \(r(1, \mathrm{j})\), select an element to repeat. The corresponding entry in the second row, \(r(2, j)\), specifies the repeat count. If \(x\) is a matrix then the columns of \(x\) are imagined to be stacked on top of each other for purposes of the selection index. A row vector is always returned.
Conceptually the result is calculated as follows:
```

y = [];
for i = 1:columns (r)
y = [y, x(r(1,i)*ones(1, r(2,i)))];
endfor

```

See also: [repmat], page 527, [cat], page 517.
xxx \(=\) repelem \((x, R)\)
xxx \(=\) repelem ( \(x, R_{-} 1, \ldots, R_{-} n\) )
Construct an array of repeated elements from \(x\) and repeat instructions \(R_{-} 1, \ldots\).
\(x\) must be a scalar, vector, or N -dimensional array.

A repeat instruction \(R_{-} j\) must either be a scalar or a vector. If the instruction is a scalar then each component of \(x\) in dimension \(j\) is repeated \(R_{-j}\) times. If the instruction is a vector then it must have the same number of elements as the corresponding dimension \(j\) of \(x\). In this case, the \(k\) th component of dimension \(j\) is repeated \(R_{-} j(k)\) times.

If \(x\) is a scalar or vector then repelem may be called with just a single repeat instruction \(R\) and repelem will return a vector with the same orientation as the input.

If \(x\) is a matrix then at least two \(R_{-}\)js must be specified.

Note: Using repelem with a vector x and a vector for \(R_{-} j\) is equivalent to Run Length Decoding.

Examples:
```

A = [lllllll
B = [2 1 0 1 2];
repelem (A, B)
\#
A = magic (3)
A A =
8 1 6
3 5 7
B1 = [llll
B2 = 2;
repelem (A, B1, B2)
\#
3
3
4
4

```

More \(R_{-}\)may be specified than the number of dimensions of x. Any excess \(R_{-} j\) must be scalars (because \(x\) 's size in those dimensions is only 1 ), and \(x\) will be replicated in those dimensions accordingly.
```

A = [11 2 3 4 5];
B1 = 2;
B2 = [2 [1 3 0 2];
B3 = 3;
repelem (A, B1, B2, B3)
ans(:,:,1) =
1
ans(:,:,2) =
1
ans(:,:,3) =
1

```
\(R_{-j}\) must be specified in order. A placeholder of 1 may be used for dimensions which do not need replication.
```

repelem ([-1, 0; 0, 1], 1, 2, 1, 2)
$\Rightarrow \operatorname{ans}(:,:, 1,1)=$
$\begin{array}{llll}-1 & -1 & 0 & 0\end{array}$
$0 \quad 0 \quad 1 \quad 1$
$\operatorname{ans}(:,:, 1,2)=$
$\begin{array}{llll}-1 & -1 & 0 & 0\end{array}$
$0 \quad 0 \quad 1 \quad 1$

```

If fewer \(R_{-} j\) are given than the number of dimensions in \(x\), repelem will assume \(R_{-} j\) is 1 for those dimensions.
```

A = cat (3, [-1 0; 0 1], [-1 0; 0 1])
a ans(:,:,1) =
-1 0
0}
ans(:,:,2) =
-1 0
0}
repelem (A,2,3)
ans(:,:,1) =
-1
-1
0
ans(:,:,2) =

```
\begin{tabular}{rrrrrr}
-1 & -1 & -1 & 0 & 0 & 0 \\
-1 & -1 & -1 & 0 & 0 & 0 \\
0 & 0 & 0 & 1 & 1 & 1 \\
0 & 0 & 0 & 1 & 1 & 1
\end{tabular}
repelem preserves the class of \(x\), and works with strings, cell arrays, NA, and NAN inputs. If any \(R_{-j}\) is 0 the output will be an empty array.
```

repelem ("Octave", 2, 3)
\# OOOccctttaaavvveee
000ccctttaaavvveee
repelem ([1 2 3; 1 2 3], 2, 0)
[ [] (4x0)

```

See also: [cat], page 517, [kron], page 610, [repmat], page 527.
The functions linspace and logspace make it very easy to create vectors with evenly or logarithmically spaced elements. See Section 4.2 [Ranges], page 56.
linspace (start, end)
linspace (start, end, n)
Return a row vector with \(n\) linearly spaced elements between start and end.
If the number of elements is greater than one, then the endpoints start and end are always included in the range. If start is greater than end, the elements are stored in decreasing order. If the number of points is not specified, a value of 100 is used.
The linspace function returns a row vector when both start and end are scalars. If one, or both, inputs are vectors, then linspace transforms them to column vectors and returns a matrix where each row is an independent sequence between start(row_n), end(row_n).
For compatibility with matlab, return the second argument (end) when only a single value ( \(n=1\) ) is requested.

See also: [colon], page 898, [logspace], page 530.
logspace ( \(a, b\) )
logspace ( \(a, b, n\) )
logspace (a, pi, n)
Return a row vector with \(n\) elements logarithmically spaced from \(10^{a}\) to \(10^{b}\).
If \(n\) is unspecified it defaults to 50 .
If \(b\) is equal to \(\pi\), the points are between \(10^{a}\) and \(\pi\), not \(10^{a}\) and \(10^{\pi}\), in order to be compatible with the corresponding matlab function.
Also for compatibility with MATLAB, return the right-hand side of the range \(\left(10^{b}\right)\) when just a single value is requested.
See also: [linspace], page 530.
rand (n)
rand ( \(m, n, \ldots\) )
rand ([mn...])
```

v = rand ("state")
rand ("state", v)
rand ("state", "reset")
v = rand ("seed")
rand ("seed", v)
rand ("seed", "reset")
rand (..., "single")
rand (..., "double")

```

Return a matrix with random elements uniformly distributed on the interval \((0,1)\).
The arguments are handled the same as the arguments for eye.
You can query the state of the random number generator using the form
```

v = rand ("state")

```

This returns a column vector \(v\) of length 625. Later, you can restore the random number generator to the state \(v\) using the form
```

rand ("state", v)

```

You may also initialize the state vector from an arbitrary vector of length \(\leq 625\) for \(v\). This new state will be a hash based on the value of \(v\), not \(v\) itself.
By default, the generator is initialized by contributing entropy from the wall clock time, the CPU time, the current fraction of a second, the process ID and-if available - up to 1024 bits from the C++ random numbers source random_device, which might be non-deterministic (implementation specific). Note that this differs from matlab, which always initializes the state to the same state at startup. To obtain behavior comparable to matLab, initialize with a deterministic state vector in Octave's startup files (see Section 2.1.2 [Startup Files], page 18).
To compute the pseudo-random sequence, rand uses the Mersenne Twister with a period of \(2^{19937}-1\) (See M. Matsumoto and T. Nishimura, Mersenne Twister: A 623-dimensionally equidistributed uniform pseudorandom number generator, ACM Trans. on Modeling and Computer Simulation Vol. 8, No. 1, pp. 3-30, January 1998, http://www.math.sci.hiroshima-u.ac.jp/~m-mat/MT/emt.html). Do not use for cryptography without securely hashing several returned values together, otherwise the generator state can be learned after reading 624 consecutive values.
Older versions of Octave used a different random number generator. The new generator is used by default as it is significantly faster than the old generator, and produces random numbers with a significantly longer cycle time. However, in some circumstances it might be desirable to obtain the same random sequences as produced by the old generators. To do this the keyword "seed" is used to specify that the old generators should be used, as in
rand ("seed", val)
which sets the seed of the generator to val. The seed of the generator can be queried with
```

s = rand ("seed")

```

However, it should be noted that querying the seed will not cause rand to use the old generators, only setting the seed will. To cause rand to once again use the new generators, the keyword "state" should be used to reset the state of the rand.

The state or seed of the generator can be reset to a new random value using the "reset" keyword.
The class of the value returned can be controlled by a trailing "double" or "single" argument. These are the only valid classes.
See also: [randn], page 532, [rande], page 533, [randg], page 534, [randp], page 533.
randi (imax)
randi (imax, n)
randi (imax, m, n,...)
randi ([imin imax], ...)
randi (..., "class")
Return random integers in the range 1:imax.
Additional arguments determine the shape of the return matrix. When no arguments are specified a single random integer is returned. If one argument \(n\) is specified then a square matrix ( \(n \times n\) ) is returned. Two or more arguments will return a multidimensional matrix ( \(m \times n \times \ldots\) ).
The integer range may optionally be described by a two element matrix with a lower and upper bound in which case the returned integers will be on the interval [imin, imax].
The optional argument class will return a matrix of the requested type. The default is "double".
The following example returns 150 integers in the range 1-10.
```

ri = randi (10, 150, 1)

```

Implementation Note: randi relies internally on rand which uses class "double" to represent numbers. This limits the maximum integer (imax) and range (imax - imin) to the value returned by the flintmax function. For IEEE floating point numbers this value is \(2^{53}-1\).

See also: [rand], page 530.
```

randn (n)
randn (m, n, ...)
randn ([m n ...])
v = randn ("state")
randn ("state", v)
randn ("state", "reset")
v = randn ("seed")
randn ("seed", v)
randn ("seed", "reset")
randn (..., "single")
randn (..., "double")

```

Return a matrix with normally distributed random elements having zero mean and variance one.
The arguments are handled the same as the arguments for rand.
By default, randn uses the Marsaglia and Tsang "Ziggurat technique" to transform from a uniform to a normal distribution.

The class of the value returned can be controlled by a trailing "double" or "single" argument. These are the only valid classes.
Reference: G. Marsaglia and W.W. Tsang, Ziggurat Method for Generating Random
Variables, J. Statistical Software, vol 5, 2000, https://www.jstatsoft.org/v05/ i08/

See also: [rand], page 530, [rande], page 533, [randg], page 534, [randp], page 533.
```

rande (n)
rande (m, n, ...)
rande ([m n ...])
v = rande ("state")
rande ("state", v)
rande ("state", "reset")
v = rande ("seed")
rande ("seed", v)
rande ("seed", "reset")
rande (..., "single")
rande (..., "double")

```

Return a matrix with exponentially distributed random elements.
The arguments are handled the same as the arguments for rand.
By default, rande uses the Marsaglia and Tsang "Ziggurat technique" to transform from a uniform to an exponential distribution.

The class of the value returned can be controlled by a trailing "double" or "single" argument. These are the only valid classes.

Reference: G. Marsaglia and W.W. Tsang, Ziggurat Method for Generating Random Variables, J. Statistical Software, vol 5, 2000, https://www.jstatsoft.org/v05/ i08/

See also: [rand], page 530, [randn], page 532, [randg], page 534, [randp], page 533.
```

randp (1, n)
randp (l, m, n, ...)
randp (l, [m n ...])
v = randp ("state")
randp ("state", v)
randp ("state", "reset")
v = randp ("seed")
randp ("seed", v)
randp ("seed", "reset")
randp (..., "single")
randp (. .., "double")

```

Return a matrix with Poisson distributed random elements with mean value parameter given by the first argument, \(l\).
The arguments are handled the same as the arguments for rand, except for the argument \(l\).

Five different algorithms are used depending on the range of \(l\) and whether or not \(l\) is a scalar or a matrix.

For scalar \(l \leq 12\), use direct method.
W.H. Press, et al., Numerical Recipes in C, Cambridge University Press, 1992.

For scalar l>12, use rejection method.[1]
W.H. Press, et al., Numerical Recipes in C, Cambridge University Press, 1992.

For matrix \(l \leq 10\), use inversion method.[2]
E. Stadlober, et al., WinRand source code, available via FTP.

For matrix l>10, use patchwork rejection method.
E. Stadlober, et al., WinRand source code, available via FTP, or H. Zechner, Efficient sampling from continuous and discrete unimodal distributions, Doctoral Dissertation, 156pp., Technical University Graz, Austria, 1994.

For \(l>1 \mathrm{e} 8\), use normal approximation.
L. Montanet, et al., Review of Particle Properties, Physical Review D 50 p1284, 1994.

The class of the value returned can be controlled by a trailing "double" or "single" argument. These are the only valid classes.
See also: [rand], page 530, [randn], page 532, [rande], page 533, [randg], page 534.
```

randg (a, n)
randg (a,m, n, ...)
randg (a,[mn...])
v = randg ("state")
randg ("state", v)
randg ("state", "reset")
v = randg ("seed")
randg ("seed", v)
randg ("seed", "reset")
randg (..., "single")
randg (..., "double")

```

Return a matrix with gamma ( \(a, 1\) ) distributed random elements.
The arguments are handled the same as the arguments for rand, except for the argument a.
This can be used to generate many distributions:
```

gamma (a, b) for a > -1, b > 0
r = b * randg (a)
beta (a, b) for a > -1, b > -1
r1 = randg (a, 1)
r = r1 / (r1 + randg (b, 1))

```

Erlang (a, n)
\[
\mathrm{r}=\mathrm{a} * \mathrm{randg}(\mathrm{n})
\]
chisq (df) for \(d f>0\)
\[
r=2 * \operatorname{randg}(d f / 2)
\]
t (df) for \(0<\mathrm{df}<\inf\) (use randn if df is infinite)
\[
r=\operatorname{randn}() / \operatorname{sqrt}(2 * \operatorname{randg}(d f / 2) / d f)
\]

F ( \(\mathrm{n} 1, \mathrm{n} 2\) ) for \(0<\mathrm{n} 1,0<\mathrm{n} 2\)
\#\# r1 equals 1 if n1 is infinite
\(r 1=2 * \operatorname{randg}(n 1 / 2) / n 1\)
\#\# r2 equals 1 if n2 is infinite
r2 \(=2\) * randg (n2 / 2) / n2
\(r=r 1 / r 2\)
negative binomial ( \(\mathrm{n}, \mathrm{p}\) ) for \(\mathrm{n}>0,0<\mathrm{p}<=1\)
```

r = randp ((1 - p) / p * randg (n))

```
non-central chisq ( \(\mathrm{df}, \mathrm{L}\) ), for \(\mathrm{df}>=0\) and \(L>0\)
(use chisq if \(L=0\) )
\(r=r a n d p(L / 2)\)
\(r(r>0)=2 * \operatorname{randg}(r(r>0))\)
\(r(d f>0)+=2 * \operatorname{randg}(d f(d f>0) / 2)\)
Dirichlet (a1, ... ak)
\(r=(r a n d g(a 1), \ldots, r a n d g(a k))\)
\(r=r / \operatorname{sum}(r)\)
The class of the value returned can be controlled by a trailing "double" or "single" argument. These are the only valid classes.

See also: [rand], page 530, [randn], page 532, [rande], page 533, [randp], page 533.
```

rng (seed)
rng (seed, "generator")
rng ("shuffle")
rng ("shuffle", "generator")
rng ("default")
s = rng ()
rng (s)
s = rng (...)

```

Set or query the seed of the random number generator used by rand and randn.
The input seed is a scalar numeric value used to initialize the state vector of the random number generator.
The optional string generator specifies the type of random number generator to be used. Its value can be "twister", "v5uniform", or "v5normal". The "twister" keyword is described below. "v5uniform" and "v5normal" refer to older versions of Octave that used to use a different random number generator.
The state or seed of the random number generator can be reset to a new random value using the "shuffle" keyword.

The random number generator can be reset to default values using the "default" keyword. The default values are to use the Mersenne Twister generator with a seed of 0 .
The optional return value \(s\) contains the state of the random number generator at the time the function is called (i.e., before it might be modified according to the input arguments). It is encoded as a structure variable with three fields: "Type", "Seed", and "State". The random number generator can be restored to the state \(s\) using rng ( \(s\) ). This is useful when the identical sequence of pseudo-random numbers is required for an algorithm.
By default, and with the "twister" option, pseudo-random sequences are computed using the Mersenne Twister with a period of \(2^{19937}-1\) (See M. Matsumoto and T. Nishimura, Mersenne Twister: A 623-dimensionally equidistributed uniform pseudorandom number generator, ACM Trans. on Modeling and Computer Simulation Vol. 8, No. 1, pp. 3-30, January 1998, http://www.math.sci.hiroshima-u.ac.jp/ \({ }^{\sim} \mathrm{m}\)-mat/MT/emt.html). Do not use for cryptography without securely hashing several returned values together, otherwise the generator state can be learned after reading 624 consecutive values.
See also: [rand], page 530, [randn], page 532.
The generators operate in the new or old style together, it is not possible to mix the two. Initializing any generator with "state" or "seed" causes the others to switch to the same style for future calls.

The state of each generator is independent and calls to different generators can be interleaved without affecting the final result. For example,
```

rand ("state", [11, 22, 33]);
randn ("state", [44, 55, 66]);
u = rand (100, 1);
n = randn (100, 1);

```
```

rand ("state", [11, 22, 33]);

```
rand ("state", [11, 22, 33]);
randn ("state", [44, 55, 66]);
randn ("state", [44, 55, 66]);
u = zeros (100, 1);
u = zeros (100, 1);
n = zeros (100, 1);
n = zeros (100, 1);
for i = 1:100
for i = 1:100
    u(i) = rand ();
    u(i) = rand ();
    n(i) = randn ();
    n(i) = randn ();
end
```

end

```
and
produce equivalent results. When the generators are initialized in the old style with "seed" only rand and randn are independent, because the old rande, randg and randp generators make calls to rand and randn.

The generators are initialized with random states at start-up, so that the sequences of random numbers are not the same each time you run Octave. \({ }^{1}\) If you really do need to reproduce a sequence of numbers exactly, you can set the state or seed to a specific value.

\footnotetext{
\({ }^{1}\) The old versions of rand and randn obtain their initial seeds from the system clock.
}

If invoked without arguments, rand and randn return a single element of a random sequence.

The original rand and randn functions use Fortran code from Ranlib, a library of Fortran routines for random number generation, compiled by Barry W. Brown and James Lovato of the Department of Biomathematics at The University of Texas, M.D. Anderson Cancer Center, Houston, TX 77030.
randperm (n)
randperm (n, m)
Return a row vector containing a random permutation of \(1: n\).
If \(m\) is supplied, return \(m\) unique entries, sampled without replacement from 1:n.
The complexity is \(\mathrm{O}(n)\) in memory and \(\mathrm{O}(m)\) in time, unless \(m<n / 5\), in which case \(\mathrm{O}(m)\) memory is used as well. The randomization is performed using rand(). All permutations are equally likely.
See also: [perms], page 785.

\subsection*{16.4 Famous Matrices}

The following functions return famous matrix forms.
```

gallery (name)
gallery (name, args)
Create interesting matrices for testing.
c = gallery ("cauchy", x)
c = gallery ("cauchy", x, y)
Create a Cauchy matrix.
c = gallery ("chebspec", n)
c = gallery ("chebspec", n, k)
Create a Chebyshev spectral differentiation matrix.
c = gallery ("chebvand", p)
c = gallery ("chebvand", m, p)
Create a Vandermonde-like matrix for the Chebyshev polynomials.
a = gallery ("chow", n)
a = gallery ("chow", n, alpha)
a = gallery ("chow", n, alpha, delta)
Create a Chow matrix - a singular Toeplitz lower Hessenberg matrix.
c = gallery ("circul", v)
Create a circulant matrix.
a = gallery ("clement", n)
a = gallery ("clement", n, k)
Create a tridiagonal matrix with zero diagonal entries.

```
```

c = gallery ("compar", a)
c = gallery ("compar", a, k)
Create a comparison matrix.
a = gallery ("condex", n)
a = gallery ("condex", n, k)
a = gallery ("condex", n, k, theta)
Create a "counterexample" matrix to a condition estimator.
a = gallery ("cycol", [m n])
a = gallery ("cycol", n)
a = gallery (..., k)
Create a matrix whose columns repeat cyclically.
[c, d, e] = gallery ("dorr", n)
[c, d, e] = gallery ("dorr", n, theta)
a = gallery ("dorr",...)
Create a diagonally dominant, ill-conditioned, tridiagonal matrix.
a = gallery ("dramadah", n)
a = gallery ("dramadah", n, k)
Create a (0, 1) matrix whose inverse has large integer entries.
a = gallery ("fiedler", c)
Create a symmetric Fiedler matrix.
a = gallery ("forsythe", n)
a = gallery ("forsythe", n, alpha)
a = gallery ("forsythe", n, alpha, lambda)
Create a Forsythe matrix (a perturbed Jordan block).
f = gallery ("frank", n)
f = gallery ("frank", n, k)
Create a Frank matrix (ill-conditioned eigenvalues).
c = gallery ("gcdmat", n)
Create a greatest common divisor matrix.

```
    \(c\) is an \(n\)-by- \(n\) matrix whose values correspond to the greatest common divisor of its
    coordinate values, i.e., \(c(\mathrm{i}, \mathrm{j})\) correspond \(\operatorname{gcd}(\mathrm{i}, \mathrm{j})\).
a = gallery ("gearmat", n)
a = gallery ("gearmat", n, i)
a = gallery ("gearmat", n, i, j)
    Create a Gear matrix.
g = gallery ("grcar", n)
g = gallery ("grcar", n, k)
    Create a Toeplitz matrix with sensitive eigenvalues.
```

a = gallery ("hanowa", n)
a = gallery ("hanowa", n, d)
Create a matrix whose eigenvalues lie on a vertical line in the complex plane.
v = gallery ("house", x)
[v, beta] = gallery ("house", x)
Create a householder matrix.
a = gallery ("integerdata", imax, [M N ...], j)
a = gallery ("integerdata", imax, M, N, ..., j)
a = gallery ("integerdata", [imin, imax], [M N ...], j)
a = gallery ("integerdata", [imin, imax], M, N, ..., j)
a = gallery ("integerdata", ..., "class")
Create a matrix with random integers in the range [1, imax]. If imin is given then the integers are in the range [imin, imax].
The second input is a matrix of dimensions describing the size of the output. The dimensions can also be input as comma-separated arguments.
The input $j$ is an integer index in the range $\left[0,2^{\wedge} 32-1\right]$. The values of the output matrix are always exactly the same (reproducibility) for a given size input and $j$ index. The final optional argument determines the class of the resulting matrix. Possible values for class: "uint8", "uint16", "uint32", "int8", "int16", int32", "single", "double". The default is "double".
a = gallery ("invhess", x)
a = gallery ("invhess", x, y)
Create the inverse of an upper Hessenberg matrix.
a = gallery ("invol", n)
Create an involutory matrix.
a = gallery ("ipjfact", n)
a = gallery ("ipjfact", n, k)
Create a Hankel matrix with factorial elements.
a = gallery ("jordbloc", n)
a = gallery ("jordbloc", n, lambda)
Create a Jordan block.
u = gallery ("kahan", n)
u = gallery ("kahan", n, theta)
u = gallery ("kahan", n, theta, pert)
Create a Kahan matrix (upper trapezoidal).
a = gallery ("kms", n)
a = gallery ("kms", n, rho)
Create a Kac-Murdock-Szego Toeplitz matrix.
b = gallery ("krylov", a)
b = gallery ("krylov", a, x)
b = gallery ("krylov", a, x, j)
Create a Krylov matrix.

```
```

a = gallery ("lauchli", n)
a = gallery ("lauchli", n, mu)
Create a Lauchli matrix (rectangular).
a = gallery ("lehmer", n)
Create a Lehmer matrix (symmetric positive definite).
t = gallery ("lesp", n)
Create a tridiagonal matrix with real, sensitive eigenvalues.
a = gallery ("lotkin", n)
Create a Lotkin matrix.
a = gallery ("minij", n)
Create a symmetric positive definite matrix $\operatorname{MIN}(\mathrm{i}, \mathrm{j})$.
a = gallery ("moler", n)
a = gallery ("moler", n, alpha)
Create a Moler matrix (symmetric positive definite).
[a, t] = gallery ("neumann", n)
Create a singular matrix from the discrete Neumann problem (sparse).
a = gallery ("normaldata", [M N ...], j)
a = gallery ("normaldata", M, N, ..., j)
a = gallery ("normaldata", ..., "class")

```

Create a matrix with random samples from the standard normal distribution (mean \(=0, \operatorname{std}=1\) ).
The first input is a matrix of dimensions describing the size of the output. The dimensions can also be input as comma-separated arguments.
The input \(j\) is an integer index in the range \(\left[0,2^{\wedge} 32-1\right]\). The values of the output matrix are always exactly the same (reproducibility) for a given size input and \(j\) index.
The final optional argument determines the class of the resulting matrix. Possible values for class: "single", "double". The default is "double".
\(q=\) gallery ("orthog", n)
\(q=\) gallery ("orthog", n, k)
Create orthogonal and nearly orthogonal matrices.
a = gallery ("parter", n)
Create a Parter matrix (a Toeplitz matrix with singular values near pi).
p = gallery ("pei", n)
p = gallery ("pei", n, alpha)
Create a Pei matrix.
a = gallery ("poisson", n)
Create a block tridiagonal matrix from Poisson's equation (sparse).
```

a = gallery ("prolate", n)
a = gallery ("prolate", n, w)
Create a prolate matrix (symmetric, ill-conditioned Toeplitz matrix).
h = gallery ("randhess", x)
Create a random, orthogonal upper Hessenberg matrix.
a = gallery ("rando", n)
a = gallery ("rando", n, k)
Create a random matrix with elements -1, 0 or 1.
a = gallery ("randsvd", n)
a = gallery ("randsvd", n, kappa)
a = gallery ("randsvd", n, kappa, mode)
a = gallery ("randsvd", n, kappa, mode, kl)
a = gallery ("randsvd", n, kappa, mode, kl, ku)
Create a random matrix with pre-assigned singular values.
a = gallery ("redheff", n)
Create a zero and ones matrix of Redheffer associated with the Riemann hypothesis.
a = gallery ("riemann", n)
Create a matrix associated with the Riemann hypothesis.
a = gallery ("ris", n)
Create a symmetric Hankel matrix.
a = gallery ("smoke", n)
a = gallery ("smoke", n, k)
Create a complex matrix, with a "smoke ring" pseudospectrum.
t = gallery ("toeppd", n)
t = gallery ("toeppd", n, m)
t = gallery ("toeppd", n, m, w)
t = gallery ("toeppd", n, m, w, theta)
Create a symmetric positive definite Toeplitz matrix.
p = gallery ("toeppen", n)
p = gallery ("toeppen", n, a)
p = gallery ("toeppen", n, a, b)
p = gallery ("toeppen", n, a, b, c)
p = gallery ("toeppen", n, a, b, c, d)
p = gallery ("toeppen", n, a, b, c, d, e)
Create a pentadiagonal Toeplitz matrix (sparse).
a = gallery ("tridiag", x, y, z)
a = gallery ("tridiag", n)
a = gallery ("tridiag", n, c, d, e)
Create a tridiagonal matrix (sparse).

```
t = gallery ("triw", n)
t = gallery ("triw", n, alpha)
t = gallery ("triw", n, alpha, k)
Create an upper triangular matrix discussed by Kahan, Golub, and Wilkinson.
a = gallery ("uniformdata", [M N ...], j)
a = gallery ("uniformdata", M, N, ..., j)
a = gallery ("uniformdata", ..., "class")
Create a matrix with random samples from the standard uniform distribution (range \([0,1]\) ).
The first input is a matrix of dimensions describing the size of the output. The dimensions can also be input as comma-separated arguments.
The input \(j\) is an integer index in the range \(\left[0,2^{\wedge} 32-1\right]\). The values of the output matrix are always exactly the same (reproducibility) for a given size input and \(j\) index.
The final optional argument determines the class of the resulting matrix. Possible values for class: "single", "double". The default is "double".
a = gallery ("wathen", nx, ny)
a = gallery ("wathen", nx, ny, k)
Create the Wathen matrix.
[a, b] = gallery ("wilk", n)
Create various specific matrices devised/discussed by Wilkinson.

\section*{hadamard (n)}

Construct a Hadamard matrix (Hn) of size n-by-n.
The size \(n\) must be of the form \(2^{k} * p\) in which p is one of \(1,12,20\) or 28 . The returned matrix is normalized, meaning \(\operatorname{Hn}(:, 1)==1\) and \(\operatorname{Hn}(1,:)==1\).
Some of the properties of Hadamard matrices are:
- kron ( \(\mathrm{Hm}, \mathrm{Hn}\) ) is a Hadamard matrix of size m-by-n.
- \(\mathrm{Hn} * \mathrm{Hn}^{\prime}=n *\) eye ( \(n\) ).
- The rows of Hn are orthogonal.
- \(\operatorname{det}(A)<=\operatorname{abs}(\operatorname{det}(H n))\) for all \(A\) with abs \((A(i, j))<=1\).
- Multiplying any row or column by -1 and the matrix will remain a Hadamard matrix.

See also: [compan], page 798, [hankel], page 542, [toeplitz], page 544.
hankel (c)
hankel (c, r)
Return the Hankel matrix constructed from the first column \(c\), and (optionally) the last row \(r\).
If the last element of \(c\) is not the same as the first element of \(r\), the last element of \(c\) is used. If the second argument is omitted, it is assumed to be a vector of zeros with the same size as \(c\).

A Hankel matrix formed from an m-vector \(c\), and an \(n\)-vector \(r\), has the elements
\[
H(i, j)= \begin{cases}c_{i+j-1}, & i+j-1 \leq m \\ r_{i+j-m}, & \text { otherwise }\end{cases}
\]

See also: [hadamard], page 542, [toeplitz], page 544.

\section*{hilb (n)}

Return the Hilbert matrix of order \(n\).
The \(i, j\) element of a Hilbert matrix is defined as
\[
H(i, j)=\frac{1}{(i+j-1)}
\]

Hilbert matrices are close to being singular which make them difficult to invert with numerical routines. Comparing the condition number of a random matrix \(5 \times 5\) matrix with that of a Hilbert matrix of order 5 reveals just how difficult the problem is.
```

cond (rand (5))
=> 14.392
cond (hilb (5))
\# 4.7661e+05

```

See also: [invhilb], page 543 .
invhilb (n)
Return the inverse of the Hilbert matrix of order \(n\).
This can be computed exactly using
\[
\begin{aligned}
A_{i j} & =-1^{i+j}(i+j-1)\binom{n+i-1}{n-j}\binom{n+j-1}{n-i}\binom{i+j-2}{i-2}^{2} \\
& =\frac{p(i) p(j)}{(i+j-1)}
\end{aligned}
\]
where
\[
p(k)=-1^{k}\binom{k+n-1}{k-1}\binom{n}{k}
\]

The validity of this formula can easily be checked by expanding the binomial coefficients in both formulas as factorials. It can be derived more directly via the theory of Cauchy matrices. See J. W. Demmel, Applied Numerical Linear Algebra, p. 92.
Compare this with the numerical calculation of inv (hilb (n)), which suffers from the ill-conditioning of the Hilbert matrix, and the finite precision of your computer's floating point arithmetic.
See also: [hilb], page 543.
magic ( \(n\) )
Create an \(n\)-by- \(n\) magic square.
A magic square is an arrangement of the integers \(1: \mathrm{n}^{\wedge} 2\) such that the row sums, column sums, and diagonal sums are all equal to the same value.
Note: \(n\) must be a scalar greater than or equal to 3 . If you supply \(n\) less than 3 , magic returns either a nonmagic square, or else the degenerate magic squares 1 and [].
pascal (n)
pascal ( \(n, t\) )
Return the Pascal matrix of order \(n\) if \(t=0\).
The default value of \(t\) is 0 .
When \(t=1\), return the pseudo-lower triangular Cholesky factor of the Pascal matrix (The sign of some columns may be negative). This matrix is its own inverse, that is pascal ( \(n, 1\) ) ~ \(2==\) eye ( \(n\) ).
If \(t=-1\), return the true Cholesky factor with strictly positive values on the diagonal.
If \(t=2\), return a transposed and permuted version of pascal \((n, 1)\), which is the cube root of the identity matrix. That is, pascal (n, 2) ~ \(3==\) eye ( \(n\) ).
See also: [chol], page 595.
rosser ()
Return the Rosser matrix.
This is a difficult test case used to evaluate eigenvalue algorithms.
See also: [wilkinson], page 545, [eig], page 587.
toeplitz (c)
toeplitz (c, r)
Return the Toeplitz matrix constructed from the first column \(c\), and optionally the first row \(r\).
If the second argument is omitted, the first row is taken to be the same as the first column. If the first element of \(r\) is not the same as the first element of \(c\), the first element of \(c\) is used.
A Toeplitz, or diagonal-constant, matrix has the same value along each diagonal. Although it need not be square, it often is. An MxN Toeplitz matrix has the form:
\[
\left[\begin{array}{ccccc}
c_{1} & r_{2} & r_{3} & \ldots & r_{n} \\
c_{2} & c_{1} & r_{2} & \ldots & r_{n-1} \\
c_{3} & c_{2} & c_{1} & \ldots & r_{n-2} \\
\vdots & \vdots & \vdots & \ddots & \vdots \\
c_{m} & c_{m-1} & c_{m-2} & \ldots & c m-n+1
\end{array}\right]
\]

See also: [hankel], page 542.
vander (c)
vander ( \(c, n\) )
Return the Vandermonde matrix whose next to last column is \(c\).
If \(n\) is specified, it determines the number of columns; otherwise, \(n\) is taken to be equal to the length of \(c\).
A Vandermonde matrix has the form:
\[
\left[\begin{array}{ccccc}
c_{1}^{n-1} & \cdots & c_{1}^{2} & c_{1} & 1 \\
c_{2}^{n-1} & \cdots & c_{2}^{2} & c_{2} & 1 \\
\vdots & \ddots & \vdots & \vdots & \vdots \\
c_{n}^{n-1} & \cdots & c_{n}^{2} & c_{n} & 1
\end{array}\right]
\]

See also: [polyfit], page 803.

\section*{wilkinson (n)}

Return the Wilkinson matrix of order \(n\).
Wilkinson matrices are symmetric and tridiagonal with pairs of nearly, but not exactly, equal eigenvalues. They are useful in testing the behavior and performance of eigenvalue solvers.
See also: [rosser], page 544, [eig], page 587.

\section*{17 Arithmetic}

Unless otherwise noted, all of the functions described in this chapter will work for real and complex scalar, vector, or matrix arguments. Functions described as mapping functions apply the given operation individually to each element when given a matrix argument. For example:
```

sin ([1, 2; 3, 4])
\# 0.84147 0.90930
0.14112 -0.75680

```

\subsection*{17.1 Exponents and Logarithms}
\(\exp (x)\)
Compute \(e^{x}\) for each element of x .
To compute the matrix exponential, see Chapter 18 [Linear Algebra], page 585.
See also: [log], page 547.
expm1 (x)
Compute \(e^{x}-1\) accurately in the neighborhood of zero.
See also: [exp], page 547.
\(\log (x)\)
Compute the natural logarithm, \(\ln (x)\), for each element of \(x\).
To compute the matrix logarithm, see Chapter 18 [Linear Algebra], page 585.
See also: [exp], page 547, [log1p], page 547, [log2], page 547, [ \(\log 10]\), page 547, [logspace], page 530.
reallog (x)
Return the real-valued natural logarithm of each element of \(x\).
If any element results in a complex return value reallog aborts and issues an error.
See also: [log], page 547, [realpow], page 548, [realsqrt], page 548.
\(\log 1 \mathrm{p}(\mathrm{x})\)
Compute \(\ln (1+x)\) accurately in the neighborhood of zero.
See also: [log], page 547, [exp], page 547, [expm1], page 547.
\(\log 10(x)\)
Compute the base-10 logarithm of each element of \(x\).
See also: [log], page 547, [log2], page 547, [logspace], page 530, [exp], page 547.
\(y=\log 2(x)\)
\([f, e]=\log 2(x)\)
Compute the base-2 logarithm of each element of \(x\).
If called with one output, compute the base- 2 logarithm such that \(2^{y}=x\).
If called with two output arguments, split \(x\) into binary mantissa \((f)\) and exponent (e) such that \(x=f \cdot 2^{e}\) where \(\frac{1}{2} \leq|f|<1\) and \(e\) is an integer. If \(x=0, f=e=0\).

See also: [pow2], page 548, [log], page 547, [log10], page 547, [exp], page 547.
pow2 (x)
pow2 (f, e)
With one input argument, compute \(2^{x}\) for each element of \(x\).
With two input arguments, return \(f \cdot 2^{e}\).
See also: [log2], page 547, [nextpow2], page 548, [power], page 163.
n = nextpow2 (x)
Compute the exponent for the smallest power of two larger than the input.
For each element in the input array \(x\), return the first integer \(n\) such that \(2^{n} \geq|x|\).
See also: [pow2], page 548, [log2], page 547.
realpow (x, y)
Compute the real-valued, element-by-element power operator.
This is equivalent to \(x .^{\wedge} y\), except that realpow reports an error if any return value is complex.
See also: [power], page 163, [reallog], page 547, [realsqrt], page 548.
sqrt (x)
Compute the square root of each element of \(x\).
If \(x\) is negative, a complex result is returned.
To compute the matrix square root, see Chapter 18 [Linear Algebra], page 585.
See also: [realsqrt], page 548, [nthroot], page 548.
realsqrt (x)
Return the real-valued square root of each element of \(x\).
If any element results in a complex return value realsqrt aborts and issues an error.
See also: [sqrt], page 548, [realpow], page 548, [reallog], page 547.
cbrt (x)
Compute the real-valued cube root of each element of \(x\).
Unlike \(x^{\wedge}(1 / 3)\), the result will be negative if \(x\) is negative.
If any element of \(x\) is complex, cbrt aborts with an error.
See also: [nthroot], page 548.
nthroot ( \(x, n\) )
Compute the real (non-complex) \(n\)-th root of \(x\).
\(x\) must have all real entries and \(n\) must be a scalar. If \(n\) is an even integer and \(x\) has negative entries then nthroot aborts and issues an error.
Example:
```

nthroot (-1, 3)

# -1

(-1) ~ (1 / 3)
=> 0.50000 - 0.86603i

```

See also: [realsqrt], page 548, [sqrt], page 548, [cbrt], page 548.

\subsection*{17.2 Complex Arithmetic}

In the descriptions of the following functions, \(z\) is the complex number \(x+i y\), where \(i\) is defined as \(\sqrt{-1}\).
abs (z)
Compute the magnitude of \(z\).
The magnitude is defined as \(|z|=\sqrt{x^{2}+y^{2}}\).
For example:
```

abs (3 + 4i)
F 5

```

See also: [arg], page 549 .
\(\arg (z)\)
angle (z)
Compute the argument, i.e., angle of \(z\).
This is defined as, \(\theta=\operatorname{atan} 2(y, x)\), in radians.
For example:
```

arg (3 + 4i)
\# 0.92730

```

See also: [abs], page 549 .
conj (z)
Return the complex conjugate of \(z\).
The complex conjugate is defined as \(\bar{z}=x-i y\).
See also: [real], page 550, [imag], page 550.
cplxpair (z)
cplxpair (z, tol)
cplxpair (z, tol, dim)
Sort the numbers \(z\) into complex conjugate pairs ordered by increasing real part.
The negative imaginary complex numbers are placed first within each pair. All real numbers (those with abs (imag (z)) / abs (z) < tol) are placed after the complex pairs.
tol is a weighting factor in the range \([0,1)\) which determines the tolerance of the matching. The default value is \(100 *\) eps and the resulting tolerance for a given complex pair is tol \(*\) abs \((z(i))\) ).
By default the complex pairs are sorted along the first non-singleton dimension of \(z\). If \(\operatorname{dim}\) is specified, then the complex pairs are sorted along this dimension.
Signal an error if some complex numbers could not be paired. Signal an error if all complex numbers are not exact conjugates (to within tol). Note that there is no defined order for pairs with identical real parts but differing imaginary parts.
```

cplxpair (exp (2i*pi*[0:4]'/5)) == exp (2i*pi*[3; 2; 4; 1; 0]/5)

```
imag ( \(z\) )
Return the imaginary part of \(z\) as a real number.
See also: [real], page 550, [conj], page 549.
real ( \(z\) )
Return the real part of \(z\).
See also: [imag], page 550, [conj], page 549.

\subsection*{17.3 Trigonometry}

Octave provides the following trigonometric functions where angles are specified in radians. To convert from degrees to radians multiply by \(\pi / 180\) or use the deg2rad function. For example, \(\sin (30 * \mathrm{pi} / 180)\) returns the sine of 30 degrees. As an alternative, Octave provides a number of trigonometric functions which work directly on an argument specified in degrees. These functions are named after the base trigonometric function with a ' \(d\) ' suffix. As an example, sin expects an angle in radians while sind expects an angle in degrees.

Octave uses the C library trigonometric functions. It is expected that these functions are defined by the ISO/IEC 9899 Standard. This Standard is available at: http:// www.open-std.org/jtc1/sc22/wg14/www/docs/n1124.pdf. Section F.9.1 deals with the trigonometric functions. The behavior of most of the functions is relatively straightforward. However, there are some exceptions to the standard behavior. Many of the exceptions involve the behavior for -0 . The most complex case is atan2. Octave exactly implements the behavior given in the Standard. Including atan2 \(( \pm 0,-0)\) returns \(\pm \pi\).

It should be noted that matlab uses different definitions which apparently do not distinguish -0.
rad \(=\operatorname{deg} 2 r a d ~(d e g)\)
Convert degrees to radians.
The input deg must be a scalar, vector, or N-dimensional array of double or single floating point values. deg may be complex in which case the real and imaginary components are converted separately.
The output rad is the same size and shape as deg with degrees converted to radians using the conversion constant pi/180.
Example:
```

deg2rad ([0, 90, 180, 270, 360])
\#

```

See also: \([\mathrm{rad} 2 \mathrm{deg}]\), page 550 .
```

deg = rad2deg (rad)

```

Convert radians to degrees.
The input rad must be a scalar, vector, or N-dimensional array of double or single floating point values. rad may be complex in which case the real and imaginary components are converted separately.
The output deg is the same size and shape as rad with radians converted to degrees using the conversion constant 180/pi.

Example:
```

rad2deg ([0, pi/2, pi, 3/2*pi, 2*pi])
\#

```

See also: [deg2rad], page 550.
\(\sin (x)\)
Compute the sine for each element of \(x\) in radians.
See also: [asin], page 551, [sind], page 553, [sinh], page 552.
\(\cos (x)\)
Compute the cosine for each element of \(x\) in radians.
See also: [acos], page 551, [cosd], page 553, [cosh], page 552.
\(\tan (z)\)
Compute the tangent for each element of \(x\) in radians.
See also: [atan], page 551, [tand], page 553, [tanh], page 552.
\(\sec (x)\)
Compute the secant for each element of \(x\) in radians.
See also: [asec], page 551, [secd], page 553, [sech], page 552.
\(\csc (x)\)
Compute the cosecant for each element of \(x\) in radians.
See also: [acsc], page 552, [cscd], page 554, [csch], page 552.
\(\cot (x)\)
Compute the cotangent for each element of \(x\) in radians.
See also: [acot], page 552, [cotd], page 554, [coth], page 552.
\(\operatorname{asin}(x)\)
Compute the inverse sine in radians for each element of \(x\).
See also: [sin], page 551, [asind], page 554.
\(\operatorname{acos}(x)\)
Compute the inverse cosine in radians for each element of \(x\).
See also: [cos], page 551, [acosd], page 554.
\(\operatorname{atan}(x)\)
Compute the inverse tangent in radians for each element of \(x\).
See also: [tan], page 551, [atand], page 554.
\(\operatorname{asec}(x)\)
Compute the inverse secant in radians for each element of \(x\).
See also: [sec], page 551, [asecd], page 554.
\(\operatorname{acsc}(x)\)
Compute the inverse cosecant in radians for each element of \(x\).
See also: [csc], page 551, [acscd], page 554.
\(\operatorname{acot}(x)\)
Compute the inverse cotangent in radians for each element of \(x\).
See also: [cot], page 551, [acotd], page 554 .
\(\sinh (x)\)
Compute the hyperbolic sine for each element of \(x\).
See also: [asinh], page 552, [cosh], page 552, [tanh], page 552.
\(\cosh (x)\)
Compute the hyperbolic cosine for each element of \(x\).
See also: [acosh], page 552, [sinh], page 552, [tanh], page 552.
\(\tanh (x)\)
Compute hyperbolic tangent for each element of \(x\).
See also: [atanh], page 552 , [sinh], page 552 , [cosh], page 552.
\(\operatorname{sech}(x)\)
Compute the hyperbolic secant of each element of \(x\).
See also: [asech], page 552.
\(\operatorname{csch}(x)\)
Compute the hyperbolic cosecant of each element of \(x\).
See also: [acsch], page 553 .
\(\operatorname{coth}(x)\)
Compute the hyperbolic cotangent of each element of \(x\).
See also: [acoth], page 553 .
\(\operatorname{asinh}(x)\)
Compute the inverse hyperbolic sine for each element of \(x\).
See also: [sinh], page 552 .
\(\operatorname{acosh}(x)\)
Compute the inverse hyperbolic cosine for each element of \(x\).
See also: [cosh], page 552.
\(\operatorname{atanh}(x)\)
Compute the inverse hyperbolic tangent for each element of \(x\).
See also: [tanh], page 552.
\(\operatorname{asech}(\mathrm{x})\)
Compute the inverse hyperbolic secant of each element of \(x\).
See also: [sech], page 552 .

\section*{\(\operatorname{acsch}(x)\)}

Compute the inverse hyperbolic cosecant of each element of \(x\).
See also: [csch], page 552.
\(\operatorname{acoth}(x)\)
Compute the inverse hyperbolic cotangent of each element of \(x\).
See also: [coth], page 552.
\(\operatorname{atan} 2(y, x)\)
Compute atan \((y / x)\) for corresponding elements of \(y\) and \(x\).
\(y\) and \(x\) must match in size and orientation. The signs of elements of \(y\) and \(x\) are used to determine the quadrants of each resulting value.
This function is equivalent to arg (complex ( \(x, y\) )).
See also: [tan], page 551, [tand], page 553, [tanh], page 552, [atanh], page 552.
Octave provides the following trigonometric functions where angles are specified in degrees. These functions produce true zeros at the appropriate intervals rather than the small round-off error that occurs when using radians. For example:
```

cosd (90)

```
    \(\Rightarrow 0\)
\(\cos (\mathrm{pi} / 2)\)
    \(\Rightarrow 6.1230 \mathrm{e}-17\)
sind ( \(x\) )
Compute the sine for each element of \(x\) in degrees.
The function is more accurate than \(\sin\) for large values of \(x\) and for multiples of 180 degrees ( \(x / 180\) is an integer) where sind returns 0 rather than a small value on the order of eps.
See also: [asind], page 554, [sin], page 551.
cosd (x)
Compute the cosine for each element of \(x\) in degrees.
The function is more accurate than cos for large values of \(x\) and for multiples of 90 degrees ( \(x=90+180 * n\) with \(n\) an integer) where cosd returns 0 rather than a small value on the order of eps.
See also: [acosd], page 554, [cos], page 551.
tand ( \(x\) )
Compute the tangent for each element of \(x\) in degrees.
Returns zero for elements where \(x / 180\) is an integer and \(\operatorname{Inf}\) for elements where \((x-90) / 180\) is an integer.
See also: [atand], page 554, [tan], page 551.
\(\sec (x)\)
Compute the secant for each element of \(x\) in degrees.
See also: [asecd], page 554, [sec], page 551.
\(\csc (x)\)
Compute the cosecant for each element of \(x\) in degrees.
See also: [acscd], page 554, [csc], page 551.
cotd (x)
Compute the cotangent for each element of \(x\) in degrees.
See also: [acotd], page 554, [cot], page 551.
asind (x)
Compute the inverse sine in degrees for each element of \(x\).
See also: [sind], page 553, [asin], page 551.
\(\operatorname{acosd}(x)\)
Compute the inverse cosine in degrees for each element of \(x\).
See also: [cosd], page 553, [acos], page 551.
atand (x)
Compute the inverse tangent in degrees for each element of \(x\).
See also: [tand], page 553, [atan], page 551.
\(\operatorname{atan} 2 \mathrm{~d}(y, x)\)
Compute atan \((y / x)\) in degrees for corresponding elements from \(y\) and \(x\).
See also: [tand], page 553, [atan2], page 553.
\(\operatorname{asec}(x)\)
Compute the inverse secant in degrees for each element of \(x\).
See also: [secd], page 553, [asec], page 551.
\(\operatorname{acscd}(x)\)
Compute the inverse cosecant in degrees for each element of \(x\).
See also: [cscd], page 554, [acsc], page 552.
\(\operatorname{acotd}(x)\)
Compute the inverse cotangent in degrees for each element of \(x\).
See also: [cotd], page 554, [acot], page 552.
Finally, there are two trigonometric functions that calculate special arguments with increased accuracy.
\(y=\operatorname{sinp} i(x)\)
Compute sine ( \(x^{*}\) pi) for each element of \(x\) accurately.
The ordinary sin function uses IEEE floating point numbers and may produce results that are very close (within a few eps) of the correct value, but which are not exact. The sinpi function is more accurate and returns 0 exactly for integer values of \(x\) and \(+1 /-1\) for half-integer values (e.g., ..., \(-3 / 2,-1 / 2,1 / 2,3 / 2, \ldots\).

Example
comparison of \(\sin\) and sinpi for integer values of \(x\)
```

$\sin ([0,1,2,3] * p i)$
$\Rightarrow$
$0 \quad 1.2246 \mathrm{e}-16 \quad-2.4493 \mathrm{e}-16 \quad 3.6739 \mathrm{e}-16$
sinpi ([0, 1, 2, 3])
$\Rightarrow$
0000

```

See also: [cospi], page 555, [sin], page 551.
\(y=\operatorname{cospi}(x)\)
Compute cosine ( \(x *\) pi) for each element of \(x\) accurately.
The ordinary cos function uses IEEE floating point numbers and may produce results that are very close (within a few eps) of the correct value, but which are not exact. The cospi function is more accurate and returns 0 exactly for half-integer values of x (e.g., \(\ldots,-3 / 2,-1 / 2,1 / 2,3 / 2, \ldots\) ), and \(+1 /-1\) for integer values.
Example
comparison of cos and cospi for half-integer values of \(x\)
```

cos ([-3/2, -1/2, 1/2, 3/2] * pi)

# 

        -1.8370e-16 6.1232e-17 6.1232e-17 -1.8370e-16
    cospi ([-3/2, -1/2, 1/2, 3/2])

# 

        0 0 0 0
    ```

See also: [sinpi], page 554, [cos], page 551.

\subsection*{17.4 Sums and Products}
```

sum (x)
sum (x, dim)
sum (..., "native")
sum (..., "double")
sum (..., "extra")

```

Sum of elements along dimension dim.
If \(\operatorname{dim}\) is omitted, it defaults to the first non-singleton dimension.
The optional "type" input determines the class of the variable used for calculations. By default, operations on floating point inputs (double or single) are performed in their native data type, while operations on integer, logical, and character data types are performed using doubles. If the argument "native" is given, then the operation is performed in the same type as the original argument.
For example:
```

sum ([true, true])
=>2
sum ([true, true], "native")
=> true

```

If "double" is given the sum is performed in double precision even for single precision inputs.

For double precision inputs, the "extra" option will use a more accurate algorithm than straightforward summation. For single precision inputs, "extra" is the same as "double". For all other data type "extra" has no effect.

See also: [cumsum], page 556, [sumsq], page 557, [prod], page 556.
```

prod (x)
prod (x, dim)
prod (..., "native")
prod (..., "double")
Product of elements along dimension dim. as the original argument, rather than the default double type.
For example:

```
```

prod ([true, true])

```
prod ([true, true])
    =>1
    =>1
prod ([true, true], "native")
prod ([true, true], "native")
    # true
```

    # true
    ```

If dim is omitted, it defaults to the first non-singleton dimension.
The optional "type" input determines the class of the variable used for calculations. If the argument "native" is given, then the operation is performed in the same type

On the contrary, if "double" is given, the operation is performed in double precision even for single precision inputs.
See also: [cumprod], page 557, [sum], page 555.
```

cumsum (x)
cumsum (x, dim)
cumsum (..., "native")
cumsum (..., "double")

```

Cumulative sum of elements along dimension dim.
If dim is omitted, it defaults to the first non-singleton dimension. For example:
```

cumsum ([1, 2; 3, 4; 5, 6])

# 1 2

    4 6
    9 12
    ```

For an explanation of the optional parameters "native" and "double", see [sum], page 555.
See also: [sum], page 555, [cumprod], page 557.
cumprod ( \(x\) )
cumprod (x, dim)
Cumulative product of elements along dimension dim.
If dim is omitted, it defaults to the first non-singleton dimension. For example:
```

cumprod ([1, 2; 3, 4; 5, 6])

# 1 2

                3 8
                1548
    ```

See also: [prod], page 556, [cumsum], page 556.
sumsq ( \(x\) )
sumsq ( \(x, \operatorname{dim}\) )
Sum of squares of elements along dimension dim.
If dim is omitted, it defaults to the first non-singleton dimension.
This function is conceptually equivalent to computing
```

sum (x .* conj (x), dim)

```
but it uses less memory and avoids calling conj if x is real.
See also: [sum], page 555 , [prod], page 556 .

\subsection*{17.5 Utility Functions}
ceil (x)
Return the smallest integer not less than \(x\).
This is equivalent to rounding towards positive infinity.
If \(x\) is complex, return ceil (real ( \(x\) ) ) + ceil (imag ( \(x\) )) * I.
```

ceil ([-2.7, 2.7])
=> -2 3

```

See also: [floor], page 557, [round], page 558, [fix], page 557.
fix (x)
Truncate fractional portion of \(x\) and return the integer portion.
This is equivalent to rounding towards zero. If \(x\) is complex, return fix (real ( \(x\) ) )
+ fix (imag (x)) * I.
fix ([-2.7, 2.7])
\(\Rightarrow-2 \quad 2\)
See also: [ceil], page 557, [floor], page 557, [round], page 558.
floor (x)
Return the largest integer not greater than \(x\).
This is equivalent to rounding towards negative infinity. If \(x\) is complex, return floor (real (x)) + floor (imag (x)) * I.
floor ([-2.7, 2.7])
\(\Rightarrow-3 \quad 2\)
See also: [ceil], page 557, [round], page 558, [fix], page 557.
round (x)
Return the integer nearest to \(x\).
If \(x\) is complex, return round (real (x)) + round (imag (x)) * I. If there are two nearest integers, return the one further away from zero.
\[
\begin{aligned}
\text { round } & ([-2.7, \\
\Rightarrow-3 & 2.7]) \\
\Rightarrow & 3
\end{aligned}
\]

See also: [ceil], page 557, [floor], page 557, [fix], page 557, [roundb], page 558.
roundb (x)
Return the integer nearest to \(x\). If there are two nearest integers, return the even one (banker's rounding).
If \(x\) is complex, return roundb (real (x)) + roundb (imag (x)) * I.
See also: [round], page 558.
\(\max (x)\)
\(\max (x,[], \operatorname{dim})\)
[ \(\mathrm{w}, \mathrm{iw}\) ] \(=\max (\mathrm{x})\)
\(\max (x, y)\)
Find maximum values in the array \(x\).
For a vector argument, return the maximum value. For a matrix argument, return a row vector with the maximum value of each column. For a multi-dimensional array, max operates along the first non-singleton dimension.
If the optional third argument dim is present then operate along this dimension. In this case the second argument is ignored and should be set to the empty matrix.
For two inputs ( \(x\) and \(y\) ), return the pairwise maximum according to the rules for Section 19.2 [Broadcasting], page 627.
Thus,
\(\max (\max (x))\)
returns the largest element of the 2-D matrix \(x\), and
```

max (2:5, pi)
\# 3.1416 3.1416 4.0000 5.0000

```
compares each element of the range \(2: 5\) with pi, and returns a row vector of the maximum values.
For complex arguments, the magnitude of the elements are used for comparison. If the magnitudes are identical, then the results are ordered by phase angle in the range (-pi, pi]. Hence,
```

max ([[-1 i 1 1 -i])
\# -1

```
because all entries have magnitude 1, but -1 has the largest phase angle with value pi.
If called with one input and two output arguments, max also returns the first index of the maximum value(s). Thus,
```

[x, ix] = max ([1, 3, 5, 2, 5])
m x = 5
ix = 3

```

See also: [min], page 559, [cummax], page 559, [cummin], page 560.
```

min (x)

```
\(\min (x,[], \operatorname{dim})\)
[ \(\mathrm{w}, \mathrm{i} \mathrm{w}]=\min (\mathrm{x})\)
\(\min (x, y)\)

Find minimum values in the array \(x\).
For a vector argument, return the minimum value. For a matrix argument, return a row vector with the minimum value of each column. For a multi-dimensional array, min operates along the first non-singleton dimension.
If the optional third argument dim is present then operate along this dimension. In this case the second argument is ignored and should be set to the empty matrix.

For two inputs ( \(x\) and \(y\) ), return the pairwise minimum according to the rules for Section 19.2 [Broadcasting], page 627.

Thus,
```

min (min (x))

```
returns the smallest element of the 2-D matrix \(x\), and
```

min (2:5, pi)
\#}2.0000\quad3.0000 3.1416 3.1416

```
compares each element of the range \(2: 5\) with pi, and returns a row vector of the minimum values.
For complex arguments, the magnitude of the elements are used for comparison. If the magnitudes are identical, then the results are ordered by phase angle in the range (-pi, pi]. Hence,
```

min ([-1 i 1 1 -i])
\# -i

```
because all entries have magnitude 1, but -i has the smallest phase angle with value -pi/2.
If called with one input and two output arguments, min also returns the first index of the minimum value(s). Thus,
\[
\begin{aligned}
& {[x,}i x] \\
& \Rightarrow \quad \min ([1,3,0,2,0]) \\
& \quad x=0 \\
& i x=3
\end{aligned}
\]

See also: [max], page 558, [cummin], page 560, [cummax], page 559.
cummax (x)
cummax (x, dim)
\([\mathrm{w}, \mathrm{iw}]=\) cummax (...)
Return the cumulative maximum values along dimension dim.
If \(\operatorname{dim}\) is unspecified it defaults to column-wise operation. For example:
```

cummax ([[1 [10 2 % 6
m}\begin{array}{lllllll}{1}\&{3}\&{3}\&{6}\&{6}\&{6}

```

If called with two output arguments the index of the maximum value is also returned.
```

[w, iw] = cummax ([[1 [13 2 6 4 5])

# 

w = 1 1 3 3 6 6 6 6
iw = 1 2 2 2 4 4 4 4

```

See also: [cummin], page 560, [max], page 558, [min], page 559.
cummin ( \(x\) )
cummin ( \(x\), dim)
\([\mathrm{w}, \mathrm{i} \mathrm{w}]=\operatorname{cummin}(\mathrm{x})\)
Return the cumulative minimum values along dimension dim.
If \(\operatorname{dim}\) is unspecified it defaults to column-wise operation. For example:
```

cummin ([[5 4 4 6 2 3 1])
=>

```

If called with two output arguments the index of the minimum value is also returned.
```

[w, iw] = cummin ([[5 4 6 2 3 1])

# 

w = 5 5 4 4 4 2 2 2 1
iw = 1 1 2 2 2 4 4 4

```

See also: [cummax], page 559, [min], page 559, [max], page 558.
hypot ( \(x, y\) )
hypot ( \(x, y, z, \ldots\) )
Compute the element-by-element square root of the sum of the squares of \(x\) and \(y\).
This is equivalent to sqrt ( \(x . \wedge 2+y . \wedge 2\) ), but is calculated in a manner that avoids overflows for large values of \(x\) or \(y\).
hypot can also be called with more than 2 arguments; in this case, the arguments are accumulated from left to right:
```

hypot (hypot (x, y), z)
hypot (hypot (hypot (x, y), z), w), etc.

```
\(d x=\) gradient ( \(m\) )
\([d x, d y, d z, \ldots]=\) gradient (m)
[...] = gradient (m, s)
[...] = gradient (m, x, y, z, ...)
[...] = gradient (f, x0)
[...] = gradient (f, x0, s)
[...] = gradient (f, \(x 0, x, y, \ldots)\)

Calculate the gradient of sampled data or a function.
If \(m\) is a vector, calculate the one-dimensional gradient of \(m\). If \(m\) is a matrix the gradient is calculated for each dimension.
[dx, dy] = gradient \((m)\) calculates the one-dimensional gradient for \(x\) and \(y\) direction if \(m\) is a matrix. Additional return arguments can be use for multi-dimensional matrices.

A constant spacing between two points can be provided by the \(s\) parameter. If \(s\) is a scalar, it is assumed to be the spacing for all dimensions. Otherwise, separate values of the spacing can be supplied by the \(x, \ldots\) arguments. Scalar values specify an equidistant spacing. Vector values for the \(x, \ldots\) arguments specify the coordinate for that dimension. The length must match their respective dimension of \(m\).
At boundary points a linear extrapolation is applied. Interior points are calculated with the first approximation of the numerical gradient
\[
y^{\prime}(i)=1 /(x(i+1)-x(i-1)) *(y(i-1)-y(i+1)) .
\]

If the first argument \(f\) is a function handle, the gradient of the function at the points in \(x 0\) is approximated using central difference. For example, gradient (@cos, 0) approximates the gradient of the cosine function in the point \(x 0=0\). As with sampled data, the spacing values between the points from which the gradient is estimated can be set via the \(s\) or \(d x, d y, \ldots\) arguments. By default a spacing of 1 is used.
See also: [diff], page 512, [del2], page 562.
\(\operatorname{dot}(x, y, \operatorname{dim})\)
Compute the dot product of two vectors.
If \(x\) and \(y\) are matrices, calculate the dot products along the first non-singleton dimension.
If the optional argument dim is given, calculate the dot products along this dimension. Implementation Note: This is equivalent to sum (conj (X) .* Y, dim), but avoids forming a temporary array and is faster. When \(X\) and \(Y\) are column vectors, the result is equivalent to \(X^{\prime} * Y\). Although, dot is defined for integer arrays, the output may differ from the expected result due to the limited range of integer objects.

See also: [cross], page 561, [divergence], page 561.
cross ( \(x, y\) )
cross ( \(x, y, d i m\) )
Compute the vector cross product of two 3-dimensional vectors \(x\) and \(y\).
If \(x\) and \(y\) are matrices, the cross product is applied along the first dimension with three elements.
The optional argument dim forces the cross product to be calculated along the specified dimension.
Example Code:
```

cross ([1, 1, 0], [0, 1, 1])
\#
1 -1 1

```

See also: [dot], page 561, [curl], page 562, [divergence], page 561.
```

div = divergence (x, y, z, fx, fy,fz)

```
div \(=\) divergence ( \(f x, f y, f z\) )
div \(=\) divergence ( \(x, y, f x, f y\) )
div \(=\) divergence ( \(f x, f y\) )

Calculate divergence of a vector field given by the arrays \(f x\), \(f y\), and \(f z\) or \(f x\), fy respectively.
\[
\operatorname{div} F(x, y, z)=\partial_{x} F+\partial_{y} F+\partial_{z} F
\]

The coordinates of the vector field can be given by the arguments \(x, y, z\) or \(x, y\) respectively.
See also: [curl], page 562, [gradient], page 560, [del2], page 562, [dot], page 561.
[cx, cy, cz, v] = curl (x, y, z, fx, fy, fz)
[cz, v] = curl (x, y,fx, fy)
[...] = curl (fx,fy,fz)
[...] = curl (fx, fy)
\(\mathrm{v}=\operatorname{curl}(. .\).
Calculate curl of vector field given by the arrays \(f x\), fy, and \(f z\) or \(f x\), fy respectively.
\[
\operatorname{curlF}(x, y, z)=\left(\frac{\partial F_{z}}{\partial y}-\frac{\partial F_{y}}{\partial z}, \frac{\partial F_{x}}{\partial z}-\frac{\partial F_{z}}{\partial x}, \frac{\partial F_{y}}{\partial x}-\frac{\partial F_{x}}{\partial y}\right)
\]

The coordinates of the vector field can be given by the arguments \(x, y, z\) or \(x, y\) respectively. \(v\) calculates the scalar component of the angular velocity vector in direction of the z-axis for two-dimensional input. For three-dimensional input the scalar rotation is calculated at each grid point in direction of the vector field at that point.
See also: [divergence], page 561, [gradient], page 560, [del2], page 562, [cross], page 561.
\(L=\operatorname{del} 2(M)\)
\(L=\operatorname{del} 2(M, h)\)
\(L=\operatorname{del} 2(M, d x, d y, \ldots)\)
Calculate the discrete Laplace operator \(\left(\nabla^{2}\right)\).
For a 2 -dimensional matrix \(M\) this is defined as
\[
L=\frac{1}{4}\left(\frac{d^{2}}{d x^{2}} M(x, y)+\frac{d^{2}}{d y^{2}} M(x, y)\right)
\]

For N-dimensional arrays the sum in parentheses is expanded to include second derivatives over the additional higher dimensions.
The spacing between evaluation points may be defined by \(h\), which is a scalar defining the equidistant spacing in all dimensions. Alternatively, the spacing in each dimension may be defined separately by \(d x\), \(d y\), etc. A scalar spacing argument defines equidistant spacing, whereas a vector argument can be used to specify variable spacing. The length of the spacing vectors must match the respective dimension of \(M\). The default spacing value is 1 .
Dimensions with fewer than 3 data points are skipped. Boundary points are calculated from the linear extrapolation of interior points.
Example: Second derivative of \(2^{*} x^{\wedge} 3\)
```

f = @(x) 2*x.^3;
dd = @(x) 12*x;
x = 1:6;
L = 4*del2 (f(x));
assert (L, dd (x));

```

See also: [gradient], page 560, [diff], page 512.
factorial ( \(n\) )
Return the factorial of \(n\) where \(n\) is a real non-negative integer.
If \(n\) is a scalar, this is equivalent to prod \((1: n)\). For vector or matrix arguments, return the factorial of each element in the array.
For non-integers see the generalized factorial function gamma. Note that the factorial function grows large quite quickly, and even with double precision values overflow will occur if \(n>171\). For such cases consider gammaln.
See also: [prod], page 556, [gamma], page 574, [gammaln], page 576.
pf = factor (q)
\([p f, n]=\) factor (q)
Return the prime factorization of \(q\).
The prime factorization is defined as prod \((p f)==q\) where every element of \(p f\) is a prime number. If \(q==1\), return 1 . The output \(p f\) is of the same numeric class as the input.
With two output arguments, return the unique prime factors \(p f\) and their multiplicities. That is, \(\operatorname{prod}\left(p f .{ }^{\wedge}\right)==q\).
Implementation Note: The input \(q\) must be less than flintmax when the input is a floating-point class (double or single).

See also: [gcd], page 563, [lcm], page 563, [isprime], page 70, [primes], page 564.
\(g=\operatorname{gcd}(a 1, a 2, \ldots)\)
\([g, \mathrm{v} 1, \ldots]=\operatorname{gcd}(\mathrm{a} 1, \mathrm{a} 2, \ldots)\)
Compute the greatest common divisor of a1, a2,....
All arguments must be the same size or scalar. For arrays, the greatest common divisor is calculated for each element individually. All elements must be ordinary or Gaussian (complex) integers. Note that for Gaussian integers, the gcd is only unique up to a phase factor (multiplication by \(1,-1\), i , or -i ), so an arbitrary greatest common divisor among the four possible is returned.
Optional return arguments \(v 1, \ldots\), contain integer vectors such that,
\(g=v_{1} a_{1}+v_{2} a_{2}+\cdots\)
Example code:
\[
\begin{aligned}
& \operatorname{gcd}([15,9],[20,18]) \\
& \quad \Rightarrow 59
\end{aligned}
\]

See also: [lcm], page 563, [factor], page 563, [isprime], page 70.
\(1 \mathrm{~cm}(x, y)\)
lcm ( \(x, y, \ldots\) )
Compute the least common multiple of \(x\) and \(y\), or of the list of all arguments.
All elements must be numeric and of the same size or scalar.
See also: [factor], page 563, [gcd], page 563, [isprime], page 70.
rem ( \(x, y\) )
Return the remainder of the division \(x / y\).
The remainder is computed using the expression
x - y .* fix (x ./ y)

An error message is printed if the dimensions of the arguments do not agree, or if either argument is complex.
Programming Notes: When calculating with floating point numbers (double, single), values within a few eps of an integer will be rounded to that integer before computation for compatibility with matlab. Any floating point integers greater than flintmax ( \(2^{\wedge} 53\) for double) will not compute correctly. For larger integer values convert the input to uint64 before calling this function.
By convention,
```

rem (x, 0) = NaN if x is a floating point variable
rem (x, 0) = 0 if x is an integer variable
rem (x, y) returns a value with the signbit from x

```

For the opposite conventions see the mod function. In general, rem is best when computing the remainder after division of two positive numbers. For negative numbers, or when the values are periodic, mod is a better choice.
See also: [mod], page 564 .
\(\bmod (x, y)\)
Compute the modulo of \(x\) and \(y\).
Conceptually this is given by
x - y .* floor (x ./ y)
and is written such that the correct modulus is returned for integer types. This function handles negative values correctly. That is, \(\bmod (-1,3)\) is 2 , not -1 , as rem ( \(-1,3\) ) returns.
An error results if the dimensions of the arguments do not agree, or if either of the arguments is complex.
Programming Notes: When calculating with floating point numbers (double, single), values within a few eps of an integer will be rounded to that integer before computation for compatibility with matlab. Any floating point integers greater than flintmax ( \(2^{\wedge} 53\) for double) will not compute correctly. For larger integer values convert the input to uint64 before calling this function.
By convention,
```

mod (x, 0) = x
mod (x, y) returns a value with the signbit from y

```

For the opposite conventions see the rem function. In general, mod is a better choice than rem when any of the inputs are negative numbers or when the values are periodic.
See also: [rem], page 564.
```

p = primes (n)

```

Return all primes up to \(n\).

The output data class (double, single, uint32, etc.) is the same as the input class of n. The algorithm used is the Sieve of Eratosthenes.

Note: For a specific number \(n\) of primes, call list_primes ( \(n\) ). Alternatively, call primes \((n * \log (k * n))(1: n)\) where \(k\) is about 5 or 6 . This works because the distance
from one prime to the next is proportional to the logarithm of the prime, on average.
On integrating, there are about \(n\) primes less than \(n * \log (5 * n)\).
See also: [list_primes], page 565, [isprime], page 70.
```

list_primes ()
list_primes (n)

```

List the first \(n\) primes.
If \(n\) is unspecified, the first 25 primes are listed.
See also: [primes], page 564, [isprime], page 70.
sign ( \(x\) )
Compute the signum function.
This is defined as
\[
\operatorname{sign}(x)= \begin{cases}1, & x>0 \\ 0, & x=0 \\ -1, & x<0\end{cases}
\]

For complex arguments, sign returns \(\mathrm{x} . /\) abs ( x ).
Note that sign ( -0.0 ) is 0 . Although IEEE 754 floating point allows zero to be signed, 0.0 and -0.0 compare equal. If you must test whether zero is signed, use the signbit function.

See also: [signbit], page 565.
signbit (x)
Return logical true if the value of \(x\) has its sign bit set and false otherwise.
This behavior is consistent with the other logical functions. See Section 4.6 [Logical Values], page 64. The behavior differs from the C language function which returns nonzero if the sign bit is set.
This is not the same as \(\mathrm{x}<0.0\), because IEEE 754 floating point allows zero to be signed. The comparison \(-0.0<0.0\) is false, but signbit ( -0.0 ) will return a nonzero value.
See also: [sign], page 565.

\subsection*{17.6 Special Functions}
[a, ierr] = airy (k, z, opt)
Compute Airy functions of the first and second kind, and their derivatives.
\begin{tabular}{|c|c|c|}
\hline K & Function & Scale factor (if "opt" is supplied) \\
\hline 0 & Ai (Z) & \(\exp ((2 / 3) * \mathrm{Z} *\) sqrt (Z) ) \\
\hline 1 & dAi \((\mathrm{Z}) / \mathrm{dZ}\) & \(\exp ((2 / 3) * Z * \operatorname{sqrt}(\mathrm{Z})\) ) \\
\hline 2 & Bi (Z) & \(\exp (-\mathrm{abs}(\) real \(((2 / 3) * \mathrm{Z} *\) sqrt (Z))) \\
\hline 3 & dBi \((\mathrm{Z}) / \mathrm{dZ}\) & \(\exp (-\mathrm{abs}(\mathrm{real}((2 / 3) * \mathrm{Z} *\) sqrt (Z))) \\
\hline
\end{tabular}

The function call airy ( \(z\) ) is equivalent to airy ( \(0, z\) ).
The result is the same size as \(z\).
If requested, ierr contains the following status information and is the same size as the result.

0 . Normal return.
1. Input error, return NaN.
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Loss of significance by argument reduction, output may be inaccurate.
5. Error-no computation, algorithm termination condition not met, return NaN.
\(J\) = besselj (alpha, x)
\(J\) = besselj (alpha, x, opt)
[J, ierr] = besselj (...)
Compute Bessel functions of the first kind.
The order of the Bessel function alpha must be real. The points for evaluation \(x\) may be complex.
If the optional argument opt is 1 or true, the result \(J\) is multiplied by exp (-abs (imag (x))).
If alpha is a scalar, the result is the same size as \(x\). If \(x\) is a scalar, the result is the same size as alpha. If alpha is a row vector and \(x\) is a column vector, the result is a matrix with length ( \(x\) ) rows and length (alpha) columns. Otherwise, alpha and \(x\) must conform and the result will be the same size.
If requested, ierr contains the following status information and is the same size as the result.

0 . Normal return.
1. Input error, return NaN.
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Loss of significance by argument reduction, output may be inaccurate.
5. Error-no computation, algorithm termination condition not met, return NaN .

See also: [bessely], page 566, [besseli], page 567, [besselk], page 568, [besselh], page 568.
\(Y=\) bessely (alpha, \(x\) )
\(Y\) = bessely (alpha, x, opt)
[Y, ierr] = bessely (...)
Compute Bessel functions of the second kind.
The order of the Bessel function alpha must be real. The points for evaluation \(x\) may be complex.

If the optional argument opt is 1 or true, the result \(Y\) is multiplied by exp (-abs (imag (x))).
If alpha is a scalar, the result is the same size as \(x\). If \(x\) is a scalar, the result is the same size as alpha. If alpha is a row vector and \(x\) is a column vector, the result is a matrix with length ( \(x\) ) rows and length (alpha) columns. Otherwise, alpha and \(x\) must conform and the result will be the same size.
If requested, ierr contains the following status information and is the same size as the result.
0 . Normal return.
1. Input error, return NaN .
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Complete loss of significance by argument reduction, return NaN .
5. Error-no computation, algorithm termination condition not met, return NaN.

See also: [besselj], page 566, [besseli], page 567, [besselk], page 568, [besselh], page 568.
\(I=\) besseli (alpha, \(x\) )
\(I\) = besseli (alpha, x, opt)
[I, ierr] = besseli (...)
Compute modified Bessel functions of the first kind.
The order of the Bessel function alpha must be real. The points for evaluation x may be complex.
If the optional argument opt is 1 or true, the result \(I\) is multiplied by \(\exp (-\mathrm{abs}(\) real (x))).
If alpha is a scalar, the result is the same size as \(x\). If \(x\) is a scalar, the result is the same size as alpha. If alpha is a row vector and \(x\) is a column vector, the result is a matrix with length ( \(x\) ) rows and length (alpha) columns. Otherwise, alpha and \(x\) must conform and the result will be the same size.
If requested, ierr contains the following status information and is the same size as the result.
0 . Normal return.
1. Input error, return NaN .
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Complete loss of significance by argument reduction, return NaN.
5. Error-no computation, algorithm termination condition not met, return NaN.

See also: [besselk], page 568, [besselj], page 566, [bessely], page 566, [besselh], page 568.
\(K=\) besselk (alpha, \(x\) )
\(K=\) besselk (alpha, x, opt)
[K, ierr] = besselk (...)
Compute modified Bessel functions of the second kind.
The order of the Bessel function alpha must be real. The points for evaluation \(x\) may be complex.
If the optional argument opt is 1 or true, the result \(K\) is multiplied by \(\exp (x)\).
If alpha is a scalar, the result is the same size as \(x\). If \(x\) is a scalar, the result is the same size as alpha. If alpha is a row vector and \(x\) is a column vector, the result is a matrix with length ( \(x\) ) rows and length (alpha) columns. Otherwise, alpha and \(x\) must conform and the result will be the same size.
If requested, ierr contains the following status information and is the same size as the result.
0 . Normal return.
1. Input error, return NaN .
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Complete loss of significance by argument reduction, return NaN .
5. Error-no computation, algorithm termination condition not met, return NaN.

See also: [besseli], page 567, [besselj], page 566, [bessely], page 566, [besselh], page 568 .
```

H = besselh (alpha, x)

```
\(H=\) besselh (alpha, \(k, x\) )
\(H=\) besselh (alpha, \(k, x, o p t)\)
[ \(H\), ierr] = besselh (...)
Compute Bessel functions of the third kind (Hankel functions).
The order of the Bessel function alpha must be real. The kind of Hankel function is specified by \(k\) and may be either first \((k=1)\) or second \((k=2)\). The default is Hankel functions of the first kind. The points for evaluation \(x\) may be complex.
If the optional argument opt is 1 or true, the result is multiplied by \(\exp (-I * x)\) for \(k=1\) or \(\exp (I * x)\) for \(k=2\).
If alpha is a scalar, the result is the same size as \(x\). If \(x\) is a scalar, the result is the same size as alpha. If alpha is a row vector and \(x\) is a column vector, the result is a matrix with length (x) rows and length (alpha) columns. Otherwise, alpha and \(x\) must conform and the result will be the same size.
If requested, ierr contains the following status information and is the same size as the result.
0 . Normal return.
1. Input error, return NaN .
2. Overflow, return Inf.
3. Loss of significance by argument reduction results in less than half of machine accuracy.
4. Complete loss of significance by argument reduction, return NaN .
5. Error-no computation, algorithm termination condition not met, return NaN.

See also: [besselj], page 566, [bessely], page 566, [besseli], page 567, [besselk], page 568.
beta ( \(a, b\) )
Compute the Beta function for real inputs \(a\) and \(b\).
The Beta function definition is
\[
B(a, b)=\frac{\Gamma(a) \Gamma(b)}{\Gamma(a+b)}
\]

The Beta function can grow quite large and it is often more useful to work with the logarithm of the output rather than the function directly. See [betaln], page 570, for computing the logarithm of the Beta function in an efficient manner.
See also: [betaln], page 570, [betainc], page 569, [betaincinv], page 569 .
betainc ( \(x, a, b\) )
betainc (x, a, b, tail)
Compute the incomplete beta function.
This is defined as
\[
I_{x}(a, b)=\frac{1}{B(a, b)} \int_{0}^{x} t^{a-1}(1-t)^{b-1} d t
\]
with real \(x\) in the range \([0,1]\). The inputs a and \(b\) must be real and strictly positive (>0). If one of the inputs is not a scalar then the other inputs must be scalar or of compatible dimensions.
By default, tail is "lower" and the incomplete beta function integrated from 0 to \(x\) is computed. If tail is "upper" then the complementary function integrated from \(x\) to 1 is calculated. The two choices are related by
betainc (x, a, b, "upper") = 1-betainc (x, a, b, "lower").
betainc uses a more sophisticated algorithm than subtraction to get numerically accurate results when the "lower" value is small.
Reference: A. Cuyt, V. Brevik Petersen, B. Verdonk, H. Waadeland, W.B. Jones, Handbook of Continued Fractions for Special Functions, ch. 18.
See also: [beta], page 569, [betaincinv], page 569, [betaln], page 570.
betaincinv ( \(y, a, b\) )
betaincinv ( \(y, a, b\), "lower")
betaincinv (y, a, b, "upper")
Compute the inverse of the normalized incomplete beta function.
The normalized incomplete beta function is defined as
\[
I_{x}(a, b)=\frac{1}{B(a, b)} \int_{0}^{x} t^{a-1}(1-t)^{b-1} d t
\]

If two inputs are scalar, then betaincinv ( \(y, a, b\) ) is returned for each of the other inputs.
If two or more inputs are not scalar, the sizes of them must agree, and betaincinv is applied element-by-element.
The variable \(y\) must be in the interval \([0,1]\), while \(a\) and \(b\) must be real and strictly positive.
By default, tail is "lower" and the inverse of the incomplete beta function integrated from 0 to \(x\) is computed. If tail is "upper" then the complementary function integrated from \(x\) to 1 is inverted.
The function is computed by standard Newton's method, by solving
\[
y-I_{x}(a, b)=0
\]

See also: [betainc], page 569, [beta], page 569, [betaln], page 570.
betaln (a, b)
Compute the natural logarithm of the Beta function for real inputs a and \(b\).
betaln is defined as
\[
\operatorname{betaln}(a, b)=\ln (B(a, b)) \equiv \ln \left(\frac{\Gamma(a) \Gamma(b)}{\Gamma(a+b)}\right) .
\]
and is calculated in a way to reduce the occurrence of underflow.
The Beta function can grow quite large and it is often more useful to work with the logarithm of the output rather than the function directly.
See also: [beta], page 569, [betainc], page 569, [betaincinv], page 569, [gammaln], page 576.
bincoeff ( \(n, ~ k)\)
Return the binomial coefficient of \(n\) and \(k\).
The binomial coefficient is defined as
\[
\binom{n}{k}=\frac{n(n-1)(n-2) \cdots(n-k+1)}{k!}
\]

For example:
\[
\begin{aligned}
& \text { bincoeff }(5,2) \\
& \Rightarrow 10
\end{aligned}
\]

In most cases, the nchoosek function is faster for small scalar integer arguments. It also warns about loss of precision for big arguments.

See also: [nchoosek], page 785.
commutation_matrix (m, n)
Return the commutation matrix \(K_{m, n}\) which is the unique \(m n \times m n\) matrix such that \(K_{m, n} \cdot \operatorname{vec}(A)=\operatorname{vec}\left(A^{T}\right)\) for all \(m \times n\) matrices \(A\).
If only one argument \(m\) is given, \(K_{m, m}\) is returned.
See Magnus and Neudecker (1988), Matrix Differential Calculus with Applications in Statistics and Econometrics.
cosint (x)
Compute the cosine integral function:
\[
\mathrm{Ci}(x)=-\int_{x}^{\infty} \frac{\cos (t)}{t} d t
\]

An equivalent definition is
\[
\operatorname{Ci}(x)=\gamma+\log (x)+\int_{0}^{x} \frac{\cos (t)-1}{t} d t
\]

Reference:
M. Abramowitz and I.A. Stegun, Handbook of Mathematical Functions, 1964.

See also: [sinint], page 576, [expint], page 573, [cos], page 551.
duplication_matrix (n)
Return the duplication matrix \(D_{n}\) which is the unique \(n^{2} \times n(n+1) / 2\) matrix such that \(D_{n} * \operatorname{vech}(A)=\operatorname{vec}(A)\) for all symmetric \(n \times n\) matrices \(A\).
See Magnus and Neudecker (1988), Matrix Differential Calculus with Applications in Statistics and Econometrics.
dawson ( \(z\) )
Compute the Dawson (scaled imaginary error) function.
The Dawson function is defined as
\[
\frac{\sqrt{\pi}}{2} e^{-z^{2}} \operatorname{erfi}(z) \equiv-i \frac{\sqrt{\pi}}{2} e^{-z^{2}} \operatorname{erf}(i z)
\]

See also: [erfc], page 572, [erf], page 572, [erfcx], page 572, [erfi], page 573, [erfinv], page 573, [erfcinv], page 573.
[sn, cn, dn, err] = ellipj (u, m)
[sn, cn, dn, err] = ellipj (u, m, tol)
Compute the Jacobi elliptic functions sn, \(c n\), and \(d n\) of complex argument \(u\) and real parameter \(m\).
If \(m\) is a scalar, the results are the same size as \(u\). If \(u\) is a scalar, the results are the same size as \(m\). If \(u\) is a column vector and \(m\) is a row vector, the results are matrices with length ( \(u\) ) rows and length ( \(m\) ) columns. Otherwise, \(u\) and \(m\) must conform in size and the results will be the same size as the inputs.
The value of \(u\) may be complex. The value of \(m\) must be \(0 \leq m \leq 1\).
The optional input tol is currently ignored (matlab uses this to allow faster, less accurate approximation).
If requested, err contains the following status information and is the same size as the result.
0 . Normal return.
1. Error-no computation, algorithm termination condition not met, return NaN .

Reference: Milton Abramowitz and Irene A Stegun, Handbook of Mathematical Functions, Chapter 16 (Sections 16.4, 16.13, and 16.15), Dover, 1965.
See also: [ellipke], page 572.
\(k=\) ellipke (m)
\(k=\) ellipke (m, tol)
[k, e] = ellipke (...)
Compute complete elliptic integrals of the first \(\mathrm{K}(m)\) and second \(\mathrm{E}(m)\) kind.
\(m\) must be a scalar or real array with \(-\operatorname{Inf} \leq m \leq 1\).
The optional input tol controls the stopping tolerance of the algorithm and defaults to eps (class (m)). The tolerance can be increased to compute a faster, less accurate approximation.
When called with one output only elliptic integrals of the first kind are returned.
Mathematical Note:
Elliptic integrals of the first kind are defined as
\[
\mathrm{K}(m)=\int_{0}^{1} \frac{d t}{\sqrt{\left(1-t^{2}\right)\left(1-m t^{2}\right)}}
\]

Elliptic integrals of the second kind are defined as
\[
\mathrm{E}(m)=\int_{0}^{1} \frac{\sqrt{1-m t^{2}}}{\sqrt{1-t^{2}}} d t
\]

Reference: Milton Abramowitz and Irene A. Stegun, Handbook of Mathematical Functions, Chapter 17, Dover, 1965.
See also: [ellipj], page 571.
\(\operatorname{erf}(z)\)
Compute the error function.
The error function is defined as
\[
\operatorname{erf}(z)=\frac{2}{\sqrt{\pi}} \int_{0}^{z} e^{-t^{2}} d t
\]

See also: [erfc], page 572, [erfcx], page 572, [erfi], page 573, [dawson], page 571, [erfinv], page 573, [erfcinv], page 573.
erfc (z)
Compute the complementary error function.
The complementary error function is defined as \(1-\operatorname{erf}(z)\).
See also: [erfcinv], page 573, [erfcx], page 572, [erfi], page 573, [dawson], page 571, [erf], page 572, [erfinv], page 573 .
\(\operatorname{erfcx}(z)\)
Compute the scaled complementary error function.
The scaled complementary error function is defined as
\[
e^{z^{2}} \operatorname{erfc}(z) \equiv e^{z^{2}}(1-\operatorname{erf}(z))
\]

See also: [erfc], page 572, [erf], page 572, [erfi], page 573, [dawson], page 571, [erfinv], page 573 , [erfcinv], page 573.
erfi (z)
Compute the imaginary error function.
The imaginary error function is defined as
\[
-i \operatorname{erf}(i z)
\]

See also: [erfc], page 572, [erf], page 572, [erfcx], page 572, [dawson], page 571, [erfinv], page 573, [erfcinv], page 573.
erfinv (x)
Compute the inverse error function.
The inverse error function is defined such that
\[
\operatorname{erf}(y)==x
\]

See also: [erf], page 572, [erfc], page 572, [erfcx], page 572, [erfi], page 573, [dawson], page 571, [erfcinv], page 573.
erfcinv ( \(x\) )
Compute the inverse complementary error function.
The inverse complementary error function is defined such that
```

erfc (y) == x

```

See also: [erfc], page 572, [erf], page 572, [erfcx], page 572, [erfi], page 573, [dawson], page 571, [erfinv], page 573.
expint (x)
Compute the exponential integral.
The exponential integral is defined as:
\[
\mathrm{E}_{1}(x)=\int_{x}^{\infty} \frac{e^{-t}}{t} d t
\]

Note: For compatibility, this function uses the matlab definition of the exponential integral. Most other sources refer to this particular value as \(E_{1}(x)\), and the exponential integral as
\[
\operatorname{Ei}(x)=-\int_{-x}^{\infty} \frac{e^{-t}}{t} d t
\]

The two definitions are related, for positive real values of \(x\), by \(E_{1}(-x)=-\operatorname{Ei}(x)-i \pi\).
References:
M. Abramowitz and I.A. Stegun, Handbook of Mathematical Functions, 1964.
N. Bleistein and R.A. Handelsman, Asymptotic expansions of integrals, 1986.

See also: [cosint], page 571, [sinint], page 576, [exp], page 547.
gamma (z)
Compute the Gamma function.
The Gamma function is defined as
\[
\Gamma(z)=\int_{0}^{\infty} t^{z-1} e^{-t} d t
\]

Programming Note: The gamma function can grow quite large even for small input values. In many cases it may be preferable to use the natural logarithm of the gamma function (gammaln) in calculations to minimize loss of precision. The final result is then \(\exp\) (result_using_gammaln).
See also: [gammainc], page 574, [gammaln], page 576, [factorial], page 563.
gammainc ( \(x, a\) )
gammainc (x, a, tail)
Compute the normalized incomplete gamma function.
This is defined as
\[
\gamma(x, a)=\frac{1}{\Gamma(a)} \int_{0}^{x} t^{a-1} e^{-t} d t
\]
with the limiting value of 1 as \(x\) approaches infinity. The standard notation is \(P(a, x)\), e.g., Abramowitz and Stegun (6.5.1).

If \(a\) is scalar, then gammainc \((x, a)\) is returned for each element of \(x\) and vice versa.
If neither \(x\) nor \(a\) is scalar then the sizes of \(x\) and a must agree, and gammainc is applied element-by-element. The elements of a must be non-negative.
By default, tail is "lower" and the incomplete gamma function integrated from 0 to \(x\) is computed. If tail is "upper" then the complementary function integrated from \(x\) to infinity is calculated.
If tail is "scaledlower", then the lower incomplete gamma function is multiplied by \(\Gamma(a+1) \exp (x) x^{-a}\). If tail is "scaledupper", then the upper incomplete gamma function is multiplied by the same quantity.
References:
M. Abramowitz and I.A. Stegun, Handbook of mathematical functions, Dover publications, Inc., 1972.
W. Gautschi, A computational procedure for incomplete gamma functions, ACM Trans. Math Software, pp. 466-481, Vol 5, No. 4, 2012.
W. H. Press, S. A. Teukolsky, W. T. Vetterling, and B. P. Flannery, Numerical Recipes in Fortran 77, ch. 6.2, Vol 1, 1992.

See also: [gamma], page 574, [gammaincinv], page 574, [gammaln], page 576.
gammaincinv ( \(y, a)\)
gammaincinv (y, a, tail)
Compute the inverse of the normalized incomplete gamma function.
The normalized incomplete gamma function is defined as
\[
\gamma(x, a)=\frac{1}{\Gamma(a)} \int_{0}^{x} t^{a-1} e^{-t} d t
\]
and gammaincinv (gammainc \((x, a), a)=x\) for each non-negative value of \(x\). If \(a\) is scalar then gammaincinv \((y, a)\) is returned for each element of \(y\) and vice versa.
If neither \(y\) nor a is scalar then the sizes of \(y\) and a must agree, and gammaincinv is applied element-by-element. The variable \(y\) must be in the interval \([0,1]\) while a must be real and positive.
By default, tail is "lower" and the inverse of the incomplete gamma function integrated from 0 to \(x\) is computed. If tail is "upper", then the complementary function integrated from \(x\) to infinity is inverted.
The function is computed with Newton's method by solving
\[
y-\gamma(x, a)=0
\]

Reference: A. Gil, J. Segura, and N. M. Temme, Efficient and accurate algorithms for the computation and inversion of the incomplete gamma function ratios, SIAM J. Sci. Computing, pp. A2965-A2981, Vol 34, 2012.
See also: [gammainc], page 574, [gamma], page 574, [gammaln], page 576.
\(1=\) legendre ( \(n, x\) )
1 = legendre ( \(n, x\), normalization)
Compute the associated Legendre function of degree \(n\) and order \(m=0 \ldots n\).
The value \(n\) must be a real non-negative integer.
\(x\) is a vector with real-valued elements in the range \([-1,1]\).
The optional argument normalization may be one of "unnorm", "sch", or "norm". The default if no normalization is given is "unnorm".
When the optional argument normalization is "unnorm", compute the associated Legendre function of degree \(n\) and order \(m\) and return all values for \(m=0 \ldots n\). The return value has one dimension more than \(x\).
The associated Legendre function of degree \(n\) and order \(m\) :
\[
P_{n}^{m}(x)=(-1)^{m}\left(1-x^{2}\right)^{m / 2} \frac{d^{m}}{d x^{m}} P_{n}(x)
\]
with Legendre polynomial of degree n:
\[
P(x)=\frac{1}{2^{n} n!}\left(\frac{d^{n}}{d x^{n}}\left(x^{2}-1\right)^{n}\right)
\]
legendre (3, \([-1.0,-0.9,-0.8]\) ) returns the matrix:
\begin{tabular}{c|c|c|c}
x & \(\mid\) & -1.0 & -0.9 \\
\hline--0.8 \\
\(\mathrm{~m}=0\) & -1.00000 & -0.47250 & -0.08000 \\
\(\mathrm{~m}=1\) & 0.00000 & -1.99420 & -1.98000 \\
\(\mathrm{~m}=2\) & 0.00000 & -2.56500 & -4.32000 \\
\(\mathrm{~m}=3\) & & 0.00000 & -1.24229
\end{tabular}\(|-3.24000\)

When the optional argument normalization is "sch", compute the Schmidt seminormalized associated Legendre function. The Schmidt semi-normalized associated Legendre function is related to the unnormalized Legendre functions by the following:

For Legendre functions of degree \(n\) and order 0 :
\[
S P_{n}^{0}(x)=P_{n}^{0}(x)
\]

For Legendre functions of degree n and order m :
\[
S P_{n}^{m}(x)=P_{n}^{m}(x)(-1)^{m}\left(\frac{2(n-m)!}{(n+m)!}\right)^{0.5}
\]

When the optional argument normalization is "norm", compute the fully normalized associated Legendre function. The fully normalized associated Legendre function is related to the unnormalized associated Legendre functions by the following:
For Legendre functions of degree \(n\) and order \(m\)
\[
N P_{n}^{m}(x)=P_{n}^{m}(x)(-1)^{m}\left(\frac{(n+0.5)(n-m)!}{(n+m)!}\right)^{0.5}
\]
gammaln (x)
lgamma (x)
Return the natural logarithm of the gamma function of \(x\).
See also: [gamma], page 574, [gammainc], page 574.
psi (z)
psi (k, z)
Compute the psi (polygamma) function.
The polygamma functions are the \(k\) th derivative of the logarithm of the gamma function. If unspecified, \(k\) defaults to zero. A value of zero computes the digamma function, a value of 1 , the trigamma function, and so on.
The digamma function is defined:
\[
\Psi(z)=\frac{d(\log (\Gamma(z)))}{d x}
\]

When computing the digamma function (when \(k\) equals zero), \(z\) can have any value real or complex value. However, for polygamma functions ( \(k\) higher than 0 ), z must be real and non-negative.
See also: [gamma], page 574, [gammainc], page 574, [gammaln], page 576.
sinint (x)
Compute the sine integral function:
\[
\operatorname{Si}(x)=\int_{0}^{x} \frac{\sin (t)}{t} d t
\]

Reference: M. Abramowitz and I.A. Stegun, Handbook of Mathematical Functions, 1964.

See also: [cosint], page 571, [expint], page 573 , [sin], page 551.

\subsection*{17.7 Rational Approximations}
\(s=r a t(x)\)
\(s=r a t(x, t o l)\)
\([n, d]=\operatorname{rat}(\ldots)\)
Find a rational approximation of \(x\) to within the tolerance defined by tol.
If unspecified, the default tolerance is \(1 \mathrm{e}-6 * \operatorname{norm}(x(:), 1)\).
When called with one output argument, return a string containing a continued fraction expansion (multiple terms).

When called with two output arguments, return numeric matrices for the numerator and denominator of a fractional representation of \(x\) such that \(x=n . / d\).

For example:
```

s = rat (pi)
m s = 3 + 1/(7 + 1/16)
[n, d] = rat (pi)

# n = 355

d = 113
n / d - pi

# 0.00000026676

```

Programming Note: With one output rat produces a string which is a continued fraction expansion. To produce a string which is a simple fraction (one numerator, one denominator) use rats.

See also: [rats], page 577, [format], page 272.
\(s=r a t s(x)\)
\(s=\operatorname{rats}(x\), len)
Convert x into a rational approximation represented as a string.
A rational approximation to a floating point number is a simple fraction with numerator \(N\) and denominator \(D\) such that \(x=N / D\).

The optional second argument defines the maximum length of the string representing the elements of x. By default, len is 9 .

If the length of the smallest possible rational approximation exceeds len, an asterisk \(\left.{ }^{*}\right)\) padded with spaces will be returned instead.

Example conversion from matrix to string, and back again.
```

r = rats (hilb (4));
x = str2num (r)

```

See also: [rat], page 577, [format], page 272.

\subsection*{17.8 Coordinate Transformations}
[theta, \(r\) ] = cart2pol ( \(x, y\) )
[theta, \(r, z]=\) cart2pol (x, y, z)
[theta, r] = cart2pol (C)
[theta, \(r, z]=\) cart2pol (C)
Transform Cartesian coordinates to polar or cylindrical coordinates.
The inputs \(x, y(\), and \(z)\) must be the same shape, or scalar. If called with a single matrix argument then each row of \(C\) represents the Cartesian coordinate pair ( \(x, y\) ) or triplet ( \(x, y, z\) ).
The outputs theta, r (, and \(z\) ) match the shape of the inputs. For a matrix input \(C\) the outputs will be column vectors with rows corresponding to the rows of the input matrix.
theta describes the angle relative to the positive x -axis measured in the xy-plane.
\(r\) is the distance to the z -axis \((0,0, \mathrm{z})\).
\(z\), if present, is unchanged by the transformation.
The coordinate transformation is computed using:
\[
\begin{gathered}
\theta=\arctan \left(\frac{y}{x}\right) \\
r=\sqrt{x^{2}+y^{2}} \\
z=z
\end{gathered}
\]

Note: For matlab compatibility, this function no longer returns a full coordinate matrix when called with a single return argument.
See also: [pol2cart], page 578, [cart2sph], page 579, [sph2cart], page 579.
[x, y] = pol2cart (theta, r)
[x, y, z] = pol2cart (theta, r, z)
[x, y] = pol2cart (P)
\([x, y, z]=\operatorname{pol} 2 \operatorname{cart}(P)\)
Transform polar or cylindrical coordinates to Cartesian coordinates.
The inputs theta, \(r,(\operatorname{and} z)\) must be the same shape, or scalar. If called with a single matrix argument then each row of \(P\) represents the polar coordinate pair (theta, \(r\) ) or the cylindrical triplet (theta, \(r, z\) ).
The outputs \(x, y(\), and \(z)\) match the shape of the inputs. For a matrix input \(P\) the outputs will be column vectors with rows corresponding to the rows of the input matrix.
theta describes the angle relative to the positive x -axis measured in the xy-plane.
\(r\) is the distance to the z -axis \((0,0, \mathrm{z})\).
\(z\), if present, is unchanged by the transformation.
The coordinate transformation is computed using:
\[
x=r \cos \theta
\]
\[
\begin{gathered}
y=r \sin \theta \\
z=z
\end{gathered}
\]

Note: For matlab compatibility, this function no longer returns a full coordinate matrix when called with a single return argument.

See also: [cart2pol], page 578, [sph2cart], page 579, [cart2sph], page 579.
[theta, phi, r] \(=\operatorname{cart2sph}(x, y, z)\)
[theta, phi, r] = cart2sph (C)
Transform Cartesian coordinates to spherical coordinates.
The inputs \(x, y\), and \(z\) must be the same shape, or scalar. If called with a single matrix argument then each row of \(C\) must represent a Cartesian coordinate triplet ( \(x, y, z\) ).
The outputs theta, phi, r match the shape of the inputs. For a matrix input \(C\) the outputs will be column vectors with rows corresponding to the rows of the input matrix.
theta describes the azimuth angle relative to the positive x -axis measured in the xy-plane.
phi is the elevation angle measured relative to the xy-plane.
\(r\) is the distance to the origin \((0,0,0)\).
The coordinate transformation is computed using:
\[
\begin{gathered}
\theta=\arctan \left(\frac{y}{x}\right) \\
\phi=\arctan \left(\frac{z}{\sqrt{x^{2}+y^{2}}}\right) \\
r=\sqrt{x^{2}+y^{2}+z^{2}}
\end{gathered}
\]

Note: For matlab compatibility, this function no longer returns a full coordinate matrix when called with a single return argument.
See also: [sph2cart], page 579, [cart2pol], page 578, [pol2cart], page 578.
[x, y, z] = sph2cart (theta, phi, r)
\([x, y, z]=\operatorname{sph} 2 \operatorname{cart}(S)\)
Transform spherical coordinates to Cartesian coordinates.
The inputs theta, phi, and \(r\) must be the same shape, or scalar. If called with a single matrix argument then each row of \(S\) must represent a spherical coordinate triplet (theta, phi, r).
The outputs \(x, y, z\) match the shape of the inputs. For a matrix input \(S\) the outputs are column vectors with rows corresponding to the rows of the input matrix.
theta describes the azimuth angle relative to the positive x -axis measured in the xy-plane.
\(p h i\) is the elevation angle measured relative to the xy-plane.
\(r\) is the distance to the origin \((0,0,0)\).

The coordinate transformation is computed using:
\[
\begin{gathered}
x=r \cos \phi \cos \theta \\
y=r \cos \phi \sin \theta \\
z=r \sin \phi
\end{gathered}
\]

Note: For matlab compatibility, this function no longer returns a full coordinate matrix when called with a single return argument.
See also: [cart2sph], page 579, [pol2cart], page 578, [cart2pol], page 578.

\subsection*{17.9 Mathematical Constants}
e
e (n)
e \((n, m)\)
e ( \(n, m, k, \ldots\) )
e (..., class)
Return a scalar, matrix, or N-dimensional array whose elements are all equal to the base of natural logarithms.
The constant \(e\) satisfies the equation \(\log (e)=1\).
When called with no arguments, return a scalar with the value \(e\).
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
The optional argument class specifies the return type and may be either "double" or "single".

See also: [log], page 547, [exp], page 547, [pi], page 580, [I], page 581.
pi
pi (n)
pi (n, m)
pi ( \(n, m, k, \ldots\) )
pi (..., class)
Return a scalar, matrix, or N-dimensional array whose elements are all equal to the ratio of the circumference of a circle to its diameter \((\pi)\).
Internally, pi is computed as ' \(4.0 *\) atan (1.0)'.
When called with no arguments, return a scalar with the value of \(\pi\).
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.

The optional argument class specifies the return type and may be either "double" or "single".
See also: [e], page 580, [I], page 581.
I
I (n)
I \((n, m)\)
I ( \(n, m, k, \ldots\) )
I (..., class)
Return a scalar, matrix, or N-dimensional array whose elements are all equal to the pure imaginary unit, defined as \(\sqrt{-1}\).
\(I\), and its equivalents \(i, j\), and \(J\), are functions so any of the names may be reused for other purposes (such as i for a counter variable).
When called with no arguments, return a scalar with the value \(i\).
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [e], page 580, [pi], page 580, [log], page 547, [exp], page 547.
```

Inf
Inf (n)
Inf (n, m)
Inf (n,m,k,...)
Inf (..., class)

```

Return a scalar, matrix or N -dimensional array whose elements are all equal to the IEEE representation for positive infinity.
Infinity is produced when results are too large to be represented using the IEEE floating point format for numbers. Two common examples which produce infinity are division by zero and overflow.
```

[ 1/0 e^800 ]

# Inf Inf

```

When called with no arguments, return a scalar with the value 'Inf'.
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [isinf], page 512, [NaN], page 582.
```

val = NaN
val = NaN (n)
val = NaN (n,m)
val = NaN (n,m,k,···)
val = NaN (..., "like", var))
val = NaN (..., class)

```

Return a scalar, matrix, or N-dimensional array whose elements are all equal to the IEEE symbol NaN (Not a Number).
NaN is the result of operations which do not produce a well defined numerical result. Common operations which produce a NaN are arithmetic with infinity \((\infty-\infty)\), zero divided by zero ( \(0 / 0\) ), and any operation involving another NaN value ( \(5+\mathrm{NaN}\) ).
Note that NaN always compares not equal to \(\mathrm{NaN}(\mathrm{NaN}!=\mathrm{NaN})\). This behavior is specified by the IEEE standard for floating point arithmetic. To find NaN values, use the isnan function.
When called with no arguments, return a scalar with the value ' NaN '.
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
If a variable var is specified after "like", the output val will have the same data type, complexity, and sparsity as var.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [isnan], page 512, [Inf], page 581.

\section*{eps}
eps (x)
eps ( \(n, m\) )
eps ( \(n, m, k, \ldots)\)
eps (..., class)
Return a scalar, matrix or N-dimensional array whose elements are all eps, the machine precision.
More precisely, eps is the relative spacing between any two adjacent numbers in the machine's floating point system. This number is obviously system dependent. On machines that support IEEE floating point arithmetic, eps is approximately \(2.2204 \times\) \(10^{-16}\) for double precision and \(1.1921 \times 10^{-7}\) for single precision.
When called with no arguments, return a scalar with the value eps (1.0).
Given a single argument \(x\), return the distance between \(x\) and the next largest value.
When called with more than one argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions. The optional argument class specifies the return type and may be either "double" or "single".
See also: [realmax], page 583, [realmin], page 583, [intmax], page 59, [flintmax], page 60.
```

realmax

```
realmax ( \(n\) )
realmax ( \(n, m\) )
realmax ( \(n, m, k, \ldots\) )
realmax (..., class)
Return a scalar, matrix, or N-dimensional array whose elements are all equal to the largest floating point number that is representable.
The actual value is system dependent. On machines that support IEEE floating point arithmetic, realmax is approximately \(1.7977 \times 10^{308}\) for double precision and \(3.4028 \times 10^{38}\) for single precision.
When called with no arguments, return a scalar with the value realmax ("double"). When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [realmin], page 583, [intmax], page 59, [flintmax], page 60, [eps], page 582.
realmin
realmin ( \(n\) )
realmin ( \(n, m\) )
realmin ( \(n, m, k, \ldots\) )
realmin (..., class)
Return a scalar, matrix, or N-dimensional array whose elements are all equal to the smallest normalized floating point number that is representable.
The actual value is system dependent. On machines that support IEEE floating point arithmetic, realmin is approximately \(2.2251 \times 10^{-308}\) for double precision and \(1.1755 \times 10^{-38}\) for single precision.
When called with no arguments, return a scalar with the value realmin ("double").
When called with a single argument, return a square matrix with the dimension specified.
When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions.
The optional argument class specifies the return type and may be either "double" or "single".
See also: [realmax], page 583, [intmin], page 60, [eps], page 582.

\section*{18 Linear Algebra}

This chapter documents the linear algebra functions provided in Octave. Reference material for many of these functions may be found in Golub and Van Loan, Matrix Computations, 2nd Ed., Johns Hopkins, 1989, and in the lapack Users' Guide, SIAM, 1992. The lapack Users' Guide is available at: http://www.netlib.org/lapack/lug/

A common text for engineering courses is G. Strang, Linear Algebra and Its Applications, 4th Edition. It has become a widespread reference for linear algebra. An alternative is P. Lax Linear Algebra and Its Applications, and also is a good choice. It claims to be suitable for high school students with substantial mathematical interests as well as firstyear undergraduates.

\subsection*{18.1 Techniques Used for Linear Algebra}

Octave includes a polymorphic solver that selects an appropriate matrix factorization depending on the properties of the matrix itself. Generally, the cost of determining the matrix type is small relative to the cost of factorizing the matrix itself. In any case the matrix type is cached once it is calculated so that it is not re-determined each time it is used in a linear equation.

The selection tree for how the linear equation is solved or a matrix inverse is formed is given by:
1. If the matrix is upper or lower triangular sparse use a forward or backward substitution using the LAPACK xTRTRS function, and goto 4.
2. If the matrix is square, Hermitian with a real positive diagonal, attempt Cholesky factorization using the LAPACK xPOTRF function.
3. If the Cholesky factorization failed or the matrix is not Hermitian with a real positive diagonal, and the matrix is square, factorize using the LAPACK xGETRF function.
4. If the matrix is not square, or any of the previous solvers flags a singular or near singular matrix, find a least squares solution using the LAPACK xGELSD function.

The user can force the type of the matrix with the matrix_type function. This overcomes the cost of discovering the type of the matrix. However, it should be noted that identifying the type of the matrix incorrectly will lead to unpredictable results, and so matrix_type should be used with care.

It should be noted that the test for whether a matrix is a candidate for Cholesky factorization, performed above, and by the matrix_type function, does not make certain that the matrix is Hermitian. However, the attempt to factorize the matrix will quickly detect a non-Hermitian matrix.

\subsection*{18.2 Basic Matrix Functions}
\(A A=\) balance \((A)\)
\(A A=\) balance ( \(A\), opt)
[DD, AA] = balance (A, opt)
\([D, P, A A]=\) balance (A, opt)
[CC, \(D D, A A, B B]=\) balance \((A, B, o p t)\)
Balance the matrix \(A\) to reduce numerical errors in future calculations.
Compute \(A A=D D \backslash A * D D\) in which \(A A\) is a matrix whose row and column norms are roughly equal in magnitude, and \(D D=P * D\), in which \(P\) is a permutation matrix and \(D\) is a diagonal matrix of powers of two. This allows the equilibration to be computed without round-off. Results of eigenvalue calculation are typically improved by balancing first.
If two output values are requested, balance returns the diagonal \(D\) and the permutation \(P\) separately as vectors. In this case, \(D D=\operatorname{eye}(\mathrm{n})(:, P) * \operatorname{diag}(D)\), where \(n\) is the matrix size.
If four output values are requested, compute \(A A=C C * A * D D\) and \(B B=C C * B * D D\), in which \(A A\) and \(B B\) have nonzero elements of approximately the same magnitude and \(C C\) and \(D D\) are permuted diagonal matrices as in \(D D\) for the algebraic eigenvalue problem.
The eigenvalue balancing option opt may be one of:
"noperm", "S"
Scale only; do not permute.
"noscal", "P"
Permute only; do not scale.
Algebraic eigenvalue balancing uses standard LAPACK routines.
Generalized eigenvalue problem balancing uses Ward's algorithm (SIAM Journal on Scientific and Statistical Computing, 1981).
bw = bandwidth (A, type)
[lower, upper] = bandwidth (A)
Compute the bandwidth of \(A\).
The type argument is the string "lower" for the lower bandwidth and "upper" for the upper bandwidth. If no type is specified return both the lower and upper bandwidth of \(A\).
The lower/upper bandwidth of a matrix is the number of subdiagonals/superdiagonals with nonzero entries.
See also: [isbanded], page 69, [isdiag], page 70, [istril], page 70, [istriu], page 70.
cond (A)
cond ( \(A, p\) )
Compute the \(p\)-norm condition number of a matrix with respect to inversion.
cond ( \(A\) ) is defined as \(\|A\|_{p} *\left\|A^{-1}\right\|_{p}\).
By default, \(p=2\) is used which implies a (relatively slow) singular value decomposition. Other possible selections are \(p=1\), Inf, "fro" which are generally faster. For a full discussion of possible \(p\) values, see [norm], page 592.
The condition number of a matrix quantifies the sensitivity of the matrix inversion operation when small changes are made to matrix elements. Ideally the condition number will be close to 1 . When the number is large this indicates small changes (such as underflow or round-off error) will produce large changes in the resulting
output. In such cases the solution results from numerical computing are not likely to be accurate.
See also: [condest], page 683, [rcond], page 594, [condeig], page 587, [norm], page 592, [svd], page 607.
\(c=\) condeig (a)
[v, lambda, c] = condeig (a)
Compute condition numbers of a matrix with respect to eigenvalues.
The condition numbers are the reciprocals of the cosines of the angles between the left and right eigenvectors; Large values indicate that the matrix has multiple distinct eigenvalues.
The input a must be a square numeric matrix.
The outputs are:
- \(c\) is a vector of condition numbers for the eigenvalues of a.
- \(v\) is the matrix of right eigenvectors of a. The result is equivalent to calling [ \(v\), lambda] = eig (a).
- lambda is the diagonal matrix of eigenvalues of a. The result is equivalent to calling [ \(\mathrm{v}, \mathrm{lambda}\) ] \(=\) eig (a).
Example
```

a = [1, 2; 3, 4];
c = condeig (a)
\# c =
1.0150
1.0150

```

See also: [eig], page 587, [cond], page 586, [balance], page 585.
\(\operatorname{det}(A)\)
\([d, r c o n d]=\operatorname{det}(A)\)
Compute the determinant of \(A\).
Return an estimate of the reciprocal condition number if requested.
Programming Notes: Routines from Lapack are used for full matrices and code from UMFPACK is used for sparse matrices.
The determinant should not be used to check a matrix for singularity. For that, use any of the condition number functions: cond, condest, rcond.
See also: [cond], page 586, [condest], page 683, [rcond], page 594.
```

lambda = eig (A)
lambda $=\operatorname{eig}(A, B)$
[V, lambda] = eig (A)
[V, lambda] = eig (A, B)
$[V$, lambda, $W$ ] $=\operatorname{eig}(A)$
$[V, \operatorname{lambda}, W]=\operatorname{eig}(A, B)$
[...] = eig (A, balanceOption)
[...] = eig (A, B, algorithm)

```
[...] = eig (..., eigvalOption)
Compute the eigenvalues (lambda) and optionally the right eigenvectors ( \(V\) ) and the left eigenvectors \((W)\) of a matrix or pair of matrices.
The flag balanceOption can be one of:
"balance" (default)
Preliminary balancing is on.
"nobalance"
Disables preliminary balancing.
The flag eigvalOption can be one of:
"matrix" Return the eigenvalues in a diagonal matrix. (default if 2 or 3 outputs are requested)
"vector" Return the eigenvalues in a column vector. (default if only 1 output is requested, e.g., lambda \(=\operatorname{eig}(A))\)
The flag algorithm can be one of:
"chol" Use the Cholesky factorization of B . (default if \(A\) is symmetric (Hermitian) and \(B\) is symmetric (Hermitian) positive definite)
"qz" Use the QZ algorithm. (used whenever \(A\) or \(B\) are not symmetric)
\begin{tabular}{lllll} 
& & no flag & chol & qz \\
both are symmetric & & "chol" & "chol" & "qz" \\
at least one is not & "qz" & "qz" & "qz"
\end{tabular}
symmetric
The eigenvalues returned by eig are not ordered.
See also: [eigs], page 686, [svd], page 607.
\(G=\) givens ( \(x, y\) )
[c,s] = givens (x, y)
Compute the Givens rotation matrix \(G\).
The Givens matrix is a \(2 \times 2\) orthogonal matrix
\[
G=\left[\begin{array}{cc}
c & s \\
-s^{\prime} & c
\end{array}\right]
\]
such that
\[
G\left[\begin{array}{l}
x \\
y
\end{array}\right]=\left[\begin{array}{l}
* \\
0
\end{array}\right]
\]
with \(x\) and \(y\) scalars.
If two output arguments are requested, return the factors \(c\) and \(s\) rather than the Givens rotation matrix.
For example:
```

givens (1, 1)
=> 0.70711

```

Note: The Givens matrix represents a counterclockwise rotation of a 2-D plane and can be used to introduce zeros into a matrix prior to complete factorization.

See also: [planerot], page 589, [qr], page 599.
\(S=\operatorname{gsvd}(A, B)\)
\([U, V, X, C, S]=\operatorname{gsvd}(A, B)\)
\([U, V, X, C, S]=\operatorname{gsvd}(A, B, 0)\)
Compute the generalized singular value decomposition of \((A, B)\).
The generalized singular value decomposition is defined by the following relations:
\[
\begin{aligned}
A & =U C X^{\dagger} \\
B & =V S X^{\dagger} \\
C^{\dagger} C+S^{\dagger} S & =\text { eye }(\operatorname{columns}(A))
\end{aligned}
\]

The function gsvd normally returns just the vector of generalized singular values
\[
\sqrt{\frac{\operatorname{diag}\left(C^{\dagger} C\right)}{\operatorname{diag}\left(S^{\dagger} S\right)}}
\]

If asked for five return values, it also computes \(U, V, X\), and \(C\).
If the optional third input is present, gsvd constructs the "economy-sized" decomposition where the number of columns of \(U, V\) and the number of rows of \(C, S\) is less than or equal to the number of columns of \(A\). This option is not yet implemented.
Programming Note: the code is a wrapper to the corresponding LAPACK dggsvd and zggsvd routines. If matrices \(A\) and \(B\) are both rank deficient then Lapack will return an incorrect factorization. Programmers should avoid this combination.

See also: [svd], page 607.
[G, y] = planerot (x)
Compute the Givens rotation matrix for the two-element column vector x .
The Givens matrix is a \(2 \times 2\) orthogonal matrix
\[
G=\left[\begin{array}{cc}
c & s \\
-s^{\prime} & c
\end{array}\right]
\]
such that
\[
G\left[\begin{array}{l}
x(1) \\
x(2)
\end{array}\right]=\left[\begin{array}{l}
* \\
0
\end{array}\right]
\]

Note: The Givens matrix represents a counterclockwise rotation of a 2-D plane and can be used to introduce zeros into a matrix prior to complete factorization.
See also: [givens], page 588, [qr], page 599.
\(x=\operatorname{inv}(A)\)
[x, rcond] = inv (A)
[...] = inverse (...)
Compute the inverse of the square matrix \(A\).
Return an estimate of the reciprocal condition number if requested, otherwise warn of an ill-conditioned matrix if the reciprocal condition number is small.
In general it is best to avoid calculating the inverse of a matrix directly. For example, it is both faster and more accurate to solve systems of equations \(\left(A^{*} x=b\right)\) with \(\mathrm{y}=\) \(A \backslash b\), rather than \(\mathrm{y}=\operatorname{inv}(A) * b\).
If called with a sparse matrix, then in general \(x\) will be a full matrix requiring significantly more storage. Avoid forming the inverse of a sparse matrix if possible.
inverse is an alias and may be used identically in place of inv.
See also: [ldivide], page 162, [rdivide], page 163, [pinv], page 593.
\(\mathrm{x}=\operatorname{linsolve}(\mathrm{A}, \mathrm{b})\)
\(x\) = linsolve (A, b, opts)
[x, R] = linsolve (...)
Solve the linear system \(\mathrm{A} * \mathrm{x}=\mathrm{b}\).
With no options, this function is equivalent to the left division operator ( \(\mathrm{x}=\mathrm{A} \backslash \mathrm{b}\) ) or the matrix-left-divide function ( \(\mathrm{x}=\mathrm{mldivide}(\mathrm{A}, \mathrm{b})\) ).
Octave ordinarily examines the properties of the matrix \(A\) and chooses a solver that best matches the matrix. By passing a structure opts to linsolve you can inform Octave directly about the matrix \(A\). In this case Octave will skip the matrix examination and proceed directly to solving the linear system.

Warning: If the matrix \(A\) does not have the properties listed in the opts structure then the result will not be accurate AND no warning will be given. When in doubt, let Octave examine the matrix and choose the appropriate solver as this step takes little time and the result is cached so that it is only done once per linear system.
Possible opts fields (set value to true/false):
LT \(\quad A\) is lower triangular
UT \(\quad A\) is upper triangular
UHESS \(\quad A\) is upper Hessenberg (currently makes no difference)
SYM \(\quad A\) is symmetric or complex Hermitian (currently makes no difference)
POSDEF \(A\) is positive definite
RECT \(\quad A\) is general rectangular (currently makes no difference)
TRANSA Solve \(\mathrm{A}^{\prime} * \mathrm{x}=\mathrm{b}\) if true rather than \(\mathrm{A} * \mathrm{x}=\mathrm{b}\)
The optional second output \(R\) is the inverse condition number of \(A\) (zero if matrix is singular).
See also: [mldivide], page 162, [matrix_type], page 591, [rcond], page 594.
```

type = matrix_type (A)
type = matrix_type (A, "nocompute")
A = matrix_type (A, type)
A = matrix_type (A, "upper", perm)
A = matrix_type (A, "lower", perm)
A = matrix_type (A, "banded", nl, nu)

```

Identify the matrix type or mark a matrix as a particular type.
This allows more rapid solutions of linear equations involving \(A\) to be performed.
Called with a single argument, matrix_type returns the type of the matrix and caches it for future use.
Called with more than one argument, matrix_type allows the type of the matrix to be defined.
If the option "nocompute" is given, the function will not attempt to guess the type if it is still unknown. This is useful for debugging purposes.
The possible matrix types depend on whether the matrix is full or sparse, and can be one of the following
"unknown"
Remove any previously cached matrix type, and mark type as unknown.
"full" Mark the matrix as full.
"positive definite"
Probable full positive definite matrix.
"diagonal"
Diagonal matrix. (Sparse matrices only)
"permuted diagonal"
Permuted Diagonal matrix. The permutation does not need to be specifically indicated, as the structure of the matrix explicitly gives this. (Sparse matrices only)
"upper" Upper triangular. If the optional third argument perm is given, the matrix is assumed to be a permuted upper triangular with the permutations defined by the vector perm.
"lower" Lower triangular. If the optional third argument perm is given, the matrix is assumed to be a permuted lower triangular with the permutations defined by the vector perm.
"banded"
"banded positive definite"
Banded matrix with the band size of \(n l\) below the diagonal and \(n u\) above it. If \(n l\) and \(n u\) are 1 , then the matrix is tridiagonal and treated with specialized code. In addition the matrix can be marked as probably a positive definite. (Sparse matrices only)
"singular"
The matrix is assumed to be singular and will be treated with a minimum norm solution.

Note that the matrix type will be discovered automatically on the first attempt to solve a linear equation involving \(A\). Therefore matrix_type is only useful to give Octave hints of the matrix type. Incorrectly defining the matrix type will result in incorrect results from solutions of linear equations; it is entirely the responsibility of the user to correctly identify the matrix type.
Also, the test for positive definiteness is a low-cost test for a Hermitian matrix with a real positive diagonal. This does not guarantee that the matrix is positive definite, but only that it is a probable candidate. When such a matrix is factorized, a Cholesky factorization is first attempted, and if that fails the matrix is then treated with an LU factorization. Once the matrix has been factorized, matrix_type will return the correct classification of the matrix.
norm (A)
norm ( \(A, p\) )
norm ( \(A, p, o p t\) )
Compute the p-norm of the matrix \(A\).
If the second argument is not given, \(\mathrm{p}=2\) is used.
If \(A\) is a matrix (or sparse matrix):
\(p=1 \quad\) 1-norm, the largest column sum of the absolute values of \(A\).
\(p=2 \quad\) Largest singular value of \(A\).
\(p=\operatorname{Inf}\) or "inf"
Infinity norm, the largest row sum of the absolute values of \(A\).
\(p=\) "fro"
Frobenius norm of \(A\), sqrt (sum (diag \(\left.\left(A^{\prime} * A\right)\right)\) ).
other \(p, p>1\)
maximum norm ( \(\mathrm{A} * \mathrm{x}, \mathrm{p}\) ) such that norm ( \(\mathrm{x}, \mathrm{p}\) ) \(=1\)
If \(A\) is a vector or a scalar:
\[
\begin{aligned}
& p=\operatorname{Inf} \text { or } \operatorname{linf} " \\
& \max (\operatorname{abs}(A)) . \\
& p=-\operatorname{Inf} \min (\operatorname{abs}(A)) . \\
& p=\text { fro" }
\end{aligned}
\]

Frobenius norm of \(A\), sqrt (sumsq (abs (A))).
\(p=0 \quad\) Hamming norm-the number of nonzero elements.
other \(p, p>1\)
p-norm of \(A\), (sum (abs (A) .^ \(p\) ) ^ ( \(1 / p\) ).
other \(p\) p < 1
the p-pseudonorm defined as above.
If opt is the value "rows", treat each row as a vector and compute its norm. The result is returned as a column vector. Similarly, if opt is "columns" or "cols" then compute the norms of each column and return a row vector.
See also: [normest], page 682, [normest1], page 682, [vecnorm], page 594, [cond], page 586, [svd], page 607.
\(Z=\operatorname{null}(A)\)
\(Z=\) null (A, tol)
Return an orthonormal basis \(Z\) of the null space of \(A\).
The dimension of the null space \(Z\) is taken as the number of singular values of \(A\) not greater than tol. If the argument tol is missing, it is computed as
```

max (size (A)) * max (svd (A, O)) * eps

```

See also: [orth], page 593, [svd], page 607.
orth (A)
orth (A, tol)
Return an orthonormal basis of the range space of \(A\).
The dimension of the range space is taken as the number of singular values of \(A\) greater than tol. If the argument tol is missing, it is computed as
```

max (size (A)) * max (svd (A)) * eps

```

See also: [null], page 593.
\([y, h]=\operatorname{mgorth}(x, v)\)
Orthogonalize a given column vector \(x\) with respect to a set of orthonormal vectors comprising the columns of \(v\) using the modified Gram-Schmidt method.
On exit, \(y\) is a unit vector such that:
\[
\begin{aligned}
& \text { norm }(y)=1 \\
& v^{\prime} * y=0 \\
& x=[v, y] * h^{\prime}
\end{aligned}
\]
pinv ( \(x\) )
pinv ( \(x\), tol)
Return the Moore-Penrose pseudoinverse of \(x\).
Singular values less than tol are ignored.
If the second argument is omitted, it is taken to be
```

tol = max ([rows(x), columns(x)]) * norm (x) * eps

```

See also: [inv], page 590, [ldivide], page 162.
rank (A)
rank (A, tol)
Compute the rank of matrix \(A\), using the singular value decomposition.
The rank is taken to be the number of singular values of \(A\) that are greater than the specified tolerance tol. If the second argument is omitted, it is taken to be
```

tol = max (size (A)) * sigma(1) * eps;

```
where eps is machine precision and sigma(1) is the largest singular value of \(A\).
The rank of a matrix is the number of linearly independent rows or columns and equals the dimension of the row and column space. The function orth may be used to compute an orthonormal basis of the column space.
For testing if a system \(A * x=b\) of linear equations is solvable, one can use
rank (A) == rank ([llll)

In this case, \(x=A \backslash b\) finds a particular solution \(x\). The general solution is \(x\) plus the null space of matrix \(A\). The function null may be used to compute a basis of the null space.
Example:
```

A = [lllll
4 6
7 9];
rank (A)
\# 2

```

In this example, the number of linearly independent rows is only 2 because the final row is a linear combination of the first two rows:
\[
A(3,:)=-A(1,:)+2 * A(2,:)
\]

See also: [null], page 593, [orth], page 593, [sprank], page 685, [svd], page 607, [eps], page 582.
\(c=r \operatorname{cond}(A)\)
Compute the 1 -norm estimate of the reciprocal condition number as returned by LAPACK.
If the matrix is well-conditioned then \(c\) will be near 1 and if the matrix is poorly conditioned it will be close to 0 .
The matrix \(A\) must not be sparse. If the matrix is sparse then condest (A) or rcond (full (A)) should be used instead.
See also: [cond], page 586, [condest], page 683.
trace (A)
Compute the trace of \(A\), the sum of the elements along the main diagonal.
The implementation is straightforward: sum (diag (A)).
See also: [eig], page 587.
rref (A)
rref (A, tol)
[r, k] = rref (...)
Return the reduced row echelon form of \(A\).
tol defaults to eps \(* \max (\operatorname{size}(A)) *\) norm ( \(A\), inf).
The optional return argument \(k\) contains the vector of "bound variables", which are those columns on which elimination has been performed.
\(\mathrm{n}=\mathrm{vecnorm}(\mathrm{A})\)
\(n=\operatorname{vecnorm}(A, p)\)
\(n=\operatorname{vecnorm}\) ( \(A, p\), dim)
Return the vector p-norm of the elements of array \(A\) along dimension dim.
The p-norm of a vector is defined as
\[
\|A\|_{p}=\left[\sum_{i=1}^{N}\left|A_{i}\right|^{p}\right]^{1 / p}
\]

The input \(p\) must be a positive scalar. If omitted it defaults to 2 (Euclidean norm or distance). Other special values of \(p\) are 1 (Manhattan norm, sum of absolute values) and \(\operatorname{Inf}\) (absolute value of largest element).
The input dim specifies the dimension of the array on which the function operates and must be a positive integer. If omitted the first non-singleton dimension is used.

See also: [norm], page 592.

\subsection*{18.3 Matrix Factorizations}
\(R=\operatorname{chol}(A)\)
\([R, p]=\operatorname{chol}(A)\)
\([R, p, Q]=\operatorname{chol}(A)\)
[R, \(p, Q]=\operatorname{chol}(A\), vector")
[L, ...] = chol (..., "lower")
[R, ...] = chol (..., "upper")
Compute the upper Cholesky factor, \(R\), of the real symmetric or complex Hermitian positive definite matrix \(A\).
The upper Cholesky factor \(R\) is computed by using the upper triangular part of matrix \(A\) and is defined by \(R^{T} R=A\).
Calling chol using the optional "upper" flag has the same behavior. In contrast, using the optional "lower" flag, chol returns the lower triangular factorization, computed by using the lower triangular part of matrix \(A\), such that \(L L^{T}=A\).
Called with one output argument chol fails if matrix \(A\) is not positive definite. Note that if matrix \(A\) is not real symmetric or complex Hermitian then the lower triangular part is considered to be the (complex conjugate) transpose of the upper triangular part, or vice versa, given the "lower" flag.
Called with two or more output arguments \(p\) flags whether the matrix \(A\) was positive definite and chol does not fail. A zero value of \(p\) indicates that matrix \(A\) is positive definite and \(R\) gives the factorization. Otherwise, \(p\) will have a positive value.
If called with three output arguments matrix \(A\) must be sparse and a sparsity preserving row/column permutation is applied to matrix \(A\) prior to the factorization. That is \(R\) is the factorization of \(A(Q, Q)\) such that \(R^{T} R=Q^{T} A Q\).
The sparsity preserving permutation is generally returned as a matrix. However, given the optional flag "vector", \(Q\) will be returned as a vector such that \(R^{T} R=A(Q, Q)\).
In general the lower triangular factorization is significantly faster for sparse matrices.
See also: [hess], page 597, [lu], page 597, [qr], page 599, [qz], page 602, [schur], page 604, [svd], page 607, [ichol], page 695, [cholinv], page 595, [chol2inv], page 596, [cholupdate], page 596, [cholinsert], page 596, [choldelete], page 596, [cholshift], page 597.
cholinv (A)
Compute the inverse of the symmetric positive definite matrix \(A\) using the Cholesky factorization.
See also: [chol], page 595, [chol2inv], page 596, [inv], page 590.
chol2inv (U)
Invert a symmetric, positive definite square matrix from its Cholesky decomposition, \(U\).
Note that \(U\) should be an upper-triangular matrix with positive diagonal elements. chol2inv ( \(U\) ) provides inv ( \(U^{\prime} * U\) ) but it is much faster than using inv.
See also: [chol], page 595, [cholinv], page 595, [inv], page 590.
[R1, info] = cholupdate ( \(R, \mathrm{u}, \mathrm{op}\) )
Update or downdate a Cholesky factorization.
Given an upper triangular matrix \(R\) and a column vector \(u\), attempt to determine another upper triangular matrix \(R 1\) such that
- \(R 1{ }^{*} R 1=R^{*} R+u^{*} u^{\prime}\) if op is "+"
- \(R 1{ }^{*} R 1=R^{* *} R-u^{*} u\) ' if \(o p\) is "-"

If \(o p\) is "-", info is set to
- 0 if the downdate was successful,
- 1 if \(R^{*} R-u^{*} u^{\prime}\) is not positive definite,
- 2 if \(R\) is singular.

If info is not present, an error message is printed in cases 1 and 2.
See also: [chol], page 595, [cholinsert], page 596, [choldelete], page 596, [cholshift], page 597.

R1 = cholinsert ( \(R, j, u\) )
[R1, info] = cholinsert ( \(R, j, u\) )
Update a Cholesky factorization given a row or column to insert in the original factored matrix.
Given a Cholesky factorization of a real symmetric or complex Hermitian positive definite matrix \(A=R * R, R\) upper triangular, return the Cholesky factorization of \(A 1\), where \(\mathrm{A} 1(\mathrm{p}, \mathrm{p})=\mathrm{A}, \mathrm{A} 1(:, \mathrm{j})=\mathrm{A} 1(\mathrm{j},:)^{\prime}=\mathrm{u}\) and \(\mathrm{p}=[1: \mathrm{j}-1, \mathrm{j}+1: \mathrm{n}+1]\). \(\mathrm{u}(\mathrm{j})\) should be positive.
On return, info is set to
- 0 if the insertion was successful,
- 1 if \(A 1\) is not positive definite,
- 2 if \(R\) is singular.

If info is not present, an error message is printed in cases 1 and 2.
See also: [chol], page 595, [cholupdate], page 596, [choldelete], page 596, [cholshift], page 597.
\(R 1=\) choldelete \((R, j)\)
Update a Cholesky factorization given a row or column to delete from the original factored matrix.
Given a Cholesky factorization of a real symmetric or complex Hermitian positive definite matrix \(A=R * R, R\) upper triangular, return the Cholesky factorization of \(A(p, p)\), where \(p=[1: j-1, j+1: n+1]\).

See also: [chol], page 595, [cholupdate], page 596, [cholinsert], page 596, [cholshift], page 597.

R1 = cholshift (R, i, j)
Update a Cholesky factorization given a range of columns to shift in the original factored matrix.

Given a Cholesky factorization of a real symmetric or complex Hermitian positive definite matrix \(A=R * R, R\) upper triangular, return the Cholesky factorization of \(A(\mathrm{p}, \mathrm{p})\), where p is the permutation
\(p=[1: i-1, \operatorname{shift}(i: j, 1), j+1: n]\) if \(i<j\)
or
\(p=[1: j-1, \operatorname{shift}(j: i,-1), i+1: n]\) if \(j<i\).

See also: [chol], page 595, [cholupdate], page 596, [cholinsert], page 596, [choldelete], page 596.
\(H=\) hess \((A)\)
\([P, H]=\) hess (A)
Compute the Hessenberg decomposition of the matrix \(A\).
The Hessenberg decomposition is
\[
A=P H P^{T}
\]
where \(P\) is a square unitary matrix \(\left(P^{T} P=I\right)\), and \(H\) is upper Hessenberg \(\left(H_{i, j}=\right.\) \(0, \forall i>j+1)\).

The Hessenberg decomposition is usually used as the first step in an eigenvalue computation, but has other applications as well (see Golub, Nash, and Van Loan, IEEE Transactions on Automatic Control, 1979).

See also: [eig], page 587, [chol], page 595, [lu], page 597, [qr], page 599, [qz], page 602, [schur], page 604, [svd], page 607.
\([L, U]=\operatorname{lu}(A)\)
\([L, U, P]=\operatorname{lu}(A)\)
\([L, U, P, Q]=1 u(S)\)
\([L, U, P, Q, R]=\operatorname{lu}(S)\)
[...] = lu (S, thresh)
\(y=l u(\ldots)\)
[...] = lu (..., "vector")
Compute the LU decomposition of \(A\).
If \(A\) is full then subroutines from LAPACK are used, and if \(A\) is sparse then UMFPACK is used.
The result is returned in a permuted form, according to the optional return value \(P\). For example, given the matrix \(A=[1,2 ; 3,4]\),
\[
[L, U, P]=\operatorname{lu}(A)
\]
```

returns
L =
1.00000 0.00000
0.33333 1.00000
U =
3.00000 4.00000
0.00000 0.66667
P =
0 1
1 0

```

The matrix is not required to be square.
When called with two or three output arguments and a sparse input matrix, lu does not attempt to perform sparsity preserving column permutations. Called with a fourth output argument, the sparsity preserving column transformation \(Q\) is returned, such that \(P * A * Q=L * U\). This is the preferred way to call lu with sparse input matrices.
Called with a fifth output argument and a sparse input matrix, lu attempts to use a scaling factor \(R\) on the input matrix such that \(P *(R \backslash A) * Q=L * U\). This typically leads to a sparser and more stable factorization.
An additional input argument thresh that defines the pivoting threshold can be given. thresh can be a scalar, in which case it defines the UMFPACK pivoting tolerance for both symmetric and unsymmetric cases. If thresh is a 2 -element vector, then the first element defines the pivoting tolerance for the unsymmetric UMFPACK pivoting strategy and the second for the symmetric strategy. By default, the values defined by spparms are used ([0.1, 0.001]).
Given the string argument "vector", lu returns the values of \(P\) and \(Q\) as vector values, such that for full matrix, \(A(P,:)=L * U\), and \(R(P,:) * A(:, Q)=L * U\).
With two output arguments, returns the permuted forms of the upper and lower triangular matrices, such that \(A=L * U\). With one output argument \(y\), then the matrix returned by the LAPACK routines is returned. If the input matrix is sparse then the matrix \(L\) is embedded into \(U\) to give a return value similar to the full case. For both full and sparse matrices, lu loses the permutation information.

See also: [luupdate], page 598, [ilu], page 696, [chol], page 595, [hess], page 597, [qr], page 599, [qz], page 602, [schur], page 604, [svd], page 607.
\([L, U]=\) luupdate \((L, U, x, y)\)
\([L, U, P]=\) luupdate ( \(L, U, P, x, y\) )
Given an LU factorization of a real or complex matrix \(A=L^{*} U, L\) lower unit trapezoidal and \(U\) upper trapezoidal, return the \(L U\) factorization of \(A+x^{*} y\). ., where \(x\) and \(y\) are column vectors (rank-1 update) or matrices with equal number of columns (rank-k update).

Optionally, row-pivoted updating can be used by supplying a row permutation (pivoting) matrix \(P\); in that case, an updated permutation matrix is returned. Note that if \(L, U, P\) is a pivoted LU factorization as obtained by lu :
\[
[L, U, P]=\operatorname{lu}(A) ;
\]
then a factorization of \(A+x * y\). ' can be obtained either as
```

[L1, U1] = lu (L, U,P*x, y)

```
or
\[
[L 1, U 1, P 1]=\operatorname{lu}(L, U, P, x, y)
\]

The first form uses the unpivoted algorithm, which is faster, but less stable. The second form uses a slower pivoted algorithm, which is more stable.
The matrix case is done as a sequence of rank-1 updates; thus, for large enough k , it will be both faster and more accurate to recompute the factorization from scratch.

See also: [lu], page 597, [cholupdate], page 596, [qrupdate], page 601.
\([Q, R]=\operatorname{qr}(A)\)
\([Q, R, P]=\operatorname{qr}(A)\)
\(X=\operatorname{qr}(A) \#\) non-sparse \(A\)
\(R=\operatorname{qr}(A) \#\) sparse \(A\)
\(X=\mathrm{qr}(A, B) \#\) sparse \(A\)
\([C, R]=\operatorname{qr}(A, B)\)
\([\ldots]=\operatorname{qr}(\ldots, 0)\)
[...] = qr (..., "vector")
[...] = qr (..., "matrix")
Compute the QR factorization of \(A\), using standard LAPACK subroutines.
The QR factorization is \(Q R=A\) where \(Q\) is an orthogonal matrix and \(R\) is upper triangular.
For example, given the matrix \(A=[1,2 ; 3,4]\),
\[
[Q, R]=\mathrm{qr}(A)
\]
returns
\[
\begin{aligned}
& Q= \\
& -0.31623-0.94868 \\
& \text {-0.94868 } 0.31623 \\
& R= \\
& \text {-3.16228 -4.42719 } \\
& 0.00000-0.63246
\end{aligned}
\]
which multiplied together return the original matrix
\[
\left.\begin{array}{ll}
Q * R & \\
\quad \Rightarrow & \\
& 1.0000 \\
& 2.0000 \\
& 3.0000
\end{array}\right) 4.0000
\]

If just a single return value is requested then it is either \(R\), if \(A\) is sparse, or \(X\), such that \(R=\operatorname{triu}(X)\) if \(A\) is full. (Note: unlike most commands, the single return value is not the first return value when multiple values are requested.)
If a third output \(P\) is requested, then qr calculates the permuted QR factorization \(Q R=A P\) where \(Q\) is an orthogonal matrix, \(R\) is upper triangular, and \(P\) is a permutation matrix.
If \(A\) is dense, the permuted QR factorization has the additional property that the diagonal entries of \(R\) are ordered by decreasing magnitude. In other words, abs (diag ( \(R\) ) ) will be ordered from largest to smallest.
If \(A\) is sparse, \(P\) is a fill-reducing ordering of the columns of \(A\). In that case, the diagonal entries of \(R\) are not ordered by decreasing magnitude.
For example, given the matrix \(A=[1,2 ; 3,4]\),
\[
[Q, R, P]=\operatorname{qr}(A)
\]
returns
```

Q =

```
\(-0.44721-0.89443\)
\(-0.89443 \quad 0.44721\)
\(R=\)
-4.47214 -3.13050
\(0.00000 \quad 0.44721\)
\(P=\)
01
10
If the input matrix \(A\) is sparse, the sparse QR factorization is computed by using SPQR or cxsparse (e.g., if SPQR is not available). Because the matrix \(Q\) is, in general, a full matrix, it is recommended to request only one return value \(R\). In that case, the computation avoids the construction of \(Q\) and returns a sparse \(R\) such that \(R=\operatorname{chol}\left(A^{\prime} * A\right)\).
If \(A\) is dense, an additional matrix \(B\) is supplied and two return values are requested, then qr returns \(C\), where \(C=Q^{\prime} * B\). This allows the least squares approximation of \(A \backslash B\) to be calculated as
```

[C,R] = qr (A,B)
X = R\C

```

If \(A\) is a sparse MxN matrix and an additional matrix \(B\) is supplied, one or two return values are possible. If one return value \(X\) is requested and \(M<N\), then \(X\) is the minimum 2 -norm solution of \(A \backslash B\). If \(\mathrm{M}>=\mathrm{N}, X\) is the least squares approximation of \(A \backslash B\). If two return values are requested, \(C\) and \(R\) have the same meaning as in the dense case ( \(C\) is dense and \(R\) is sparse). The version with one return parameter should be preferred because it uses less memory and can handle rank-deficient matrices better.

If the final argument is the string "vector" then \(P\) is a permutation vector (of the columns of \(A\) ) instead of a permutation matrix. In this case, the defining relationship is:
\[
Q * R=A(:, P)
\]

The default, however, is to return a permutation matrix and this may be explicitly specified by using a final argument of "matrix".
If the final argument is the scalar 0 an "economy" factorization is returned. If the original matrix \(A\) has size MxN and \(\mathrm{M}>\mathrm{N}\), then the "economy" factorization will calculate just N rows in \(R\) and N columns in \(Q\) and omit the zeros in \(R\). If \(\mathrm{M} \leq\) N , there is no difference between the economy and standard factorizations. When calculating an "economy" factorization and \(A\) is dense, the output \(P\) is always a vector rather than a matrix. If \(A\) is sparse, output \(P\) is a sparse permutation matrix.
Background: The QR factorization has applications in the solution of least squares problems
\[
\min _{x}\|A x-b\|_{2}
\]
for overdetermined systems of equations (i.e., \(A\) is a tall, thin matrix).
The permuted QR factorization \([Q, R, P]=\mathrm{qr}(A)\) allows the construction of an orthogonal basis of span (A).
See also: [chol], page 595, [hess], page 597, [lu], page 597, [qz], page 602, [schur], page 604, [svd], page 607, [qrupdate], page 601, [qrinsert], page 601, [qrdelete], page 602, [qrshift], page 602.
[Q1, R1] = qrupdate ( \(Q, R, u, v\) )
Update a QR factorization given update vectors or matrices.
Given a QR factorization of a real or complex matrix \(A=Q^{*} R, Q\) unitary and \(R\) upper trapezoidal, return the QR factorization of \(A+u^{*} v^{\prime}\), where \(u\) and \(v\) are column vectors (rank-1 update) or matrices with equal number of columns (rank-k update). Notice that the latter case is done as a sequence of rank-1 updates; thus, for k large enough, it will be both faster and more accurate to recompute the factorization from scratch.
The QR factorization supplied may be either full ( Q is square) or economized ( R is square).
See also: [qr], page 599, [qrinsert], page 601, [qrdelete], page 602, [qrshift], page 602.
[Q1, R1] = qrinsert ( \(Q, R, j, x\), orient)
Update a QR factorization given a row or column to insert in the original factored matrix.
Given a QR factorization of a real or complex matrix \(A=Q^{*} R, Q\) unitary and \(R\) upper trapezoidal, return the QR factorization of \([\mathrm{A}(:, 1: j-1) \times \mathrm{A}(:, j: \mathrm{n})]\), where \(u\) is a column vector to be inserted into \(A\) (if orient is "col"), or the QR factorization of \([\mathrm{A}(1: \mathrm{j}-1,:) ; \mathrm{x} ; \mathrm{A}(:, \mathrm{j}: \mathrm{n})]\), where x is a row vector to be inserted into \(A\) (if orient is "row").
The default value of orient is "col". If orient is "col", \(u\) may be a matrix and \(j\) an index vector resulting in the QR factorization of a matrix \(B\) such that \(\mathrm{B}(:, j)\)
gives \(u\) and \(B(:, j)=[]\) gives \(A\). Notice that the latter case is done as a sequence of k insertions; thus, for k large enough, it will be both faster and more accurate to recompute the factorization from scratch.
If orient is "col", the QR factorization supplied may be either full ( Q is square) or economized ( R is square).
If orient is "row", full factorization is needed.
See also: [qr], page 599, [qrupdate], page 601, [qrdelete], page 602, [qrshift], page 602.
[Q1, R1] = qrdelete ( \(Q, R, j\), orient)
Update a QR factorization given a row or column to delete from the original factored matrix.
Given a QR factorization of a real or complex matrix \(A=Q^{*} R, Q\) unitary and \(R\) upper trapezoidal, return the QR factorization of \([\mathrm{A}(:, 1: \mathrm{j}-1), \mathrm{U}, \mathrm{A}(:, \mathrm{j}: \mathrm{n})]\), where \(u\) is a column vector to be inserted into \(A\) (if orient is "col"), or the QR factorization of \([\mathrm{A}(1: \mathrm{j}-1,:) ; \mathrm{X} ; \mathrm{A}(:, \mathrm{j}: \mathrm{n})]\), where x is a row orient is "row"). The default value of orient is "col".
If orient is "col", \(j\) may be an index vector resulting in the QR factorization of a matrix \(B\) such that \(\mathrm{A}(:, j)=[]\) gives \(B\). Notice that the latter case is done as a sequence of \(k\) deletions; thus, for \(k\) large enough, it will be both faster and more accurate to recompute the factorization from scratch.
If orient is "col", the QR factorization supplied may be either full ( Q is square) or economized ( \(R\) is square).
If orient is "row", full factorization is needed.
See also: [qr], page 599, [qrupdate], page 601, [qrinsert], page 601, [qrshift], page 602.
[Q1, R1] = qrshift ( \(Q, R, i, j\) )
Update a QR factorization given a range of columns to shift in the original factored matrix.
Given a QR factorization of a real or complex matrix \(A=Q^{*} R, Q\) unitary and \(R\) upper trapezoidal, return the QR factorization of \(A(:, \mathrm{p})\), where p is the permutation \(p=[1: i-1, \operatorname{shift}(i: j, 1), j+1: n]\) if \(i<j\)
or
\(p=[1: j-1, \operatorname{shift}(j: i,-1), i+1: n]\) if \(j<i\).

See also: [qr], page 599, [qrupdate], page 601, [qrinsert], page 601, [qrdelete], page 602.
```

lambda = qz (A,B)

```
\([A A, B B, Q, Z, V, W, l a m b d a]=q z(A, B)\)
\([A A, B B, Z]=\mathrm{qz}(A, B, o p t)\)
[AA, BB, Z, lambda] = qz (A, B, opt)
Compute the QZ decomposition of a generalized eigenvalue problem.
The generalized eigenvalue problem is defined as
\[
A x=\lambda B x
\]

There are three calling forms of the function:
1. \(l a m b d a=q z(A, B)\)

Compute the generalized eigenvalues \(\lambda\).
2. \([A A, B B, Q, Z, V, W, l a m b d a]=q z(A, B)\)

Compute QZ decomposition, generalized eigenvectors, and generalized eigenvalues.
\[
\begin{gathered}
A A=Q^{T} A Z, B B=Q^{T} B Z \\
A V=B V \operatorname{diag}(\lambda) \\
W^{T} A=\operatorname{diag}(\lambda) W^{T} B
\end{gathered}
\]
with \(Q\) and \(Z\) orthogonal (unitary for complex case).
3. \([A A, B B, Z\{, l a m b d a\}]=q z(A, B, o p t)\)

As in form 2 above, but allows ordering of generalized eigenpairs for, e.g., solution of discrete time algebraic Riccati equations. Form 3 is not available for complex matrices, and does not compute the generalized eigenvectors \(V\), \(W\), nor the orthogonal matrix \(Q\).
opt for ordering eigenvalues of the GEP pencil. The leading block of the revised pencil contains all eigenvalues that satisfy:
\begin{tabular}{ll} 
"N" & unordered (default) \\
"S" & small: leading block has all \(|\lambda|<1\) \\
"B" & big: leading block has all \(|\lambda| \geq 1\) \\
"-" & \begin{tabular}{l} 
negative real part: leading block has all eigenvalues in \\
the open left half-plane
\end{tabular} \\
"+" & \begin{tabular}{l} 
non-negative real part: leading block has all eigenvalues \\
in the closed right half-plane
\end{tabular}
\end{tabular}

Note: qz performs permutation balancing, but not scaling (see [balance], page 585), which may be lead to less accurate results than eig. The order of output arguments was selected for compatibility with matlab.

See also: [eig], page 587, [gsvd], page 589, [balance], page 585, [chol], page 595, [hess], page 597, [lu], page 597, [qr], page 599, [qzhess], page 603, [schur], page 604.
[aa, bb, q, z] = qzhess \((A, B)\)
Compute the Hessenberg-triangular decomposition of the matrix pencil ( \(A, B\) ), returning \(\mathrm{aa}=\mathrm{q} * A * z, \mathrm{bb}=\mathrm{q} * B * z\), with \(q\) and \(z\) orthogonal.
For example:
```

[aa, bb, q, z] = qzhess ([1, 2; 3, 4], [5, 6; 7, 8])
=> aa =
-3.02244 -4.41741
0.92998 0.69749
b bb =
-8.60233 -9.99730
0.00000-0.23250
q q =
-0.58124 -0.81373
-0.81373 0.58124
\# z =
Diagonal Matrix
1 0
0}

```

The Hessenberg-triangular decomposition is the first step in Moler and Stewart's QZ decomposition algorithm.
Algorithm taken from Golub and Van Loan, Matrix Computations, 2nd edition.
See also: [lu], page 597, [chol], page 595, [hess], page 597, [qr], page 599, [qz], page 602, [schur], page 604, [svd], page 607.
\(S=\operatorname{schur}(A)\)
\(S\) = schur (A, "real")
\(S=\) schur (A, "complex")
\(S=\operatorname{schur}(A, o p t)\)
\([U, S]=\operatorname{schur}(. .\).
Compute the Schur decomposition of \(A\).
The Schur decomposition of a square matrix \(A\) is defined as
\[
S=U^{T} A U
\]
where \(U\) is a unitary matrix ( \(U^{T} U\) is identity) and \(S\) is upper triangular. The eigenvalues of \(A\) (and \(S\) ) are the diagonal elements of \(S\). If the matrix \(A\) is real, then the real Schur decomposition is computed, in which the matrix \(U\) is orthogonal and \(S\) is block upper triangular with blocks of size at most \(2 \times 2\) along the diagonal.
The default for real matrices is a real Schur decomposition. A complex decomposition may be forced by passing the flag "complex".
The eigenvalues are optionally ordered along the diagonal according to the value of opt:
opt = "a" Move eigenvalues with negative real parts to the leading block of \(S\). Mnemonic: "a" for Algebraic Riccati Equations, where this ordering is useful.
opt = "d" Move eigenvalues with magnitude less than one to the leading block of S. Mnemonic: "d" for Discrete Algebraic Riccati Equations, where this ordering is useful.
opt = "u" Unordered. No particular ordering of eigenvalues (default).

The leading \(k\) columns of \(U\) always span the \(A\)-invariant subspace corresponding to the \(k\) leading eigenvalues of \(S\).

See also: [rsf2csf], page 605, [ordschur], page 605, [ordeig], page 606, [lu], page 597, [chol], page 595, [hess], page 597, [qr], page 599, [qz], page 602, [svd], page 607, [eig], page 587.
\([U, T]=\operatorname{rsf} 2 \operatorname{csf}(U R, T R)\)
Convert a real, upper quasi-triangular Schur form \(T R\) to a complex, upper triangular Schur form \(T\).
Note that the following relations hold:
\(U R \cdot T R \cdot U R^{T}=U T U^{\dagger}\) and \(U^{\dagger} U\) is the identity matrix I.
Note also that \(U\) and \(T\) are not unique.
See also: [schur], page 604.
[UR, \(S R\) ] = ordschur (U, S, select)
Reorders the real Schur factorization \((U, S)\) obtained with the schur function, so that selected eigenvalues appear in the upper left diagonal blocks of the quasi triangular Schur matrix.
The logical vector select specifies the selected eigenvalues as they appear along \(S\) 's diagonal.
For example, given the matrix \(A=[1,2 ; 3,4]\), and its Schur decomposition
\[
[U, S]=\operatorname{schur}(A)
\]
which returns
\[
\begin{aligned}
& U= \\
& \begin{array}{rr}
U= \\
& 0.82456 \\
& -0.56577 \\
S= & \\
\\
& \\
& \\
& \\
0.0 .37228 & -1.00000 \\
0.0000 & 5.37228
\end{array}
\end{aligned}
\]

It is possible to reorder the decomposition so that the positive eigenvalue is in the upper left corner, by doing:
\[
[U, S]=\operatorname{ordschur}(U, S,[0,1])
\]

See also: [schur], page 604, [ordeig], page 606, [ordqz], page 605.
[AR, \(B R, Q R, Z R]=\) ordqz ( \(A A, B B, Q, Z\), keyword \()\)
\([A R, B R, Q R, Z R]=\) ordqz (AA, BB, Q, Z, select)
Reorder the QZ decomposition of a generalized eigenvalue problem.
The generalized eigenvalue problem is defined as
\[
A x=\lambda B x
\]

Its generalized Schur decomposition is computed using the qz algorithm:
\([A A, B B, Q, Z]=\mathrm{qz}(A, B)\)
where \(A A, B B, Q\), and \(Z\) fulfill
\[
A A=Q \cdot A \cdot Z, B B=Q \cdot B \cdot Z
\]

The ordqz function computes a unitary transformation \(Q R\) and \(Z R\) such that the order of the eigenvalue on the diagonal of \(A A\) and \(B B\) is changed. The resulting reordered matrices \(A R\) and \(B R\) fulfill:
\[
A_{R}=Q_{R} \cdot A \cdot Z_{R}, B_{R}=Q_{R} \cdot B \cdot Z_{R}
\]

The function can either be called with the keyword argument which selects the eigenvalues in the top left block of \(A R\) and \(B R\) in the following way:
```

"S", "udi"
small: leading block has all }|\lambda|<
"B", "udo"
big: leading block has all }|\lambda|\geq
"-", "lhp"
negative real part: leading block has all eigenvalues in the open left half-
plane
"+", "rhp"
non-negative real part: leading block has all eigenvalues in the closed
right half-plane

```

If a logical vector select is given instead of a keyword the ordqz function reorders all eigenvalues \(k\) to the left block for which select ( \(k\) ) is true.
Note: The keywords are compatible with the ones from qr.
See also: [eig], page 587, [ordeig], page 606, [qz], page 602, [schur], page 604, [ordschur], page 605.
```

lambda = ordeig (A)

```
lambda = ordeig (A, B)

Return the eigenvalues of quasi-triangular matrices in their order of appearance in the matrix \(A\).
The quasi-triangular matrix \(A\) is usually the result of a Schur factorization. If called with a second input \(B\) then the generalized eigenvalues of the pair \(A, B\) are returned in the order of appearance of the matrix \(A-l a m b d a * B\). The pair \(A, B\) is usually the result of a QZ decomposition.
See also: [ordschur], page 605, [ordqz], page 605, [eig], page 587, [schur], page 604, [qz], page 602.
angle \(=\operatorname{subspace}(A, B)\)
Determine the largest principal angle between two subspaces spanned by the columns of matrices \(A\) and \(B\).
```

$s=\operatorname{svd}(A)$
$[U, S, V]=\operatorname{svd}(A)$
[U, S, V] = svd (A, "econ")
$[U, S, V]=\operatorname{svd}(A, O)$

```
    Compute the singular value decomposition of \(A\).
    The singular value decomposition is defined by the relation
\[
A=U S V^{\dagger}
\]

The function svd normally returns only the vector of singular values. When called with three return values, it computes \(U, S\), and \(V\). For example,
```

    svd (hilb (3))
    ```
returns
```

        ans =
    ```
            1.4083189
            0.1223271
            0.0026873
and
[u, s, v] = svd (hilb (3))
returns
        \(u=\)
\begin{tabular}{rrr}
-0.82704 & 0.54745 & 0.12766 \\
-0.45986 & -0.52829 & -0.71375 \\
-0.32330 & -0.64901 & 0.68867
\end{tabular}
        \(\mathrm{s}=\)
            \(1.40832 \quad 0.00000 \quad 0.00000\)
            \(0.00000 \quad 0.122330 .00000\)
            \(0.00000 \quad 0.00000 \quad 0.00269\)
        v =
        \(\begin{array}{lll}-0.82704 & 0.54745 & 0.12766\end{array}\)
        -0.45986 -0.52829 -0.71375
        \(-0.32330-0.64901 \quad 0.68867\)

When given a second argument that is not 0 , svd returns an economy-sized decomposition, eliminating the unnecessary rows or columns of \(U\) or \(V\).
If the second argument is exactly 0 , then the choice of decomposition is based on the matrix \(A\). If \(A\) has more rows than columns then an economy-sized decomposition is returned, otherwise a regular decomposition is calculated.
Algorithm Notes: When calculating the full decomposition (left and right singular matrices in addition to singular values) there is a choice of two routines in LAPACK.

The default routine used by Octave is gesvd. The alternative is gesdd which is 5 X faster, but may use more memory and may be inaccurate for some input matrices. There is a third routine gejsv, suitable for better accuracy at extreme scale. See the documentation for svd_driver for more information on choosing a driver.
See also: [svd_driver], page 608, [svds], page 689, [eig], page 587, [lu], page 597, [chol], page 595, [hess], page 597, [qr], page 599, [qz], page 602.
```

val = svd_driver ()

```
old_val = svd_driver (new_val)
svd_driver (new_val, "local")

Query or set the underlying Lapack driver used by svd.
Currently recognized values are "gesdd", "gesvd", and "gejsv". The default is "gesvd".
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
Algorithm Notes: The Lapack library routines gesvd and gesdd are different only when calculating the full singular value decomposition (left and right singular matrices as well as singular values). When calculating just the singular values the following discussion is not relevant.
The newer gesdd routine is based on a Divide-and-Conquer algorithm that is 5X faster than the alternative gesvd, which is based on QR factorization. However, the new algorithm can use significantly more memory. For an MxN input matrix the memory usage is of order \(\mathrm{O}\left(\min (\mathrm{M}, \mathrm{N})^{\wedge} 2\right)\), whereas the alternative is of order \(\mathrm{O}(\max (\mathrm{M}, \mathrm{N}))\). The routine gejsv uses a preconditioned Jacobi SVD algorithm. Unlike gesvd and gesdd, in gejsv, there is no bidiagonalization step that could contaminate accuracy in some extreme cases. Also, gejsv is known to be optimally accurate in some sense. However, the speed is slower (single threaded at its core) and uses more memory \((\mathrm{O}(\min (\mathrm{M}, \mathrm{N}) ~ \sim ~ 2+\mathrm{M}+\mathrm{N})\) ).
Beyond speed and memory issues, there have been instances where some input matrices were not accurately decomposed by gesdd. See currently active bug https:// savannah.gnu.org/bugs/?55564. Until these accuracy issues are resolved in a new version of the LAPACK library, the default driver in Octave has been set to "gesvd".
See also: [svd], page 607.
[housv, beta, zer] \(=\operatorname{housh}(x, j, z)\)
Compute Householder reflection vector housv to reflect \(x\) to be the \(j\)-th column of identity, i.e.,
( \(\mathrm{I}-\mathrm{beta*housv*housv')} \mathrm{x}=\) norm \((\mathrm{x}) * e(\mathrm{j})\) if \(\mathrm{x}(\mathrm{j})<0\),
(I - beta*housv*housv') \(x=-n o r m(x) * e(j)\) if \(x(j)>=0\)
Inputs
x vector
\(j \quad\) index into vector
\(z \quad\) threshold for zero (usually should be the number 0 )

Outputs (see Golub and Van Loan):
beta If beta \(=0\), then no reflection need be applied (zer set to 0 )
housv householder vector
[u, h, nu] = krylov (A, V, k, eps1, pflg)
Construct an orthogonal basis \(u\) of a block Krylov subspace.
The block Krylov subspace has the following form:
```

[v a*v a^2*v ... a^(k+1)*v]

```

The construction is made with Householder reflections to guard against loss of orthogonality.
If \(V\) is a vector, then \(h\) contains the Hessenberg matrix such that a*u \(==u * h+r k * e k '\), in which \(\mathrm{rk}=\mathrm{a} * \mathrm{u}(:, \mathrm{k})-\mathrm{u} * \mathrm{~h}(:, \mathrm{k})\), and \(\mathrm{ek}{ }^{\prime}\) is the vector \([0,0, \ldots, 1]\) of length k. Otherwise, \(h\) is meaningless.

If \(V\) is a vector and \(k\) is greater than length (A) - 1, then \(h\) contains the Hessenberg matrix such that \(a * u==u * h\).
The value of \(n u\) is the dimension of the span of the Krylov subspace (based on eps1).
If \(b\) is a vector and \(k\) is greater than \(m-1\), then \(h\) contains the Hessenberg decomposition of \(A\).

The optional parameter eps1 is the threshold for zero. The default value is \(1 \mathrm{e}-12\).
If the optional parameter pflg is nonzero, row pivoting is used to improve numerical behavior. The default value is 0 .
Reference: A. Hodel, P. Misra, Partial Pivoting in the Computation of Krylov Subspaces of Large Sparse Systems, Proceedings of the 42nd IEEE Conference on Decision and Control, December 2003.

\subsection*{18.4 Functions of a Matrix}
expm (A)
Return the exponential of a matrix.
The matrix exponential is defined as the infinite Taylor series
\[
\exp (A)=I+A+\frac{A^{2}}{2!}+\frac{A^{3}}{3!}+\cdots
\]

However, the Taylor series is not the way to compute the matrix exponential; see Moler and Van Loan, Nineteen Dubious Ways to Compute the Exponential of a Matrix, SIAM Review, 1978. This routine uses Ward's diagonal Padé approximation method with three step preconditioning (SIAM Journal on Numerical Analysis, 1977). Diagonal Padé approximations are rational polynomials of matrices \(D_{q}(A)^{-1} N_{q}(A)\) whose Taylor series matches the first \(2 q+1\) terms of the Taylor series above; direct evaluation of the Taylor series (with the same preconditioning steps) may be desirable in lieu of the Padé approximation when \(D_{q}(A)\) is ill-conditioned.
See also: [logm], page 610, [sqrtm], page 610.
\(s=\operatorname{logm}(A)\)
s = logm (A, opt_iters)
[s, iters] = logm (...)
Compute the matrix logarithm of the square matrix \(A\).
The implementation utilizes a Padé approximant and the identity
```

logm (A) = 2^k * logm (A^(1 / 2^k))

```

The optional input opt_iters is the maximum number of square roots to compute and defaults to 100 .
The optional output iters is the number of square roots actually computed.
See also: [expm], page 609, [sqrtm], page 610.
\(s=\operatorname{sqrtm}(A)\)
[s, error_estimate] = sqrtm (A)
Compute the matrix square root of the square matrix \(A\).
Ref: N.J. Higham. A New sqrtm for matlab. Numerical Analysis Report No. 336, Manchester Centre for Computational Mathematics, Manchester, England, January 1999.

See also: [expm], page 609, [logm], page 610.
kron (A, B)
kron (A1, A2, ...)
Form the Kronecker product of two or more matrices.
This is defined block by block as
\[
x=[a(i, j) * b]
\]

For example:
```

kron (1:4, ones (3, 1))
=>}\begin{array}{lllll}{1}\&{2}\&{3}\&{4}
1

```

If there are more than two input arguments \(A 1, A 2, \ldots, A n\) the Kronecker product is computed as
```

kron (kron (A1, A2), ..., An)

```

Since the Kronecker product is associative, this is well-defined.
blkmm ( \(A, B\) )
Compute products of matrix blocks.
The blocks are given as 2 -dimensional subarrays of the arrays \(A, B\). The size of \(A\) must have the form \([m, k, \ldots]\) and size of \(B\) must be \([k, n, \ldots]\). The result is then of size \([m, n, \ldots]\) and is computed as follows:
```

for i = 1:prod (size (A) (3:end))
C(:,:,i) = A(:,:,i) * B(:,:,i)
endfor

```
\(X=\) sylvester ( \(A, B, C\) )
Solve the Sylvester equation.
The Sylvester equation is defined as:
\[
A X+X B=C
\]

The solution is computed using standard LAPACK subroutines.
For example:
```

sylvester ([1, 2; 3, 4], [5, 6; 7, 8], [9, 10; 11, 12])
=> [ 0.50000, 0.66667; 0.66667, 0.50000 ]

```

\subsection*{18.5 Specialized Solvers}
```

x = bicg (A, b)
x = bicg (A, b, tol)
x = bicg (A, b, tol, maxit)
x = bicg (A, b, tol, maxit, M)
x = bicg (A, b, tol, maxit, M1, M2)
x = bicg (A, b, tol, maxit, M, [], x0)
x = bicg (A, b, tol, maxit, M1, M2, x0)
x = bicg (A, b, tol, maxit, M, [], x0, ...)
x = bicg (A, b, tol, maxit, M1, M2, x0, ...)
[x, flag, relres, iter, resvec] = bicg (A,b,...)

```

Solve the linear system of equations \(A * x=b\) by means of the Bi-Conjugate Gradient iterative method.
The input arguments are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that Afun (x, "notransp") \(=\mathrm{A} * \mathrm{x}\) and Afun (x, "transp") \(=\mathrm{A}\) ' \(* \mathrm{x}\). Additional parameters to Afun may be passed after \(x 0\).
- \(b\) is the right-hand side vector. It must be a column vector with the same number of rows as \(A\).
- \(t o l\) is the required relative tolerance for the residual error, \(b-A * x\). The iteration stops if norm ( \(b-A * x\) ) \(\leq\) tol \(*\) norm ( \(b\) ). If tol is omitted or empty, then a tolerance of \(1 \mathrm{e}-6\) is used.
- maxit is the maximum allowed number of iterations; if maxit is omitted or empty then a value of 20 is used.
- M1, M2 are the preconditioners. The preconditioner \(M\) is given as \(M=M 1 * M 2\). Both M1 and M2 can be passed as a matrix or as a function handle or inline function g such that \(\mathrm{g}(\mathrm{x}\), "notransp") \(=M 1 \backslash \mathrm{x}\) or \(\mathrm{g}(\mathrm{x}\), "notransp") \(=M 2 \backslash \mathrm{x}\) and \(g(x\), "transp" \()=M 1^{\prime} \backslash x\) or \(g(x\), "transp" \()=M 2 ' \backslash x\). If M1 is omitted or empty, then preconditioning is not applied. The preconditioned system is theoretically equivalent to applying the bicg method to the linear system inv (M1) \(* \mathrm{~A} * \operatorname{inv}\left(\mathrm{M}_{2}\right) * \mathrm{y}=\operatorname{inv}\left(\mathrm{M1}^{\prime}\right) * \mathrm{~b}\) and \(\operatorname{inv}\left(\mathrm{M}^{\prime} \mathrm{I}^{\prime}\right) * \mathrm{~A}^{\prime} * \operatorname{inv}\left(\mathrm{M}^{\prime}\right) * z=\) inv (M2') * b and then setting \(x=\operatorname{inv}(M 2) * y\).
- \(x 0\) is the initial guess. If \(x 0\) is omitted or empty then the function sets \(x 0\) to a zero vector by default.

Any arguments which follow \(x 0\) are treated as parameters, and passed in an appropriate manner to any of the functions (Afun or Mfun) or that have been given to bicg.
The output parameters are:
- \(x\) is the computed approximation to the solution of \(A * x=b\). If the algorithm did not converge, then \(x\) is the iteration which has the minimum residual.
- flag indicates the exit status:
- 0: The algorithm converged to within the prescribed tolerance.
- 1: The algorithm did not converge and it reached the maximum number of iterations.
- 2: The preconditioner matrix is singular.
- 3: The algorithm stagnated, i.e., the absolute value of the difference between the current iteration \(x\) and the previous is less than eps \(*\) norm ( \(x, 2\) ).
- 4: The algorithm could not continue because intermediate values became too small or too large for reliable computation.
- relres is the ratio of the final residual to its initial value, measured in the Euclidean norm.
- iter is the iteration which \(x\) is computed.
- resvec is a vector containing the residual at each iteration. The total number of iterations performed is given by length (resvec) - 1 .
Consider a trivial problem with a tridiagonal matrix
```

n = 20;
A = toeplitz (sparse ([1, 1], [1, 2], [2, 1] * n ^ 2, 1, n)) + ...
toeplitz (sparse (1, 2, -1, 1, n) * n / 2, ...
sparse (1, 2, 1, 1, n) * n / 2);
b = A * ones (n, 1);
restart = 5;
[M1, M2] = ilu (A); \# in this tridiag case, it corresponds to lu (A)|
M = M1 * M2;
Afun = @(x, string) strcmp (string, "notransp") * (A * x) + ...
strcmp (string, "transp") * (A' * x);
Mfun = @(x, string) strcmp (string, "notransp") * (M \ x) + ...
strcmp (string, "transp") * (M' \ x);
M1fun = @(x, string) strcmp (string, "notransp") * (M1 \ x) + ...
strcmp (string, "transp") * (M1' \ x);
M2fun = @(x, string) strcmp (string, "notransp") * (M2 \ x) + ...
strcmp (string, "transp") * (M2' \ x);

```

Example 1: simplest usage of bicg
\(\mathrm{x}=\mathrm{bicg}(\mathrm{A}, \mathrm{b})\)
Example 2: bicg with a function that computes \(A * x\) and \(A^{\prime}{ }^{*} x\)
\(\mathrm{x}=\mathrm{bicg}\) (Afun, b, [], n)

Example 3: bicg with a preconditioner matrix \(M\)
\[
\mathrm{x}=\operatorname{bicg}(\mathrm{A}, \mathrm{~b}, 1 \mathrm{e}-6, \mathrm{n}, \mathrm{M})
\]

Example 4: bicg with a function as preconditioner
\[
\mathrm{x}=\mathrm{bicg} \text { (Afun, } \mathrm{b}, 1 \mathrm{e}-6, \mathrm{n}, \mathrm{Mfun})
\]

Example 5: bicg with preconditioner matrices M1 and M2
```

x = bicg (A, b, 1e-6, n, M1, M2)

```

Example 6: bicg with functions as preconditioners
```

x = bicg (Afun, b, 1e-6, n, M1fun, M2fun)

```

Example 7: bicg with as input a function requiring an argument
```

function $y=A p(A, x$, string, $z)$
\#\# compute $\mathrm{A}^{\wedge} \mathrm{z} * \mathrm{x}$ or ( $\left.\mathrm{A}^{\wedge} \mathrm{z}\right)^{\prime}$ * x
y = x ;
if (strcmp (string, "notransp"))
for $i=1: z$
$y=A * y ;$
endfor
elseif (strcmp (string, "transp"))
for $i=1: z$
$y=A^{\prime} * y ;$
endfor
endif
endfunction
Apfun $=@(x$, string, $p$ ) Ap (A, $x$, string, $p$ );
$\mathrm{x}=\mathrm{bicg}$ (Apfun, b, [], [], [], [], [], 2);

```

Reference:
Y. Saad, Iterative Methods for Sparse Linear Systems, Second edition, 2003, SIAM.

See also: [bicgstab], page 613, [cgs], page 615, [gmres], page 617, [pcg], page 690, [qmr], page 620, [tfqmr], page 621.
\(x=\) bicgstab (A, b, tol, maxit, M1, M2, x0, ...)
\(x=\) bicgstab (A, b, tol, maxit, \(M,[], x 0, \ldots)\)
[x, flag, relres, iter, resvec] = bicgstab ( \(A, b, \ldots\) )
Solve \(\mathrm{A} \mathrm{x}=\mathrm{b}\) using the stabilized Bi-conjugate gradient iterative method.
The input parameters are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that Afun ( x ) \(=\mathrm{A} * \mathrm{x}\). Additional parameters to Afun are passed after x0.
- \(b\) is the right hand side vector. It must be a column vector with the same number of rows as \(A\).
- \(t o l\) is the required relative tolerance for the residual error, \(b-A * x\). The iteration stops if norm ( \(b-A * x\) ) \(\leq\) tol \(*\) norm ( \(b\) ). If tol is omitted or empty, then a tolerance of \(1 \mathrm{e}-6\) is used.
- maxit the maximum number of outer iterations, if not given or set to [] the default value min (20, numel (b)) is used.
- M1, M2 are the preconditioners. The preconditioner \(M\) is given as \(M=M 1 * M 2\). Both M1 and M2 can be passed as a matrix or as a function handle or inline function \(g\) such that \(g(x)=M 1 \backslash x\) or \(g(x)=M 2 \backslash x\). The technique used is the right preconditioning, i.e., it is solved \(A * \operatorname{inv}(M) * y=b\) and then \(x=\operatorname{inv}(M)\) * \(y\).
- x0 the initial guess, if not given or set to [] the default value zeros (size (b)) is used.

The arguments which follow \(x 0\) are treated as parameters, and passed in a proper way to any of the functions ( \(A\) or \(M\) ) which are passed to bicstab.

The output parameters are:
\(-x\) is the approximation computed. If the method doesn't converge then it is the iterated with the minimum residual.
- flag indicates the exit status:
- 0: iteration converged to the within the chosen tolerance
- 1: the maximum number of iterations was reached before convergence
- 2: the preconditioner matrix is singular
- 3: the algorithm reached stagnation
- 4: the algorithm can't continue due to a division by zero
- relres is the relative residual obtained with as ( \(A * x-b\) ) / norm(b).
- iter is the (possibly half) iteration which x is computed. If it is an half iteration then it is iter +0.5
- resvec is a vector containing the residual of each half and total iteration (There are also the half iterations since \(x\) is computed in two steps at each iteration). Doing (length(resvec) - 1) / 2 is possible to see the total number of (total) iterations performed.

Let us consider a trivial problem with a tridiagonal matrix
```

n = 20;
A = toeplitz (sparse ([1, 1], [1, 2], [2, 1] * n ` 2, 1, n)) + ...
toeplitz (sparse (1, 2, -1, 1, n) * n / 2, ...
sparse (1, 2, 1, 1, n) * n / 2);
b = A * ones (n, 1);
restart = 5;
[M1, M2] = ilu (A); \# in this tridiag case, it corresponds to lu (A)
M = M1 * M2;
Afun = @(x) A * x;
Mfun = @(x) M \ x;
M1fun = @(x) M1 \ x;
M2fun = @(x) M2 \ x;
Example 1: simplest usage of bicgstab
$\mathrm{x}=\mathrm{bicgstab}(\mathrm{A}, \mathrm{b},[], \mathrm{n})$

```

Example 2: bicgstab with a function which computes \(A * x\)
\[
\mathrm{x}=\text { bicgstab (Afun, } \mathrm{b},[], \mathrm{n})
\]

Example 3: bicgstab with a preconditioner matrix \(M\)
\[
\mathrm{x}=\text { bicgstab (A, b, [], 1e-06, n, M) }
\]

Example 4: bicgstab with a function as preconditioner
```

x = bicgstab (Afun, b, 1e-6, n, Mfun)

```

Example 5: bicgstab with preconditioner matrices M1 and M2
\(\mathrm{x}=\) bicgstab (A, b, [], 1e-6, n, M1, M2)
Example 6: bicgstab with functions as preconditioners
```

x = bicgstab (Afun, b, 1e-6, n, M1fun, M2fun)

```

Example 7: bicgstab with as input a function requiring an argument
```

function y = Ap (A, x, z) \# compute A^z * x
y = x;
for i = 1:z
y = A * y;
endfor
endfunction
Apfun = @(x, string, p) Ap (A, x, string, p);
x = bicgstab (Apfun, b, [], [], [], [], [], 2);

```

Example 8: explicit example to show that bicgstab uses a right preconditioner
```

[M1, M2] = ilu (A + 0.1 * eye (n)); \# factorization of A perturbed
M = M1 * M2;

## reference solution computed by bicgstab after one iteration

[x_ref, fl] = bicgstab (A, b, [], 1, M)

## right preconditioning

[y, fl] = bicgstab (A / M, b, [], 1)
x = M \ y \# compare x and x_ref

```

Reference:
Y. Saad, Iterative Methods for Sparse Linear Systems, Second edition, 2003, SIAM

See also: [bicg], page 611, [cgs], page 615, [gmres], page 617, [pcg], page 690, [qmr], page 620, [tfqmr], page 621.
\(x=\operatorname{cgs}(A, b, t o l\), maxit, M1, M2, \(x 0, \ldots)\)
\(x=\operatorname{cgs}(A, b, t o l\), maxit, \(M,[], x 0, \ldots)\)
[x, flag, relres, iter, resvec] \(=c g s(A, b, \ldots)\)
Solve \(\mathrm{Ax}=\mathrm{b}\), where \(A\) is a square matrix, using the Conjugate Gradients Squared method.

The input arguments are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that \(\operatorname{Afun}(\mathrm{x})=\mathrm{A} * \mathrm{x}\). Additional parameters to Afun are passed after x0.
- \(b\) is the right hand side vector. It must be a column vector with same number of rows of \(A\).
- tol is the relative tolerance, if not given or set to [] the default value 1e-6 is used.
- maxit the maximum number of outer iterations, if not given or set to [] the default value min (20, numel (b)) is used.
- M1, M2 are the preconditioners. The preconditioner matrix is given as M = M1 * M2. Both M1 and M2 can be passed as a matrix or as a function handle or inline function \(g\) such that \(g(x)=M 1 \backslash x\) or \(g(x)=M 2 \backslash x\). If M1 is empty or not passed then no preconditioners are applied. The technique used is the right preconditioning, i.e., it is solved \(A * \operatorname{inv}(M) * y=\mathrm{b}\) and then \(\mathrm{x}=\operatorname{inv}(M) * \mathrm{y}\).
- x0 the initial guess, if not given or set to [] the default value zeros (size (b)) is used.

The arguments which follow \(x 0\) are treated as parameters, and passed in a proper way to any of the functions ( \(A\) or \(P\) ) which are passed to cgs.
The output parameters are:
- \(x\) is the approximation computed. If the method doesn't converge then it is the iterated with the minimum residual.
- flag indicates the exit status:
- 0: iteration converged to the within the chosen tolerance
- 1: the maximum number of iterations was reached before convergence
- 2: the preconditioner matrix is singular
- 3: the algorithm reached stagnation
- 4: the algorithm can't continue due to a division by zero
- relres is the relative residual obtained with as ( \(A * x-b\) ) / norm (b).
- iter is the iteration which \(x\) is computed.
- resvec is a vector containing the residual at each iteration. Doing length(resvec) - 1 is possible to see the total number of iterations performed.

Let us consider a trivial problem with a tridiagonal matrix
```

n = 20;
A = toeplitz (sparse ([1, 1], [1, 2], [2, 1] * n ` 2, 1, n)) + ...
toeplitz (sparse (1, 2, -1, 1, n) * n / 2, ...
sparse (1, 2, 1, 1, n) * n / 2);
b = A * ones (n, 1);
restart = 5;
[M1, M2] = ilu (A); \# in this tridiag case it corresponds to chol (A)'
M = M1 * M2;
Afun = @(x) A * x;
Mfun = @(x) M \ x;
M1fun = @(x) M1 \ x;
M2fun = @(x) M2 \ x;
Example 1: simplest usage of cgs
x = cgs (A, b, [], n)

```

Example 2: cgs with a function which computes \(A * x\)
\[
\mathrm{x}=\mathrm{cgs} \text { (Afun, b, [], n) }
\]

Example 3: cgs with a preconditioner matrix \(M\)
\[
\mathrm{x}=\operatorname{cgs}(\mathrm{A}, \mathrm{~b},[], 1 \mathrm{e}-06, \mathrm{n}, \mathrm{M})
\]

Example 4: cgs with a function as preconditioner
\(\mathrm{x}=\mathrm{cgs}\) (Afun, b, 1e-6, n, Mfun)
Example 5: cgs with preconditioner matrices M1 and M2
\(\mathrm{x}=\mathrm{cgs}(\mathrm{A}, \mathrm{b},[], 1 \mathrm{e}-6, \mathrm{n}, \mathrm{M} 1, \mathrm{M} 2)\)
Example 6: cgs with functions as preconditioners
\(\mathrm{x}=\mathrm{cgs}\) (Afun, \(\mathrm{b}, 1 \mathrm{e}-6, \mathrm{n}, \mathrm{M} 1 \mathrm{fun}, \mathrm{M} 2 \mathrm{fun})\)
Example 7: cgs with as input a function requiring an argument
```

function y = Ap (A, x, z) \# compute A^z * x
y = x;
for i = 1:z
y = A * y;
endfor
endfunction
Apfun = @(x, string, p) Ap (A, x, string, p);
x = cgs (Apfun, b, [], [], [], [], [], 2);

```

EXAMPLE 8: explicit example to show that cgs uses a right preconditioner
```

[M1, M2] = ilu (A + 0.3 * eye (n)); \# factorization of A perturbed
M = M1 * M2;

## reference solution computed by cgs after one iteration

[x_ref, fl] = cgs (A, b, [], 1, M)

## right preconditioning

[y, fl] = cgs (A / M, b, [], 1)
x = M \ y \# compare x and x_ref

```

References:
Y. Saad, Iterative Methods for Sparse Linear Systems, Second edition, 2003, SIAM

See also: [pcg], page 690, [bicgstab], page 613, [bicg], page 611, [gmres], page 617, [qmr], page 620, [tfqmr], page 621.
\(x=\) gmres (A, b, restart, tol, maxit, M1, M2, x0, ..)
\(x=\) gmres (A, b, restart, tol, maxit, M, [], x0, ..)
[x, flag, relres, iter, resvec] \(=\) gmres ( \(A, b, \ldots\) )
Solve \(\mathrm{A} \mathrm{x}=\mathrm{b}\) using the Preconditioned GMRES iterative method with restart, a.k.a. PGMRES(restart).
The input arguments are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that Afun (x) \(=\mathrm{A} * \mathrm{x}\). Additional parameters to Afun are passed after x0.
- \(\quad b\) is the right hand side vector. It must be a column vector with the same numbers of rows as \(A\).
- restart is the number of iterations before that the method restarts. If it is [] or \(\mathrm{N}=\) numel (b), then the restart is not applied.
- tol is the required relative tolerance for the preconditioned residual error, inv (M) * \((b-a * x)\). The iteration stops if norm (inv (M) * \((b-a * x)) \leq t o l *\) norm (inv \((M) * B\) ). If tol is omitted or empty, then a tolerance of \(1 \mathrm{e}-6\) is used.
- maxit is the maximum number of outer iterations, if not given or set to [], then the default value \(\min (10, N / r e s t a r t)\) is used. Note that, if restart is empty, then maxit is the maximum number of iterations. If restart and maxit are not empty, then the maximum number of iterations is restart * maxit. If both restart and maxit are empty, then the maximum number of iterations is set to \(\min (10, N)\).
- M1, M2 are the preconditioners. The preconditioner \(M\) is given as M \(=\) M1 \(*\) M2. Both M1 and M2 can be passed as a matrix, function handle, or inline function g such that \(\mathrm{g}(\mathrm{x})=\mathrm{M} 1 \backslash \mathrm{x}\) or \(\mathrm{g}(\mathrm{x})=\mathrm{M} 2 \backslash \mathrm{x}\). If M1 is [] or not given, then the preconditioner is not applied. The technique used is the left-preconditioning, i.e., it is solved \(\operatorname{inv}(M) * A * x=\operatorname{inv}(M) * b\) instead of \(A * x=b\).
- x0 is the initial guess, if not given or set to [], then the default value zeros (size (b)) is used.

The arguments which follow \(x 0\) are treated as parameters, and passed in a proper way to any of the functions ( \(A\) or \(M\) or \(M 1\) or M2) which are passed to gmres.
The outputs are:
- \(x\) the computed approximation. If the method does not converge, then it is the iterated with minimum residual.
- flag indicates the exit status:

0 : iteration converged to within the specified tolerance
1 : maximum number of iterations exceeded
2 : the preconditioner matrix is singular
3 : algorithm reached stagnation (the relative difference between two consecutive iterations is less than eps)
- relres is the value of the relative preconditioned residual of the approximation \(x\).
- iter is a vector containing the number of outer iterations and inner iterations performed to compute \(x\). That is:
- iter(1): number of outer iterations, i.e., how many times the method restarted. (if restart is empty or \(N\), then it is 1 , if not \(1 \leq i t e r(1) \leq\) maxit).
- iter(2): the number of iterations performed before the restart, i.e., the method restarts when iter(2) = restart. If restart is empty or \(N\), then 1 \(\leq\) iter (2) \(\leq\) maxit.
To be more clear, the approximation \(x\) is computed at the iteration (iter(1) - 1) * restart + iter (2). Since the output \(x\) corresponds to the minimal preconditioned residual solution, the total number of iterations that the method performed is given by length (resvec) - 1 .
- resvec is a vector containing the preconditioned relative residual at each iteration, including the 0 -th iteration norm ( \(A * x 0-b\) ).

Let us consider a trivial problem with a tridiagonal matrix
```

n = 20;
A = toeplitz (sparse ([1, 1], [1, 2], [2, 1] * n ^ 2, 1, n)) + ...
toeplitz (sparse (1, 2, -1, 1, n) * n / 2, ...
sparse (1, 2, 1, 1, n) * n / 2);
b = A * ones (n, 1);
restart = 5;
[M1, M2] = ilu (A); \# in this tridiag case, it corresponds to lu (A)
M = M1 * M2;
Afun = @(x) A * x;
Mfun = @(x) M \ x;
M1fun = @(x) M1 \ x;
M2fun = @(x) M2 \ x;
Example 1: simplest usage of gmres
x = gmres (A, b, [], [], n)

```

Example 2: gmres with a function which computes \(A * x\)
```

x = gmres (Afun, b, [], [], n)

```

Example 3: usage of gmres with the restart
```

x = gmres (A, b, restart);

```

Example 4: gmres with a preconditioner matrix \(M\) with and without restart
\(\mathrm{x}=\operatorname{gmres}(\mathrm{A}, \mathrm{b},[], 1 \mathrm{e}-06, \mathrm{n}, \mathrm{M})\)
\(\mathrm{x}=\mathrm{gmres}(\mathrm{A}, \mathrm{b}\), restart, \(1 \mathrm{e}-06, \mathrm{n}, \mathrm{M})\)
Example 5: gmres with a function as preconditioner
```

x = gmres (Afun, b, [], 1e-6, n, Mfun)

```

Example 6: gmres with preconditioner matrices M1 and M2
\(\mathrm{x}=\operatorname{gmres}(\mathrm{A}, \mathrm{b},[], 1 \mathrm{e}-6, \mathrm{n}, \mathrm{M} 1, \mathrm{M} 2)\)
Example 7: gmres with functions as preconditioners
\(\mathrm{x}=\) gmres (Afun, b, 1e-6, n, M1fun, M2fun)
Example 8: gmres with as input a function requiring an argument
```

    function y = Ap (A, x, p) # compute A^p * x
        y = x;
        for i = 1:p
            y = A * y;
        endfor
    endfunction
    Apfun = @(x, p) Ap (A, x, p);
x = gmres (Apfun, b, [], [], [], [], [], [], 2);

```

Example 9: explicit example to show that gmres uses a left preconditioner
```

[M1, M2] = ilu (A + 0.1 * eye (n)); \# factorization of A perturbed
M = M1 * M2;

## reference solution computed by gmres after two iterations

[x_ref, fl] = gmres (A, b, [], [], 1, M)

## left preconditioning

[x, fl] = gmres (M \ A, M \ b, [], [], 1)
x \# compare x and x_ref

```

Reference:
Y. Saad, Iterative Methods for Sparse Linear Systems, Second edition, 2003, SIAM

See also: [bicg], page 611, [bicgstab], page 613, [cgs], page 615, [pcg], page 690, [pcr], page 693, [qmr], page 620, [tfqmr], page 621.
\(x=q m r(A, b, r t o l\), maxit, M1, M2, x0)
\(\mathrm{x}=\mathrm{qmr}(A, b, r t o l\), maxit, \(P)\)
[x, flag, relres, iter, resvec] = qmr ( \(A, b, \ldots\) )
Solve \(\mathrm{A} \mathrm{x}=\mathrm{b}\) using the Quasi-Minimal Residual iterative method (without lookahead).
- rtol is the relative tolerance, if not given or set to [] the default value 1e-6 is used.
- maxit the maximum number of outer iterations, if not given or set to [] the default value \(\min\) ( 20 , numel (b)) is used.
- x0 the initial guess, if not given or set to [] the default value zeros (size (b)) is used.
\(A\) can be passed as a matrix or as a function handle or inline function \(f\) such that \(\mathrm{f}\left(\mathrm{x}\right.\), "notransp") \(=\mathrm{A} * \mathrm{x}\) and \(\mathrm{f}\left(\mathrm{x}\right.\), "transp") \(=\mathrm{A}^{\prime} * \mathrm{x}\).
The preconditioner \(P\) is given as \(\mathrm{P}=\mathrm{M} 1 * \mathrm{M} 2\). Both M1 and M2 can be passed as a matrix or as a function handle or inline function \(g\) such that \(g(x\), "notransp") \(=\mathrm{M} 1 \backslash\) x or \(\mathrm{g}(\mathrm{x}\), "notransp" \()=\mathrm{M} 2 \backslash \mathrm{x}\) and \(\mathrm{g}(\mathrm{x}\), "transp" \()=\mathrm{M} 1\) ' \(\backslash \mathrm{x}\) or \(\mathrm{g}(\mathrm{x}\), "transp") \(=\) M2 ' \(\backslash x\).
If called with more than one output parameter
- flag indicates the exit status:
- 0: iteration converged to the within the chosen tolerance
- 1: the maximum number of iterations was reached before convergence
- 3: the algorithm reached stagnation
(the value 2 is unused but skipped for compatibility).
- relres is the final value of the relative residual.
- iter is the number of iterations performed.
- resvec is a vector containing the residual norms at each iteration.

\section*{References:}
1. R. Freund and N. Nachtigal, QMR: a quasi-minimal residual method for nonHermitian linear systems, Numerische Mathematik, 1991, 60, pp. 315-339.
2. R. Barrett, M. Berry, T. Chan, J. Demmel, J. Donato, J. Dongarra, V. Eijkhour, R. Pozo, C. Romine, and H. van der Vorst, Templates for the solution of linear systems: Building blocks for iterative methods, SIAM, 2nd ed., 1994.

See also: [bicg], page 611, [bicgstab], page 613, [cgs], page 615, [gmres], page 617, [pcg], page 690.
\(x=\operatorname{tfqmr}(A, b, t o l, \operatorname{maxit}, M 1, M 2, x 0, \ldots)\)
\(x=\operatorname{tfqmr}(A, b, t o l, \operatorname{maxit}, M,[], x 0, \ldots)\)

Solve A \(\mathrm{x}=\mathrm{b}\) using the Transpose-Tree qmr method, based on the cgs.
The input parameters are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that Afun ( x ) \(=\mathrm{A} * \mathrm{x}\). Additional parameters to Afun are passed after x0.
- \(b\) is the right hand side vector. It must be a column vector with the same number of rows as \(A\).
- tol is the relative tolerance, if not given or set to [] the default value 1e-6 is used.
- maxit the maximum number of outer iterations, if not given or set to [] the default value min (20, numel (b)) is used. To be compatible, since the method as different behaviors in the iteration number is odd or even, is considered as iteration in tfqmr the entire odd-even cycle. That is, to make an entire iteration, the algorithm performs two sub-iterations: the odd one and the even one.
- M1, M2 are the preconditioners. The preconditioner M is given as M = M1 * M2. Both M1 and M2 can be passed as a matrix or as a function handle or inline function \(g\) such that \(g(x)=M 1 \backslash x\) or \(g(x)=M 2 \backslash x\). The technique used is the right-preconditioning, i.e., it is solved \(\mathrm{A} * \operatorname{inv}(\mathrm{M}) * \mathrm{y}=\mathrm{b}\) and then \(\mathrm{x}=\operatorname{inv}(\mathrm{M}) * \mathrm{y}\), instead of \(\mathrm{Ax}=\mathrm{b}\).
- x0 the initial guess, if not given or set to [] the default value zeros (size (b)) is used.

The arguments which follow \(x 0\) are treated as parameters, and passed in a proper way to any of the functions ( \(A\) or \(M\) ) which are passed to tfqmr.
The output parameters are:
- \(x\) is the approximation computed. If the method doesn't converge then it is the iterated with the minimum residual.
- flag indicates the exit status:
- 0: iteration converged to the within the chosen tolerance
- 1: the maximum number of iterations was reached before convergence
- 2: the preconditioner matrix is singular
- 3: the algorithm reached stagnation
- 4: the algorithm can't continue due to a division by zero
- relres is the relative residual obtained as ( \(A * x-b\) ) / norm (b).
- iter is the iteration which \(x\) is computed.
- resvec is a vector containing the residual at each iteration (including norm (b-A x 0 )). Doing length (resvec) - 1 is possible to see the total number of iterations performed.

Let us consider a trivial problem with a tridiagonal matrix
```

$\mathrm{n}=20$;
$\mathrm{A}=$ toeplitz (sparse ([1, 1], [1, 2], [2, 1] * $\left.\mathrm{n}^{-} 2,1, \mathrm{n}\right)$ ) +...
toeplitz (sparse (1, 2, $-1,1, \mathrm{n}) * \mathrm{n} / 2, \ldots$
sparse (1, 2, 1, 1, n) * n / 2);
$\mathrm{b}=\mathrm{A} *$ ones ( $\mathrm{n}, 1$ );
restart = 5;
[M1, M2] = ilu (A); \# in this tridiag case it corresponds to chol (A)
$\mathrm{M}=\mathrm{M} 1 * \mathrm{M} 2$;
Afun $=@(x)$ A * $x ;$
Mfun $=@(x)$ M $\backslash x$;
M1fun $=@(x)$ M1 $\backslash x ;$
M2fun $=@(x)$ M2 $\backslash x$;

```

Example 1: simplest usage of tfqmr
\(\mathrm{x}=\mathrm{tfqmr}(\mathrm{A}, \mathrm{b},[], \mathrm{n})\)
Example 2: tfqmr with a function which computes \(A * x\)
\(\mathrm{x}=\mathrm{tfqmr}\) (Afun, b, [], n)
Example 3: tfqmr with a preconditioner matrix \(M\)
\(\mathrm{x}=\mathrm{tfqmr}(\mathrm{A}, \mathrm{b},[], 1 \mathrm{e}-06, \mathrm{n}, \mathrm{M})\)
Example 4: tfqmr with a function as preconditioner
\(\mathrm{x}=\mathrm{tfqmr}\) (Afun, b, \(1 \mathrm{e}-6, \mathrm{n}, \mathrm{Mfun}\) )
Example 5: tfqmr with preconditioner matrices M1 and M2
\(\mathrm{x}=\mathrm{tfqmr}(\mathrm{A}, \mathrm{b},[], 1 \mathrm{e}-6, \mathrm{n}, \mathrm{M} 1, \mathrm{M} 2)\)
Example 6: tfmqr with functions as preconditioners
```

x = tfqmr (Afun, b, 1e-6, n, M1fun, M2fun)

```

Example 7: tfqmr with as input a function requiring an argument
```

function y = Ap (A, x, z) \# compute A^z * x
y = x;
for i = 1:z
y = A * y;
endfor
endfunction
Apfun = @(x, string, p) Ap (A, x, string, p);
x = tfqmr (Apfun, b, [], [], [], [], [], 2);

```

Example 8: explicit example to show that tfqmr uses a right preconditioner
```

[M1, M2] = ilu (A + 0.3 * eye (n)); \# factorization of A perturbed
M = M1 * M2;

## reference solution computed by tfqmr after one iteration

[x_ref, fl] = tfqmr (A, b, [], 1, M)

## right preconditioning

[y, fl] = tfqmr (A / M, b, [], 1)
x = M \ y \# compare x and x_ref

```

Reference:
Y. Saad, Iterative Methods for Sparse Linear Systems, Second edition, 2003, SIAM

See also: [bicg], page 611, [bicgstab], page 613, [cgs], page 615, [gmres], page 617, [pcg], page 690, [qmr], page 620, [pcr], page 693.

\section*{19 Vectorization and Faster Code Execution}

Vectorization is a programming technique that uses vector operations instead of element-byelement loop-based operations. Besides frequently producing more succinct Octave code, vectorization also allows for better optimization in the subsequent implementation. The optimizations may occur either in Octave's own Fortran, C, or C++ internal implementation, or even at a lower level depending on the compiler and external numerical libraries used to build Octave. The ultimate goal is to make use of your hardware's vector instructions if possible or to perform other optimizations in software.

Vectorization is not a concept unique to Octave, but it is particularly important because Octave is a matrix-oriented language. Vectorized Octave code will see a dramatic speed up (10X-100X) in most cases.

This chapter discusses vectorization and other techniques for writing faster code.

\subsection*{19.1 Basic Vectorization}

To a very good first approximation, the goal in vectorization is to write code that avoids loops and uses whole-array operations. As a trivial example, consider
```

for i = 1:n
for j = 1:m
c(i,j) = a(i,j) + b(i,j);
endfor
endfor

```
compared to the much simpler
```

c = a + b;

```

This isn't merely easier to write; it is also internally much easier to optimize. Octave delegates this operation to an underlying implementation which, among other optimizations, may use special vector hardware instructions or could conceivably even perform the additions in parallel. In general, if the code is vectorized, the underlying implementation has more freedom about the assumptions it can make in order to achieve faster execution.

This is especially important for loops with "cheap" bodies. Often it suffices to vectorize just the innermost loop to get acceptable performance. A general rule of thumb is that the "order" of the vectorized body should be greater or equal to the "order" of the enclosing loop.

As a less trivial example, instead of
```

for i = 1:n-1
a(i) = b(i+1) - b(i);
endfor

```
write
```

a = b(2:n) - b(1:n-1);

```

This shows an important general concept about using arrays for indexing instead of looping over an index variable. See Section 8.1 [Index Expressions], page 151. Also use
boolean indexing generously. If a condition needs to be tested, this condition can also be written as a boolean index. For instance, instead of
```

for i = 1:n
if (a(i) > 5)
a(i) -= 20
endif
endfor

```
write
```

a(a>5) -= 20;

```
which exploits the fact that a > 5 produces a boolean index.
Use elementwise vector operators whenever possible to avoid looping (operators like .* and.\(^{\wedge}\) ). See Section 8.3 [Arithmetic Ops], page 161.

Also exploit broadcasting in these elementwise operators both to avoid looping and unnecessary intermediate memory allocations. See Section 19.2 [Broadcasting], page 627.

Use built-in and library functions if possible. Built-in and compiled functions are very fast. Even with an m-file library function, chances are good that it is already optimized, or will be optimized more in a future release.

For instance, even better than
\[
a=b(2: n)-b(1: n-1) ;
\]
is
\[
\mathrm{a}=\operatorname{diff}(\mathrm{b}) ;
\]

Most Octave functions are written with vector and array arguments in mind. If you find yourself writing a loop with a very simple operation, chances are that such a function already exists. The following functions occur frequently in vectorized code:
- Index manipulation
- find
- sub2ind
- ind2sub
- sort
- unique
- lookup
- ifelse / merge
- Repetition
- repmat
- repelems
- Vectorized arithmetic
- sum
- prod
- cumsum
- cumprod
- sumsq
- diff
- dot
- cummax
- cummin
- Shape of higher dimensional arrays
- reshape
- resize
- permute
- squeeze
- deal

\subsection*{19.2 Broadcasting}

Broadcasting refers to how Octave binary operators and functions behave when their matrix or array operands or arguments differ in size. Since version 3.6.0, Octave now automatically broadcasts vectors, matrices, and arrays when using elementwise binary operators and functions. Broadly speaking, smaller arrays are "broadcast" across the larger one, until they have a compatible shape. The rule is that corresponding array dimensions must either
1. be equal, or

\section*{2. one of them must be 1 .}

In case all dimensions are equal, no broadcasting occurs and ordinary element-by-element arithmetic takes place. For arrays of higher dimensions, if the number of dimensions isn't the same, then missing trailing dimensions are treated as 1 . When one of the dimensions is 1 , the array with that singleton dimension gets copied along that dimension until it matches the dimension of the other array. For example, consider
```

x = [1 2 3;
4 6;
7 8 9];
y = [10 20 30];
x + y

```

Without broadcasting, \(\mathrm{x}+\mathrm{y}\) would be an error because the dimensions do not agree. However, with broadcasting it is as if the following operation were performed:
```

x = [1 [1 2 3
456
7 8 9];
y = [lllllll
10 20 30
10 20 30];
x + y

# 

    14 25 36
    17 28 39
    ```

That is, the smaller array of size [13] gets copied along the singleton dimension (the number of rows) until it is [3 3]. No actual copying takes place, however. The internal implementation reuses elements along the necessary dimension in order to achieve the desired effect without copying in memory.

Both arrays can be broadcast across each other, for example, all pairwise differences of the elements of a vector with itself:
\[
\begin{array}{lrrr}
y-y^{\prime} \\
\Rightarrow & 0 & 10 & 20 \\
& -10 & 0 & 10 \\
& -20 & -10 & 0
\end{array}
\]

Here the vectors of size [13] and [31] both get broadcast into matrices of size [3 3] before ordinary matrix subtraction takes place.

A special case of broadcasting that may be familiar is when all dimensions of the array being broadcast are 1, i.e., the array is a scalar. Thus for example, operations like \(x-42\) and \(\max (x, 2)\) are basic examples of broadcasting.

For a higher-dimensional example, suppose img is an RGB image of size [m \(n\) 3] and we wish to multiply each color by a different scalar. The following code accomplishes this with broadcasting,
```

img .*= permute ([0.8, 0.9, 1.2], [1, 3, 2]);

```

Note the usage of permute to match the dimensions of the \([0.8,0.9,1.2]\) vector with img.

For functions that are not written with broadcasting semantics, bsxfun can be useful for coercing them to broadcast.

\section*{bsxfun ( \(f, A, B\) )}

Apply a binary function \(f\) element-by-element to two array arguments \(A\) and \(B\), expanding singleton dimensions in either input argument as necessary.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be capable of accepting two column-vector arguments of equal length, or one column vector argument and a scalar.
The dimensions of \(A\) and \(B\) must be equal or singleton. The singleton dimensions of the arrays will be expanded to the same dimensionality as the other array.

See also: [arrayfun], page 631, [cellfun], page 633.

Broadcasting is only applied if either of the two broadcasting conditions hold. As usual, however, broadcasting does not apply when two dimensions differ and neither is 1 :
\[
\begin{aligned}
& \mathrm{x}=\left[\begin{array}{lll}
1 & 2 & 3 \\
4 & 5 & 6
\end{array}\right] ; \\
& \mathrm{y}=\left[\begin{array}{cc}
10 & 20 \\
30 & 40
\end{array}\right] ; \\
& \mathrm{x}+\mathrm{y}
\end{aligned}
\]

This will produce an error about nonconformant arguments.
Besides common arithmetic operations, several functions of two arguments also broadcast. The full list of functions and operators that broadcast is

Here is a real example of the power of broadcasting. The Floyd-Warshall algorithm is used to calculate the shortest path lengths between every pair of vertices in a graph. A naive implementation for a graph adjacency matrix of order \(n\) might look like this:
```

for k = 1:n
for i = 1:n
for j = 1:n
dist(i,j) = min (dist(i,j), dist(i,k) + dist(k,j));
endfor
endfor
endfor

```

Upon vectorizing the innermost loop, it might look like this:
```

for k = 1:n
for i = 1:n
dist(i,:) = min (dist(i,:), dist(i,k) + dist(k,:));
endfor
endfor

```

Using broadcasting in both directions, it looks like this:
```

for k = 1:n
dist = min (dist, dist(:,k) + dist(k,:));
endfor

```

The relative time performance of the three techniques for a given graph with 100 vertices is 7.3 seconds for the naive code, 87 milliseconds for the singly vectorized code, and 1.3 milliseconds for the fully broadcast code. For a graph with 1000 vertices, vectorization takes 11.7 seconds while broadcasting takes only 1.15 seconds. Therefore in general it is worth writing code with broadcasting semantics for performance.

However, beware of resorting to broadcasting if a simpler operation will suffice. For matrices a and \(b\), consider the following:
```

c = sum (permute (a, [1, 3, 2]) .* permute (b, [3, 2, 1]), 3);

```

This operation broadcasts the two matrices with permuted dimensions across each other during elementwise multiplication in order to obtain a larger 3-D array, and this array is then summed along the third dimension. A moment of thought will prove that this operation is simply the much faster ordinary matrix multiplication, \(c=a * b ;\).

A note on terminology: "broadcasting" is the term popularized by the Numpy numerical environment in the Python programming language. In other programming languages and environments, broadcasting may also be known as binary singleton expansion (BSX, in MATLAB, and the origin of the name of the bsxfun function), recycling ( R programming language), single-instruction multiple data (SIMD), or replication.

\subsection*{19.2.1 Broadcasting and Legacy Code}

The new broadcasting semantics almost never affect code that worked in previous versions of Octave. Consequently, all code inherited from matlab that worked in previous versions of Octave should still work without change in Octave. The only exception is code such as
```

try
c = a.*b;
catch
c = a.*a;
end_try_catch

```
that may have relied on matrices of different size producing an error. Because such operation is now valid Octave syntax, this will no longer produce an error. Instead, the following code should be used:
```

if (isequal (size (a), size (b)))
c = a .* b;
else
c = a .* a;
endif

```

\subsection*{19.3 Function Application}

As a general rule, functions should already be written with matrix arguments in mind and should consider whole matrix operations in a vectorized manner. Sometimes, writing functions in this way appears difficult or impossible for various reasons. For those situations, Octave provides facilities for applying a function to each element of an array, cell, or struct.
```

arrayfun (func, A)
x = arrayfun (func, A)
x = arrayfun (func, A, b, ...)
[x, y, ...] = arrayfun (func, A,...)
arrayfun (..., "UniformOutput", val)
arrayfun (..., "ErrorHandler", errfunc)

```

Execute a function on each element of an array.
This is useful for functions that do not accept array arguments. If the function does accept array arguments it is better to call the function directly.

The first input argument func can be a string, a function handle, an inline function, or an anonymous function. The input argument \(A\) can be a logic array, a numeric array, a string array, a structure array, or a cell array. By a call of the function arrayfun all elements of \(A\) are passed on to the named function func individually.

The named function can also take more than two input arguments, with the input arguments given as third input argument \(b\), fourth input argument \(c, \ldots\) If given more than one array input argument then all input arguments must have the same sizes, for example:
```

arrayfun (@atan2, [1, 0], [0, 1])
=> [ 1.57080 0.00000 ]

```

If the parameter val after a further string input argument "UniformOutput" is set true (the default), then the named function func must return a single element which then will be concatenated into the return value and is of type matrix. Otherwise, if that parameter is set to false, then the outputs are concatenated in a cell array. For example:
```

arrayfun (@(x,y) x:y, "abc", "def", "UniformOutput", false)

# 

    {
        [1,1] = abcd
        [1,2] = bcde
        [1,3] = cdef
    }
    ```

If more than one output arguments are given then the named function must return the number of return values that also are expected, for example:
```

[A, B, C] = arrayfun (@find, [10; 0], "UniformOutput", false)

# 

A =
{
[1,1] = 1
[2,1] = [] (0x0)
}
B =
{
[1,1] = 1
[2,1] = [] (0x0)
}
C =
{
[1,1] = 10
[2,1] = [] (0x0)
}

```

If the parameter errfunc after a further string input argument "ErrorHandler" is another string, a function handle, an inline function, or an anonymous function, then errfunc defines a function to call in the case that func generates an error. The definition of the function must be of the form
```

function [...] = errfunc (s,...)

```
where there is an additional input argument to errfunc relative to func, given by s. This is a structure with the elements "identifier", "message", and "index" giving, respectively, the error identifier, the error message, and the index of the array elements that caused the error. The size of the output argument of errfunc must have the same size as the output argument of func, otherwise a real error is thrown. For example:
```

function y = ferr (s, x), y = "MyString"; endfunction
arrayfun (@str2num, [1234],
"UniformOutput", false, "ErrorHandler", @ferr)

# 

    {
        [1,1] = MyString
    }
    ```

See also: [spfun], page 632, [cellfun], page 633, [structfun], page 635.
\(y=\operatorname{spfun}(f, S)\)
Compute \(\mathrm{f}(S)\) for the nonzero elements of \(S\).
The input function \(f\) is applied only to the nonzero elements of the input matrix \(S\) which is typically sparse. The function \(f\) can be passed as a string, function handle, or inline function.
The output \(y\) is a sparse matrix with the same sparsity structure as the input \(S\). spfun preserves sparsity structure which is different than simply applying the function \(f\) to the sparse matrix \(S\) when \(f(0)!=0\).

Example
Sparsity preserving spfun versus normal function application
```

S = pi * speye (2,2)
S =
Compressed Column Sparse (rows = 2, cols = 2, nnz = 2 [50%])
(1, 1) -> 3.1416
(2, 2) -> 3.1416
y = spfun (@cos, S)
y =
Compressed Column Sparse (rows = 2, cols = 2, nnz = 2 [50%])
(1, 1) -> -1
(2, 2) -> -1
y = cos (S)
y =
Compressed Column Sparse (rows = 2, cols = 2, nnz = 4 [100%])
(1, 1) -> -1
(2, 1) -> 1
(1, 2) -> 1
(2, 2) -> -1

```

See also: [arrayfun], page 631, [cellfun], page 633, [structfun], page 635.
```

cellfun (name, C)
cellfun ("size", C, k)
cellfun ("isclass", C, class)
cellfun (func, C)
cellfun (func, C, D)
[a, ...] = cellfun (...)
cellfun (..., "ErrorHandler", errfunc)
cellfun (..., "UniformOutput", val)

```

Evaluate the function named name on the elements of the cell array \(C\).
Elements in \(C\) are passed on to the named function individually. The function name can be one of the functions
isempty Return 1 for empty elements.
islogical
Return 1 for logical elements.
```

isnumeric

```

Return 1 for numeric elements.
isreal Return 1 for real elements.
length Return a vector of the lengths of cell elements.
ndims Return the number of dimensions of each element.
numel
prodofsize
Return the number of elements contained within each cell element. The number is the product of the dimensions of the object at each cell element.
size Return the size along the \(k\)-th dimension.
isclass Return 1 for elements of class.
Additionally, cellfun accepts an arbitrary function func in the form of an inline function, function handle, or the name of a function (in a character string). The function can take one or more arguments, with the inputs arguments given by \(C, D\), etc. Equally the function can return one or more output arguments. For example:
```

cellfun ("atan2", {1, 0}, {0, 1})
\# [ 1.57080 0.00000 ]

```

The number of output arguments of cellfun matches the number of output arguments of the function. The outputs of the function will be collected into the output arguments of cellfun like this:
```

function [a, b] = twoouts (x)
a = x;
b = x*x;
endfunction
[aa, bb] = cellfun (@twoouts, {1, 2, 3})
\#
aa =
123
bb}
149

```

Note that per default the output argument(s) are arrays of the same size as the input arguments. Input arguments that are singleton (1x1) cells will be automatically expanded to the size of the other arguments.
If the parameter "UniformOutput" is set to true (the default), then the function must return scalars which will be concatenated into the return array(s). If "UniformOutput" is false, the outputs are concatenated into a cell array (or cell arrays). For example:
```

cellfun ("tolower", {"Foo", "Bar", "FooBar"},
"UniformOutput", false)
=> {"foo", "bar", "foobar"}

```

Given the parameter "ErrorHandler", then errfunc defines a function to call in case func generates an error. The form of the function is
function [...] = errfunc (s, ...)
where there is an additional input argument to errfunc relative to func, given by s. This is a structure with the elements "identifier", "message", and "index" giving respectively the error identifier, the error message, and the index into the input arguments of the element that caused the error. For example:
```

function y = foo (s, x), y = NaN; endfunction
cellfun ("factorial", {-1,2}, "ErrorHandler", @foo)

# [NaN 2]

```

Use cellfun intelligently. The cellfun function is a useful tool for avoiding loops. It is often used with anonymous function handles; however, calling an anonymous function involves an overhead quite comparable to the overhead of an m-file function. Passing a handle to a built-in function is faster, because the interpreter is not involved in the internal loop. For example:
```

a = {...}
v = cellfun (@(x) det (x), a); \# compute determinants
v = cellfun (@det, a); \# faster

```

See also: [arrayfun], page 631, [structfun], page 635, [spfun], page 632.
```

structfun (func, S)

```
[A, ...] = structfun (...)
structfun (..., "ErrorHandler", errfunc)
structfun (..., "UniformOutput", val)

Evaluate the function named name on the fields of the structure \(S\). The fields of \(S\) are passed to the function func individually.
structfun accepts an arbitrary function func in the form of an inline function, function handle, or the name of a function (in a character string). In the case of a character string argument, the function must accept a single argument named \(x\), and it must return a string value. If the function returns more than one argument, they are returned as separate output variables.
If the parameter "UniformOutput" is set to true (the default), then the function must return a single element which will be concatenated into the return value. If "UniformOutput" is false, the outputs are placed into a structure with the same fieldnames as the input structure.
```

s.name1 = "John Smith";
s.name2 = "Jill Jones";
structfun (@(x) regexp (x, '(\w+)\$', "matches"){1}, s,
"UniformOutput", false)
\# scalar structure containing the fields:
name1 = Smith
name2 = Jones

```

Given the parameter "ErrorHandler", errfunc defines a function to call in case func generates an error. The form of the function is
```

function [...] = errfunc (se, ...)

```
where there is an additional input argument to errfunc relative to func, given by se. This is a structure with the elements "identifier", "message" and "index", giving respectively the error identifier, the error message, and the index into the input
arguments of the element that caused the error. For an example on how to use an error handler, see [cellfun], page 633.
See also: [cellfun], page 633, [arrayfun], page 631, [spfun], page 632.
Consistent with earlier advice, seek to use Octave built-in functions whenever possible for the best performance. This advice applies especially to the four functions above. For example, when adding two arrays together element-by-element one could use a handle to the built-in addition function @plus or define an anonymous function @(x,y) x+y. But, the anonymous function is \(60 \%\) slower than the first method. See Section 34.4.2 [Operator Overloading], page 900, for a list of basic functions which might be used in place of anonymous ones.

\subsection*{19.4 Accumulation}

Whenever it's possible to categorize according to indices the elements of an array when performing a computation, accumulation functions can be useful.
accumarray (subs, vals, sz, func, fillval, issparse)
accumarray (subs, vals, ...)
Create an array by accumulating the elements of a vector into the positions defined by their subscripts.
The subscripts are defined by the rows of the matrix subs and the values by vals. Each row of subs corresponds to one of the values in vals. If vals is a scalar, it will be used for each of the row of subs. If subs is a cell array of vectors, all vectors must be of the same length, and the subscripts in the kth vector must correspond to the \(k\) th dimension of the result.
The size of the matrix will be determined by the subscripts themselves. However, if \(s z\) is defined it determines the matrix size. The length of \(s z\) must correspond to the number of columns in subs. An exception is if subs has only one column, in which case \(s z\) may be the dimensions of a vector and the subscripts of subs are taken as the indices into it.
The default action of accumarray is to sum the elements with the same subscripts. This behavior can be modified by defining the func function. This should be a function or function handle that accepts a column vector and returns a scalar. The result of the function should not depend on the order of the subscripts.
The elements of the returned array that have no subscripts associated with them are set to zero. Defining fillval to some other value allows these values to be defined. This behavior changes, however, for certain values of func. If func is @min (respectively, @max) then the result will be filled with the minimum (respectively, maximum) integer if vals is of integral type, logical false (respectively, logical true) if vals is of logical type, zero if fillval is zero and all values are non-positive (respectively, non-negative), and NaN otherwise.
By default accumarray returns a full matrix. If issparse is logically true, then a sparse matrix is returned instead.
The following accumarray example constructs a frequency table that in the first column counts how many occurrences each number in the second column has, taken
from the vector x . Note the usage of unique for assigning to all repeated elements of \(x\) the same index (see [unique], page 791).
```

x = [91, 92, 90, 92, 90, 89, 91, 89, 90, 100, 100, 100];
[u, ~, j] = unique (x);
[accumarray(j', 1), u']
\# 2 89
3 90
2 91
2 92
3 100

```

Another example, where the result is a multi-dimensional 3-D array and the default value (zero) appears in the output:
```

accumarray ([1, 1, 1;
2, 1, 2;
2, 3, 2;
2, 1, 2;
2, 3, 2], 101:105)
=> ans(:,:,1) = [101, 0, 0; 0, 0, 0]
=> ans(:,:,2) = [0, 0, 0; 206, 0, 208]

```

The sparse option can be used as an alternative to the sparse constructor (see [sparse], page 664). Thus
sparse (i, j, sv)
can be written with accumarray as
accumarray ([i, j], sv', [], [], 0, true)
For repeated indices, sparse adds the corresponding value. To take the minimum instead, use min as an accumulator function:
```

accumarray ([i, j], sv', [], @min, 0, true)

```

The complexity of accumarray in general for the non-sparse case is generally \(\mathrm{O}(\mathrm{M}+\mathrm{N})\), where N is the number of subscripts and M is the maximum subscript (linearized in multi-dimensional case). If func is one of @sum (default), @max, @min or @(x) \{x\}, an optimized code path is used. Note that for general reduction function the interpreter overhead can play a major part and it may be more efficient to do multiple accumarray calls and compute the results in a vectorized manner.
See also: [accumdim], page 637, [unique], page 791, [sparse], page 664.
accumdim (subs, vals, dim, n, func, fillval)
Create an array by accumulating the slices of an array into the positions defined by their subscripts along a specified dimension.

The subscripts are defined by the index vector subs. The dimension is specified by dim. If not given, it defaults to the first non-singleton dimension. The length of subs must be equal to size (vals, dim).
The extent of the result matrix in the working dimension will be determined by the subscripts themselves. However, if \(n\) is defined it determines this extent.

The default action of accumdim is to sum the subarrays with the same subscripts. This behavior can be modified by defining the func function. This should be a function or function handle that accepts an array and a dimension, and reduces the array along this dimension. As a special exception, the built-in min and max functions can be used directly, and accumdim accounts for the middle empty argument that is used in their calling.
The slices of the returned array that have no subscripts associated with them are set to zero. Defining fillval to some other value allows these values to be defined.
An example of the use of accumdim is:
\[
\left.\left.\begin{array}{l}
\text { accumdim }\left([1,2,1,2,1], \begin{array}{rrr}
{[7,} & -10, & 4 ; \\
-5, & -12, & 8 ; \\
-12, & 2, & 8 ; \\
& -10, & 9,
\end{array}-3 ;\right. \\
\\
\hline-5,
\end{array}-3,-13\right]\right)
\]

See also: [accumarray], page 636.

\subsection*{19.5 Miscellaneous Techniques}

Here are some other ways of improving the execution speed of Octave programs.
- Avoid computing costly intermediate results multiple times. Octave currently does not eliminate common subexpressions. Also, certain internal computation results are cached for variables. For instance, if a matrix variable is used multiple times as an index, checking the indices (and internal conversion to integers) is only done once.
- Be aware of lazy copies (copy-on-write). When a copy of an object is created, the data is not immediately copied, but rather shared. The actual copying is postponed until the copied data needs to be modified. For example:
```

a = zeros (1000); \# create a 1000x1000 matrix
b = a; \# no copying done here
b(1) = 1; \# copying done here

```

Lazy copying applies to whole Octave objects such as matrices, cells, struct, and also individual cell or struct elements (not array elements).
Additionally, index expressions also use lazy copying when Octave can determine that the indexed portion is contiguous in memory. For example:
```

a = zeros (1000); \# create a 1000x1000 matrix
b = a(:,10:100); \# no copying done here
b = a(10:100,:); \# copying done here

```

This applies to arrays (matrices), cell arrays, and structs indexed using '()'. Index expressions generating comma-separated lists can also benefit from shallow copying in some cases. In particular, when \(a\) is a struct array, expressions like \(\{a \cdot x\},\{a(:, 2) \cdot x\}\) will use lazy copying, so that data can be shared between a struct array and a cell array.
Most indexing expressions do not live longer than their parent objects. In rare cases, however, a lazily copied slice outlasts its parent, in which case it becomes orphaned, still occupying unnecessarily more memory than needed. To provide a remedy working
in most real cases, Octave checks for orphaned lazy slices at certain situations, when a value is stored into a "permanent" location, such as a named variable or cell or struct element, and possibly economizes them. For example:
```

a = zeros (1000); \# create a 1000x1000 matrix
b = a(:,10:100); \# lazy slice
a = []; \# the original "a" array is still allocated
c{1} = b; \# b is reallocated at this point

```
- Avoid deep recursion. Function calls to m-file functions carry a relatively significant overhead, so rewriting a recursion as a loop often helps. Also, note that the maximum level of recursion is limited.
- Avoid resizing matrices unnecessarily. When building a single result matrix from a series of calculations, set the size of the result matrix first, then insert values into it. Write
```

result = zeros (big_n, big_m)
for i = over:and_over
ridx = ...
cidx = ...
result(ridx, cidx) = new_value ();
endfor

```
instead of
```

result = [];
for i = ever:and_ever
result = [ result, new_value() ];
endfor

```

Sometimes the number of items can not be computed in advance, and stack-like operations are needed. When elements are being repeatedly inserted or removed from the end of an array, Octave detects it as stack usage and attempts to use a smarter memory management strategy by pre-allocating the array in bigger chunks. This strategy is also applied to cell and struct arrays.
```

a = [];
while (condition)
a(end+1) = value; \# "push" operation
a(end) = []; \# "pop" operation
endwhile

```
- Avoid calling eval or feval excessively. Parsing input or looking up the name of a function in the symbol table are relatively expensive operations.
If you are using eval merely as an exception handling mechanism, and not because you need to execute some arbitrary text, use the try statement instead. See Section 10.9 [The try Statement], page 188.
- Use ignore_function_time_stamp when appropriate. If you are calling lots of functions, and none of them will need to change during your run, set the variable ignore_
function_time_stamp to "all". This will stop Octave from checking the time stamp of a function file to see if it has been updated while the program is being run.

\subsection*{19.6 Examples}

The following are examples of vectorization questions asked by actual users of Octave and their solutions.
- For a vector A , the following loop
```

n = length (A) - 1;
B = zeros (n, 2);
for i = 1:n
\#\# this will be two columns, the first is the difference and
\#\# the second the mean of the two elements used for the diff.
B(i,:) = [A(i+1)-A(i), (A(i+1) + A(i))/2];
endfor

```
can be turned into the following one-liner:
\[
B=[\operatorname{diff}(A)(:), 0.5 *(A(1: \text { end }-1)+A(2: \text { end }))(:)]
\]

Note the usage of colon indexing to flatten an intermediate result into a column vector. This is a common vectorization trick.

\section*{20 Nonlinear Equations}

\subsection*{20.1 Solvers}

Octave can solve sets of nonlinear equations of the form
\[
f(x)=0
\]
using the function fsolve, which is based on the minpack subroutine hybrd. This is an iterative technique so a starting point must be provided. This also has the consequence that convergence is not guaranteed even if a solution exists.
```

fsolve (fcn, x0)
fsolve (fcn, x0, options)
[x, fval, info, output, fjac] = fsolve (...)

```

Solve a system of nonlinear equations defined by the function fcn.
fun is a function handle, inline function, or string containing the name of the function to evaluate. fcn should accept a vector (array) defining the unknown variables, and return a vector of left-hand sides of the equations. Right-hand sides are defined to be zeros. In other words, this function attempts to determine a vector \(x\) such that \(f c n\) ( \(x\) ) gives (approximately) all zeros.
\(x 0\) is an initial guess for the solution. The shape of \(x 0\) is preserved in all calls to fcn, but otherwise is treated as a column vector.
options is a structure specifying additional parameters which control the algorithm. Currently, fsolve recognizes these options: "AutoScaling", "ComplexEqn", "FinDiffType", "FunValCheck", "Jacobian", "MaxFunEvals", "MaxIter", "OutputFcn", "TolFun", "TolX", "TypicalX", and "Updating".
If "AutoScaling" is "on", the variables will be automatically scaled according to the column norms of the (estimated) Jacobian. As a result, "TolFun" becomes scalingindependent. By default, this option is "off" because it may sometimes deliver unexpected (though mathematically correct) results.
If "ComplexEqn" is "on", fsolve will attempt to solve complex equations in complex variables, assuming that the equations possess a complex derivative (i.e., are holomorphic). If this is not what you want, you should unpack the real and imaginary parts of the system to get a real system.
If "Jacobian" is "on", it specifies that fcn-when called with 2 output argumentsalso returns the Jacobian matrix of right-hand sides at the requested point.
"MaxFunEvals" proscribes the maximum number of function evaluations before optimization is halted. The default value is 100 * number_of_variables, i.e., 100 * length ( \(x 0\) ). The value must be a positive integer.
If "Updating" is "on", the function will attempt to use Broyden updates to update the Jacobian, in order to reduce the number of Jacobian calculations. If your user function always calculates the Jacobian (regardless of number of output arguments) then this option provides no advantage and should be disabled.
"TolX" specifies the termination tolerance in the unknown variables, while "TolFun" is a tolerance for equations. Default is \(1 \mathrm{e}-6\) for both "TolX" and "TolFun".
For a description of the other options, see [optimset], page 753. To initialize an options structure with default values for fsolve use options = optimset ("fsolve").
The first output \(x\) is the solution while the second output fval contains the value of the function fcn evaluated at \(x\) (ideally a vector of all zeros).
The third output info reports whether the algorithm succeeded and may take one of the following values:

1 Converged to a solution point. Relative residual error is less than specified by TolFun.

2 Last relative step size was less than TolX.
3 Last relative decrease in residual was less than TolFun.
\(0 \quad\) Iteration limit (either MaxIter or MaxFunEvals) exceeded.
-1 Stopped by OutputFcn.
-2 The Jacobian became excessively small and the search stalled.
-3 The trust region radius became excessively small.
output is a structure containing runtime information about the fsolve algorithm. Fields in the structure are:
```

iterations

```

Number of iterations through loop.
```

successful

```

Number of successful iterations.
funcCount
Number of function evaluations.
The final output fjac contains the value of the Jacobian evaluated at \(x\).
Note: If you only have a single nonlinear equation of one variable, using fzero is usually a much better idea.
Note about user-supplied Jacobians: As an inherent property of the algorithm, a Jacobian is always requested for a solution vector whose residual vector is already known, and it is the last accepted successful step. Often this will be one of the last two calls, but not always. If the savings by reusing intermediate results from residual calculation in Jacobian calculation are significant, the best strategy is to employ OutputFcn: After a vector is evaluated for residuals, if OutputFcn is called with that vector, then the intermediate results should be saved for future Jacobian evaluation, and should be kept until a Jacobian evaluation is requested or until OutputFen is called with a different vector, in which case they should be dropped in favor of this most recent vector. A short example how this can be achieved follows:
```

function [fval, fjac] = user_func (x, optimvalues, state)
persistent sav = [], sav0 = [];
if (nargin == 1)

```
```

        ## evaluation call
        if (nargout == 1)
            savO.x = x; # mark saved vector
            ## calculate fval, save results to sav0.
    elseif (nargout == 2)
            ## calculate fjac using sav.
    endif
    else
\#\# outputfcn call.
if (all (x == sav0.x))
sav = sav0;
endif
\#\# maybe output iteration status, etc.
endif
endfunction

## ...

fsolve (@user_func, x0, optimset ("OutputFcn", @user_func, ...))

```

See also: [fzero], page 644, [optimset], page 753.
The following is a complete example. To solve the set of equations
\[
\begin{aligned}
-2 x^{2}+3 x y+4 \sin (y)-6 & =0 \\
3 x^{2}-2 x y^{2}+3 \cos (x)+4 & =0
\end{aligned}
\]
you first need to write a function to compute the value of the given function. For example:
```

function y = f (x)
y = zeros (2, 1);
y(1) = -2*x(1)^2 + 3*x(1)*x(2) + 4*sin(x(2)) - 6;
y(2) = 3*x(1)^2 - 2*x(1)*x(2)^2 + 3*\operatorname{cos}(x(1)) + 4;
endfunction

```

Then, call fsolve with a specified initial condition to find the roots of the system of equations. For example, given the function \(f\) defined above,
```

[x, fval, info] = fsolve (@f, [1; 2])

```
results in the solution
```

x =
0.57983
2.54621
fval =
-5.7184e-10
5.5460e-10
info = 1

```

A value of info = 1 indicates that the solution has converged.
When no Jacobian is supplied (as in the example above) it is approximated numerically. This requires more function evaluations, and hence is less efficient. In the example above we could compute the Jacobian analytically as
\[
\left[\begin{array}{cc}
\frac{\partial f_{1}}{\partial x_{1}} & \frac{\partial f_{1}}{\partial x_{2}} \\
\frac{\partial f_{2}}{\partial x_{1}} & \frac{\partial f_{2}}{\partial x_{2}}
\end{array}\right]=\left[\begin{array}{cc}
3 x_{2}-4 x_{1} & 4 \cos \left(x_{2}\right)+3 x_{1} \\
-2 x_{2}^{2}-3 \sin \left(x_{1}\right)+6 x_{1} & -4 x_{1} x_{2}
\end{array}\right]
\]
and compute it with the following Octave function
```

function $[y, j a c]=f(x)$
$\mathrm{y}=\mathrm{zeros}(2,1)$;
$\mathrm{y}(1)=-2 * \mathrm{x}(1)^{\wedge} 2+3 * \mathrm{x}(1) * \mathrm{x}(2)+4 * \sin (\mathrm{x}(2))-6$;
$y(2)=3 * x(1)^{\wedge} 2-2 * x(1) * x(2)^{\wedge} 2+3 * \cos (x(1))+4 ;$
if (nargout == 2)
jac = zeros (2, 2);
$\operatorname{jac}(1,1)=3 * x(2)-4 * x(1)$;
$\operatorname{jac}(1,2)=4 * \cos (x(2))+3 * x(1)$;
$\operatorname{jac}(2,1)=-2 * x(2) \wedge 2-3 * \sin (x(1))+6 * x(1)$;
$\operatorname{jac}(2,2)=-4 * x(1) * x(2)$;
endif
endfunction

```

The Jacobian can then be used with the following call to fsolve:
```

[x, fval, info] = fsolve (@f, [1; 2], optimset ("jacobian", "on"));

```
which gives the same solution as before.
fzero (fun, x0)
fzero (fun, x0, options)
[x, fval, info, output] = fzero (...)
Find a zero of a univariate function.
fun is a function handle, inline function, or string containing the name of the function to evaluate.
x0 should be a two-element vector specifying two points which bracket a zero. In other words, there must be a change in sign of the function between \(x 0(1)\) and \(x 0(2)\). More mathematically, the following must hold
```

sign (fun(xO(1))) * sign (fun(xO(2))) <= 0

```

If \(x 0\) is a single scalar then several nearby and distant values are probed in an attempt to obtain a valid bracketing. If this is not successful, the function fails.
options is a structure specifying additional options. Currently, fzero recognizes these options: "Display", "FunValCheck", "MaxFunEvals", "MaxIter", "OutputFcn", and "TolX".
"MaxFunEvals" proscribes the maximum number of function evaluations before the search is halted. The default value is Inf. The value must be a positive integer.
"MaxIter" proscribes the maximum number of algorithm iterations before the search is halted. The default value is Inf. The value must be a positive integer.
"TolX" specifies the termination tolerance for the solution x . The default value is eps.
For a description of the other options, see [optimset], page 753. To initialize an options structure with default values for fzero use options = optimset ("fzero").
On exit, the function returns \(x\), the approximate zero point, and fval, the function evaluated at \(x\).
The third output info reports whether the algorithm succeeded and may take one of the following values:
- 1 The algorithm converged to a solution.
- 0 Maximum number of iterations or function evaluations has been reached.
- -1 The algorithm has been terminated by a user OutputFcn.
- -5 The algorithm may have converged to a singular point.
output is a structure containing runtime information about the fzero algorithm. Fields in the structure are:
- iterations Number of iterations through loop.
- funcCount Number of function evaluations.
- algorithm The string "bisection, interpolation".
- bracketx A two-element vector with the final bracketing of the zero along the x -axis.
- brackety A two-element vector with the final bracketing of the zero along the \(y\)-axis.

See also: [optimset], page 753, [fsolve], page 641.

\subsection*{20.2 Minimizers}

Often it is useful to find the minimum value of a function rather than just the zeroes where it crosses the x -axis. fminbnd is designed for the simpler, but very common, case of a univariate function where the interval to search is bounded. For unbounded minimization of a function with potentially many variables use fminunc or fminsearch. The two functions use different internal algorithms and some knowledge of the objective function is required. For functions which can be differentiated, fminunc is appropriate. For functions with discontinuities, or for which a gradient search would fail, use fminsearch. See Chapter 25 [Optimization], page 741, for minimization with the presence of constraint functions. Note that searches can be made for maxima by simply inverting the objective function \(\left(F_{\max }=\right.\) \(-F_{m i n}\) ).
\(x=\) fminbnd (fun, \(a, b)\)
\(x=\) fminbnd (fun, \(a, b\), options)
[x, fval, info, output] = fminbnd (...)
Find a minimum point of a univariate function.
fun is a function handle, inline function, or string containing the name of the function to evaluate.
The starting interval is specified by a (left boundary) and \(b\) (right boundary). The endpoints must be finite.
options is a structure specifying additional parameters which control the algorithm. Currently, fminbnd recognizes these options: "Display", "FunValCheck", "MaxFunEvals", "MaxIter", "OutputFcn", "TolX".
"MaxFunEvals" proscribes the maximum number of function evaluations before optimization is halted. The default value is 500 . The value must be a positive integer.
"MaxIter" proscribes the maximum number of algorithm iterations before optimization is halted. The default value is 500 . The value must be a positive integer.
"TolX" specifies the termination tolerance for the solution \(x\). The default is 1e-4.
For a description of the other options, see [optimset], page 753. To initialize an options structure with default values for fminbnd use options \(=\) optimset ("fminbnd").
On exit, the function returns \(x\), the approximate minimum point, and fval, the function evaluated \(x\).
The third output info reports whether the algorithm succeeded and may take one of the following values:
- 1 The algorithm converged to a solution.
- 0 Iteration limit (either MaxIter or MaxFunEvals) exceeded.
- -1 The algorithm was terminated by a user OutputFcn.

Programming Notes: The search for a minimum is restricted to be in the finite interval bound by \(a\) and \(b\). If you have only one initial point to begin searching from then you will need to use an unconstrained minimization algorithm such as fminunc or fminsearch. fminbnd internally uses a Golden Section search strategy.
See also: [fzero], page 644, [fminunc], page 646, [fminsearch], page 647, [optimset], page 753.
fminunc ( \(f \subset n, x 0\) )
fminunc (fcn, \(x 0\), options)
[x, fval, info, output, grad, hess] = fminunc (fcn,...)
Solve an unconstrained optimization problem defined by the function fcn.
fminunc attempts to determine a vector \(x\) such that \(f c n(x)\) is a local minimum.
fun is a function handle, inline function, or string containing the name of the function to evaluate. fcn should accept a vector (array) defining the unknown variables, and return the objective function value, optionally with gradient.
\(x 0\) determines a starting guess. The shape of \(x 0\) is preserved in all calls to fcn, but otherwise is treated as a column vector.
options is a structure specifying additional parameters which control the algorithm. Currently, fminunc recognizes these options: "AutoScaling", "FinDiffType", "FunValCheck", "GradObj", "MaxFunEvals", "MaxIter", "OutputFcn", "TolFun", "TolX", "TypicalX".
If "AutoScaling" is "on", the variables will be automatically scaled according to the column norms of the (estimated) Jacobian. As a result, "TolFun" becomes scalingindependent. By default, this option is "off" because it may sometimes deliver unexpected (though mathematically correct) results.

If "GradObj" is "on", it specifies that fcn-when called with two output argumentsalso returns the Jacobian matrix of partial first derivatives at the requested point.
"MaxFunEvals" proscribes the maximum number of function evaluations before optimization is halted. The default value is 100 * number_of_variables, i.e., 100 * length ( \(x 0\) ). The value must be a positive integer.
"MaxIter" proscribes the maximum number of algorithm iterations before optimization is halted. The default value is 400 . The value must be a positive integer.
"TolX" specifies the termination tolerance for the unknown variables \(x\), while "TolFun" is a tolerance for the objective function value fval. The default is \(1 \mathrm{e}-6\) for both options.
For a description of the other options, see [optimset], page 753.
On return, \(x\) is the location of the minimum and fval contains the value of the objective function at \(x\).
info may be one of the following values:
1 Converged to a solution point. Relative gradient error is less than specified by TolFun.

2 Last relative step size was less than TolX.
3 Last relative change in function value was less than TolFun.
0 Iteration limit exceeded - either maximum number of algorithm iterations MaxIter or maximum number of function evaluations MaxFunEvals.
-1 Algorithm terminated by OutputFcn.
-3 The trust region radius became excessively small.
Optionally, fminunc can return a structure with convergence statistics (output), the output gradient (grad) at the solution \(x\), and approximate Hessian (hess) at the solution x .
Application Notes: If the objective function is a single nonlinear equation of one variable then using fminbnd is usually a better choice.
The algorithm used by fminunc is a gradient search which depends on the objective function being differentiable. If the function has discontinuities it may be better to use a derivative-free algorithm such as fminsearch.
See also: [fminbnd], page 645, [fminsearch], page 647, [optimset], page 753.
\(x=\) fminsearch (fun, \(x 0\) )
\(x=\) fminsearch (fun, \(x 0\), options)
\(x=\) fminsearch (problem)
[x, fval, exitflag, output] = fminsearch (...)
Find a value of \(x\) which minimizes the multi-variable function fun.
fun is a function handle, inline function, or string containing the name of the function to evaluate.
The search begins at the point x0 and iterates using the Nelder \& Mead Simplex algorithm (a derivative-free method). This algorithm is better-suited to functions which have discontinuities or for which a gradient-based search such as fminunc fails.

Options for the search are provided in the parameter options using the function optimset. Currently, fminsearch accepts the options: "Display", "FunValCheck","MaxFunEvals", "MaxIter", "OutputFcn", "TolFun", "TolX".
"MaxFunEvals" proscribes the maximum number of function evaluations before optimization is halted. The default value is 200 * number_of_variables, i.e., 200 * length ( \(x 0\) ). The value must be a positive integer.
"MaxIter" proscribes the maximum number of algorithm iterations before optimization is halted. The default value is \(200 *\) number_of_variables, i.e., \(200 *\) length ( \(x 0\) ). The value must be a positive integer.
For a description of the other options, see [optimset], page 753. To initialize an options structure with default values for fminsearch use options = optimset ("fminsearch").
fminsearch may also be called with a single structure argument with the following fields:
objective
The objective function.
\(x 0 \quad\) The initial point.
solver Must be set to "fminsearch".
options A structure returned from optimset or an empty matrix to indicate that defaults should be used.

The field options is optional. All others are required.
On exit, the function returns \(x\), the minimum point, and fval, the function value at the minimum.
The third output exitflag reports whether the algorithm succeeded and may take one of the following values:
1 if the algorithm converged (size of the simplex is smaller than TolX AND the step in function value between iterations is smaller than TolFun).

0 if the maximum number of iterations or the maximum number of function evaluations are exceeded.
-1 if the iteration is stopped by the "OutputFcn".
The fourth output is a structure output containing runtime about the algorithm. Fields in the structure are funcCount containing the number of function calls to fun, iterations containing the number of iteration steps, algorithm with the name of the search algorithm (always: "Nelder-Mead simplex direct search"), and message with the exit message.
Example:
```

fminsearch (@(x) (x(1)-5).^2+(x(2)-8).^4, [0;0])

```

Note: If you need to find the minimum of a single variable function it is probably better to use fminbnd.
See also: [fminbnd], page 645, [fminunc], page 646, [optimset], page 753.

The function humps is a useful function for testing zero and extrema finding functions.
\(y=\) humps ( \(x\) )
\([x, y]=\) humps ( \(x\) )
Evaluate a function with multiple minima, maxima, and zero crossings.
The output \(y\) is the evaluation of the rational function:
\[
y=-\frac{1200 x^{4}-2880 x^{3}+2036 x^{2}-348 x-88}{200 x^{4}-480 x^{3}+406 x^{2}-138 x+17}
\]
\(x\) may be a scalar, vector or array. If \(x\) is omitted, the default range [0:0.05:1] is used. When called with two output arguments, \([x, y], x\) will contain the input values, and \(y\) will contain the output from humps.
Programming Notes: humps has two local maxima located near \(x=0.300\) and 0.893 , a local minimum near \(x=0.637\), and zeros near \(x=-0.132\) and 1.300. humps is a useful function for testing algorithms which find zeros or local minima and maxima. Try demo humps to see a plot of the humps function.
See also: [fzero], page 644, [fminbnd], page 645, [fminunc], page 646, [fminsearch], page 647.

\section*{21 Diagonal and Permutation Matrices}

\subsection*{21.1 Creating and Manipulating Diagonal/Permutation Matrices}

A diagonal matrix is defined as a matrix that has zero entries outside the main diagonal; that is, \(D_{i j}=0\) if \(i \neq j\) Most often, square diagonal matrices are considered; however, the definition can equally be applied to non-square matrices, in which case we usually speak of a rectangular diagonal matrix.

A permutation matrix is defined as a square matrix that has a single element equal to unity in each row and each column; all other elements are zero. That is, there exists a permutation (vector) \(p\) such that \(P_{i j}=1\) if \(j=p_{i}\) and \(P_{i j}=0\) otherwise.

Octave provides special treatment of real and complex rectangular diagonal matrices, as well as permutation matrices. They are stored as special objects, using efficient storage and algorithms, facilitating writing both readable and efficient matrix algebra expressions in the Octave language. The special treatment may be disabled by using the functions optimize_diagonal_matrix and optimize_permutation_matrix.
val = optimize_diagonal_matrix ()
old_val = optimize_diagonal_matrix (new_val)
optimize_diagonal_matrix (new_val, "local")
Query or set whether a special space-efficient format is used for storing diagonal matrices.

The default value is true. If this option is set to false, Octave will store diagonal matrices as full matrices.

When called from inside a function with the "local" option, the setting is changed locally for the function and any subroutines it calls. The original setting is restored when exiting the function.

See also: [optimize_range], page 56, [optimize_permutation_matrix], page 651.
```

val = optimize_permutation_matrix ()
old_val = optimize_permutation_matrix (new_val)
optimize_permutation_matrix (new_val, "local")

```

Query or set whether a special space-efficient format is used for storing permutation matrices.

The default value is true. If this option is set to false, Octave will store permutation matrices as full matrices.
When called from inside a function with the "local" option, the setting is changed locally for the function and any subroutines it calls. The original setting is restored when exiting the function.

See also: [optimize_range], page 56, [optimize_diagonal_matrix], page 651.
The space savings are significant as demonstrated by the following code.
```

$\mathrm{x}=\operatorname{diag}(\operatorname{rand}(10,1))$;
xf = full (x);
sizeof (x)
$\Rightarrow 80$
sizeof (xf)
$\Rightarrow 800$

```

\subsection*{21.1.1 Creating Diagonal Matrices}

The most common and easiest way to create a diagonal matrix is using the built-in function diag. The expression diag (v), with \(v\) a vector, will create a square diagonal matrix with elements on the main diagonal given by the elements of \(v\), and size equal to the length of \(v\). diag ( \(\mathrm{v}, \mathrm{m}, \mathrm{n}\) ) can be used to construct a rectangular diagonal matrix. The result of these expressions will be a special diagonal matrix object, rather than a general matrix object.

Diagonal matrix with unit elements can be created using eye. Some other built-in functions can also return diagonal matrices. Examples include balance or inv.

Example:
```

    diag (1:4)
    
# 

Diagonal Matrix
1 0}0
0}20
0}003
0}000
diag (1:3,5,3)

# 

Diagonal Matrix

```
    100
    020
    003
    000
    000

\subsection*{21.1.2 Creating Permutation Matrices}

For creating permutation matrices, Octave does not introduce a new function, but rather overrides an existing syntax: permutation matrices can be conveniently created by indexing an identity matrix by permutation vectors. That is, if \(q\) is a permutation vector of length \(n\), the expression
```

P = eye (n) (:, q);

```
will create a permutation matrix - a special matrix object.
eye ( \(n\) ) ( \(q,:\) )
will also work (and create a row permutation matrix), as well as
```

eye (n) (q1, q2).

```

For example:
```

    eye (4) ([1,3,2,4],:)
    
# 

Permutation Matrix

```
\begin{tabular}{llll}
1 & 0 & 0 & 0 \\
0 & 0 & 1 & 0 \\
0 & 1 & 0 & 0 \\
0 & 0 & 0 & 1
\end{tabular}
    eye (4) (:, [1, 3, 2, 4])
\(\Rightarrow\)
Permutation Matrix
\begin{tabular}{llll}
1 & 0 & 0 & 0 \\
0 & 0 & 1 & 0 \\
0 & 1 & 0 & 0 \\
0 & 0 & 0 & 1
\end{tabular}

Mathematically, an identity matrix is both diagonal and permutation matrix. In Octave, eye ( \(n\) ) returns a diagonal matrix, because a matrix can only have one class. You can convert this diagonal matrix to a permutation matrix by indexing it by an identity permutation, as shown below. This is a special property of the identity matrix; indexing other diagonal matrices generally produces a full matrix.
```

    eye (3)
    
# 

Diagonal Matrix
1 0}
0 1 0
0}00
eye(3)(1:3,:)

# 

Permutation Matrix

```
\begin{tabular}{lll}
1 & 0 & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{tabular}

Some other built-in functions can also return permutation matrices. Examples include inv or lu.

\subsection*{21.1.3 Explicit and Implicit Conversions}

The diagonal and permutation matrices are special objects in their own right. A number of operations and built-in functions are defined for these matrices to use special, more efficient
code than would be used for a full matrix in the same place. Examples are given in further sections.

To facilitate smooth mixing with full matrices, backward compatibility, and compatibility with matlab, the diagonal and permutation matrices should allow any operation that works on full matrices, and will either treat it specially, or implicitly convert themselves to full matrices.

Instances include matrix indexing, except for extracting a single element or a leading submatrix, indexed assignment, or applying most mapper functions, such as exp.

An explicit conversion to a full matrix can be requested using the built-in function full. It should also be noted that the diagonal and permutation matrix objects will cache the result of the conversion after it is first requested (explicitly or implicitly), so that subsequent conversions will be very cheap.

\subsection*{21.2 Linear Algebra with Diagonal/Permutation Matrices}

As has been already said, diagonal and permutation matrices make it possible to use efficient algorithms while preserving natural linear algebra syntax. This section describes in detail the operations that are treated specially when performed on these special matrix objects.

\subsection*{21.2.1 Expressions Involving Diagonal Matrices}

Assume \(D\) is a diagonal matrix. If \(M\) is a full matrix, then \(\mathrm{D} * \mathrm{M}\) will scale the rows of \(M\). That means, if \(S=D * M\), then for each pair of indices \(\mathrm{i}, \mathrm{j}\) it holds
\[
S_{i j}=D_{i i} M_{i j}
\]

Similarly, M*D will do a column scaling.
The matrix \(D\) may also be rectangular, m-by-n where \(m!=n\). If \(m<n\), then the expression \(D * M\) is equivalent to
```

D(:,1:m) * M(1:m,:),

```
i.e., trailing \(n-m\) rows of \(M\) are ignored. If \(m>n\), then \(D * M\) is equivalent to
[D(1:n,:) * M; zeros(m-n, columns (M))],
i.e., null rows are appended to the result. The situation for right-multiplication \(\mathrm{M} * \mathrm{D}\) is analogous.

The expressions D \M and M / D perform inverse scaling. They are equivalent to solving a diagonal (or rectangular diagonal) in a least-squares minimum-norm sense. In exact arithmetic, this is equivalent to multiplying by a pseudoinverse. The pseudoinverse of a rectangular diagonal matrix is again a rectangular diagonal matrix with swapped dimensions, where each nonzero diagonal element is replaced by its reciprocal. The matrix division algorithms do, in fact, use division rather than multiplication by reciprocals for better numerical accuracy; otherwise, they honor the above definition. Note that a diagonal matrix is never truncated due to ill-conditioning; otherwise, it would not be of much use for scaling. This is typically consistent with linear algebra needs. A full matrix that only happens to be diagonal (and is thus not a special object) is of course treated normally.

Multiplication and division by diagonal matrices work efficiently also when combined with sparse matrices, i.e., \(\mathrm{D} * \mathrm{~S}\), where \(D\) is a diagonal matrix and \(S\) is a sparse matrix scales
the rows of the sparse matrix and returns a sparse matrix. The expressions \(\mathrm{S} * \mathrm{D}\), \(\mathrm{D} \backslash \mathrm{S}, \mathrm{S} / \mathrm{D}\) work analogically.

If \(D 1\) and \(D 2\) are both diagonal matrices, then the expressions
D1 + D2
D1 - D2
D1 * D2
D1 / D2
D1 \D2
again produce diagonal matrices, provided that normal dimension matching rules are obeyed. The relations used are same as described above.

Also, a diagonal matrix \(D\) can be multiplied or divided by a scalar, or raised to a scalar power if it is square, producing diagonal matrix result in all cases.

A diagonal matrix can also be transposed or conjugate-transposed, giving the expected result. Extracting a leading submatrix of a diagonal matrix, i.e., \(D(1: m, 1: n)\), will produce a diagonal matrix, other indexing expressions will implicitly convert to full matrix.

Adding a diagonal matrix to a full matrix only operates on the diagonal elements. Thus,
```

A = A + eps * eye (n)

```
is an efficient method of augmenting the diagonal of a matrix. Subtraction works analogically.

When involved in expressions with other element-by-element operators, .*, ./, . \({ }^{\text {or }} .^{\wedge}\), an implicit conversion to full matrix will take place. This is not always strictly necessary but chosen to facilitate better consistency with matlab.

\subsection*{21.2.2 Expressions Involving Permutation Matrices}

If \(P\) is a permutation matrix and \(M\) a matrix, the expression \(\mathrm{P} * \mathrm{M}\) will permute the rows of M. Similarly, \(M * P\) will yield a column permutation. Matrix division \(P \backslash M\) and \(M / P\) can be used to do inverse permutation.

The previously described syntax for creating permutation matrices can actually help an user to understand the connection between a permutation matrix and a permuting vector. Namely, the following holds, where \(I=\) eye ( \(n\) ) is an identity matrix:
```

    I(p,:) * M = (I*M) (p,:) = M(p,:)
    Similarly,
M* I(:,p) = (M*I) (:,p) = M(:,p)

```

The expressions \(I(p,:)\) and \(I(:, p)\) are permutation matrices.
A permutation matrix can be transposed (or conjugate-transposed, which is the same, because a permutation matrix is never complex), inverting the permutation, or equivalently, turning a row-permutation matrix into a column-permutation one. For permutation matrices, transpose is equivalent to inversion, thus \(\mathrm{P} \backslash \mathrm{M}\) is equivalent to \(\mathrm{P}^{\prime} * \mathrm{M}\). Transpose of a permutation matrix (or inverse) is a constant-time operation, flipping only a flag internally, and thus the choice between the two above equivalent expressions for inverse permuting is completely up to the user's taste.

Multiplication and division by permutation matrices works efficiently also when combined with sparse matrices, i.e., \(\mathrm{P} * \mathrm{~S}\), where \(P\) is a permutation matrix and \(S\) is a sparse
matrix permutes the rows of the sparse matrix and returns a sparse matrix. The expressions S*P, P\S, S/P work analogically.

Two permutation matrices can be multiplied or divided (if their sizes match), performing a composition of permutations. Also a permutation matrix can be indexed by a permutation vector (or two vectors), giving again a permutation matrix. Any other operations do not generally yield a permutation matrix and will thus trigger the implicit conversion.

\subsection*{21.3 Functions That Are Aware of These Matrices}

This section lists the built-in functions that are aware of diagonal and permutation matrices on input, or can return them as output. Passed to other functions, these matrices will in general trigger an implicit conversion. (Of course, user-defined dynamically linked functions may also work with diagonal or permutation matrices).

\subsection*{21.3.1 Diagonal Matrix Functions}
inv and pinv can be applied to a diagonal matrix, yielding again a diagonal matrix. det will use an efficient straightforward calculation when given a diagonal matrix, as well as cond. The following mapper functions can be applied to a diagonal matrix without converting it to a full one: abs, real, imag, conj, sqrt. A diagonal matrix can also be returned from the balance and svd functions. The sparse function will convert a diagonal matrix efficiently to a sparse matrix.

\subsection*{21.3.2 Permutation Matrix Functions}
inv and pinv will invert a permutation matrix, preserving its specialness. det can be applied to a permutation matrix, efficiently calculating the sign of the permutation (which is equal to the determinant).

A permutation matrix can also be returned from the built-in functions lu and \(q\) r, if a pivoted factorization is requested.

The sparse function will convert a permutation matrix efficiently to a sparse matrix. The find function will also work efficiently with a permutation matrix, making it possible to conveniently obtain the permutation indices.

\subsection*{21.4 Examples of Usage}

The following can be used to solve a linear system \(\mathrm{A} * \mathrm{x}=\mathrm{b}\) using the pivoted LU factorization:
```

[L, U, P] = lu (A); \#\# now L*U = P*A
x = U \ (L \ P) * b;

```

This is one way to normalize columns of a matrix \(X\) to unit norm:
```

s = norm (X, "columns");
X /= diag (s);

```

The same can also be accomplished with broadcasting (see Section 19.2 [Broadcasting], page 627):
```

s = norm (X, "columns");
X ./= s;

```

The following expression is a way to efficiently calculate the sign of a permutation, given by a permutation vector \(p\). It will also work in earlier versions of Octave, but slowly.
```

det (eye (length (p))(p, :))

```

Finally, here's how to solve a linear system \(\mathrm{A} * \mathrm{x}=\mathrm{b}\) with Tikhonov regularization (ridge regression) using SVD (a skeleton only):
```

m = rows (A); n = columns (A);
[U, S, V] = svd (A);

## determine the regularization factor alpha

## alpha = ...

## transform to orthogonal basis

b = U'*b;

## Use the standard formula, replacing A with S.

## S is diagonal, so the following will be very fast and accurate.

x = (S'*S + alpha^2 * eye (n)) \ (S' * b);

## transform to solution basis

x = V*x;

```

\subsection*{21.5 Differences in Treatment of Zero Elements}

Making diagonal and permutation matrices special matrix objects in their own right and the consequent usage of smarter algorithms for certain operations implies, as a side effect, small differences in treating zeros. The contents of this section apply also to sparse matrices, discussed in the following chapter. (see Chapter 22 [Sparse Matrices], page 659)

The IEEE floating point standard defines the result of the expressions \(0 *\) Inf and \(0 * \mathrm{NaN}\) as NaN. This is widely agreed to be a good compromise. Numerical software dealing with structured and sparse matrices (including Octave) however, almost always makes a distinction between a "numerical zero" and an "assumed zero". A "numerical zero" is a zero value occurring in a place where any floating-point value could occur. It is normally stored somewhere in memory as an explicit value. An "assumed zero", on the contrary, is a zero matrix element implied by the matrix structure (diagonal, triangular) or a sparsity pattern; its value is usually not stored explicitly anywhere, but is implied by the underlying data structure.

The primary distinction is that an assumed zero, when multiplied by any number, or divided by any nonzero number, yields always a zero, even when, e.g., multiplied by Inf or divided by NaN. The reason for this behavior is that the numerical multiplication is not actually performed anywhere by the underlying algorithm; the result is just assumed to be zero. Equivalently, one can say that the part of the computation involving assumed zeros is performed symbolically, not numerically.

This behavior not only facilitates the most straightforward and efficient implementation of algorithms, but also preserves certain useful invariants, like:
- scalar * diagonal matrix is a diagonal matrix
- sparse matrix / scalar preserves the sparsity pattern
- permutation matrix * matrix is equivalent to permuting rows
all of these natural mathematical truths would be invalidated by treating assumed zeros as numerical ones.

Note that matlab does not strictly follow this principle and converts assumed zeros to numerical zeros in certain cases, while not doing so in other cases. As of today, there are no intentions to mimic such behavior in Octave.

Examples of effects of assumed zeros vs. numerical zeros:
```

Inf * eye (3)

# 

    Inf 0 0
        0 Inf 0
        0 0 Inf
    Inf * speye (3)

# 

Compressed Column Sparse (rows = 3, cols = 3, nnz = 3 [33%])
(1, 1) -> Inf
(2, 2) -> Inf
(3, 3) -> Inf
Inf * full (eye (3))
=>
Inf NaN NaN
NaN Inf NaN
NaN NaN Inf
diag (1:3) * [NaN; 1; 1]

# 

    NaN
        2
        3
    sparse (1:3,1:3,1:3) * [NaN; 1; 1]

# 

    NaN
        2
        3
    [1,0,0;0,2,0;0,0,3] * [NaN; 1; 1]

# 

    NaN
    NaN
    NaN
    ```

\section*{22 Sparse Matrices}

\subsection*{22.1 Creation and Manipulation of Sparse Matrices}

The size of mathematical problems that can be treated at any particular time is generally limited by the available computing resources. Both, the speed of the computer and its available memory place limitation on the problem size.

There are many classes of mathematical problems which give rise to matrices, where a large number of the elements are zero. In this case it makes sense to have a special matrix type to handle this class of problems where only the nonzero elements of the matrix are stored. Not only does this reduce the amount of memory to store the matrix, but it also means that operations on this type of matrix can take advantage of the a priori knowledge of the positions of the nonzero elements to accelerate their calculations.

A matrix type that stores only the nonzero elements is generally called sparse. It is the purpose of this document to discuss the basics of the storage and creation of sparse matrices and the fundamental operations on them.

\subsection*{22.1.1 Storage of Sparse Matrices}

It is not strictly speaking necessary for the user to understand how sparse matrices are stored. However, such an understanding will help to get an understanding of the size of sparse matrices. Understanding the storage technique is also necessary for those users wishing to create their own oct-files.

There are many different means of storing sparse matrix data. What all of the methods have in common is that they attempt to reduce the complexity and storage given a priori knowledge of the particular class of problems that will be solved. A good summary of the available techniques for storing sparse matrix is given by Saad \({ }^{1}\). With full matrices, knowledge of the point of an element of the matrix within the matrix is implied by its position in the computers memory. However, this is not the case for sparse matrices, and so the positions of the nonzero elements of the matrix must equally be stored.

An obvious way to do this is by storing the elements of the matrix as triplets, with two elements being their position in the array (rows and column) and the third being the data itself. This is conceptually easy to grasp, but requires more storage than is strictly needed.

The storage technique used within Octave is the compressed column format. It is similar to the Yale format. \({ }^{2}\) In this format the position of each element in a row and the data are stored as previously. However, if we assume that all elements in the same column are stored adjacent in the computers memory, then we only need to store information on the number of nonzero elements in each column, rather than their positions. Thus assuming that the matrix has more nonzero elements than there are columns in the matrix, we win in terms of the amount of memory used.

In fact, the column index contains one more element than the number of columns, with the first element always being zero. The advantage of this is a simplification in the code, in

\footnotetext{
1 Y. Saad "SPARSKIT: A basic toolkit for sparse matrix computation", 1994, https://www-users.cs. umn.edu/~saad/software/SPARSKIT/paper.ps
2 https://en.wikipedia.org/wiki/Sparse_matrix\#Yale_format
}
that there is no special case for the first or last columns. A short example, demonstrating this in C is.
```

for (j = 0; j < nc; j++)
for (i = cidx(j); i < cidx(j+1); i++)
printf ("nonzero element (%i,%i) is %d\n",
ridx(i), j, data(i));

```

A clear understanding might be had by considering an example of how the above applies to an example matrix. Consider the matrix
\begin{tabular}{llll}
1 & 2 & 0 & 0 \\
0 & 0 & 0 & 3 \\
0 & 0 & 0 & 4
\end{tabular}

The nonzero elements of this matrix are
\[
\begin{array}{ll}
(1,1) & \Rightarrow 1 \\
(1,2) & \Rightarrow 2 \\
(2,4) & \Rightarrow 3 \\
(3,4) & \Rightarrow 4
\end{array}
\]

This will be stored as three vectors cidx, ridx and data, representing the column indexing, row indexing and data respectively. The contents of these three vectors for the above matrix will be
```

cidx = [0, 1, 2, 2, 4]
ridx = [0, 0, 1, 2]
data = [1, 2, 3, 4]

```

Note that this is the representation of these elements with the first row and column assumed to start at zero, while in Octave itself the row and column indexing starts at one. Thus the number of elements in the \(i\)-th column is given by cidx ( \(i+1\) ) - cidx (i).

Although Octave uses a compressed column format, it should be noted that compressed row formats are equally possible. However, in the context of mixed operations between mixed sparse and dense matrices, it makes sense that the elements of the sparse matrices are in the same order as the dense matrices. Octave stores dense matrices in column major ordering, and so sparse matrices are equally stored in this manner.

A further constraint on the sparse matrix storage used by Octave is that all elements in the rows are stored in increasing order of their row index, which makes certain operations faster. However, it imposes the need to sort the elements on the creation of sparse matrices. Having disordered elements is potentially an advantage in that it makes operations such as concatenating two sparse matrices together easier and faster, however it adds complexity and speed problems elsewhere.

\subsection*{22.1.2 Creating Sparse Matrices}

There are several means to create sparse matrix.
Returned from a function
There are many functions that directly return sparse matrices. These include speye, sprand, diag, etc.

Constructed from matrices or vectors
The function sparse allows a sparse matrix to be constructed from three vectors representing the row, column and data. Alternatively, the function spconvert uses a three column matrix format to allow easy importation of data from elsewhere.

Created and then filled
The function sparse or spalloc can be used to create an empty matrix that is then filled by the user
From a user binary program
The user can directly create the sparse matrix within an oct-file.
There are several basic functions to return specific sparse matrices. For example the sparse identity matrix, is a matrix that is often needed. It therefore has its own function to create it as speye ( \(n\) ) or speye ( \(r, c\) ), which creates an \(n\)-by- \(n\) or \(r\)-by- \(c\) sparse identity matrix.

Another typical sparse matrix that is often needed is a random distribution of random elements. The functions sprand and sprandn perform this for uniform and normal random distributions of elements. They have exactly the same calling convention, where sprand ( \(r, c, d\) ), creates an \(r\)-by- \(c\) sparse matrix with a density of filled elements of \(d\).

Other functions of interest that directly create sparse matrices, are diag or its generalization spdiags, that can take the definition of the diagonals of the matrix and create the sparse matrix that corresponds to this. For example,
```

s = diag (sparse (randn (1,n)), -1);

```
creates a sparse \((n+1)\)-by- \((n+1)\) sparse matrix with a single diagonal defined.
```

B = spdiags (A)
[B, d] = spdiags (A)
B = spdiags (A,d)
A = spdiags (v,d,A)
A = spdiags (v,d,m,n)

```

A generalization of the function diag.
Called with a single input argument, the nonzero diagonals \(d\) of \(A\) are extracted.
With two arguments the diagonals to extract are given by the vector \(d\).
The other two forms of spdiags modify the input matrix by replacing the diagonals. They use the columns of \(v\) to replace the diagonals represented by the vector \(d\). If the sparse matrix \(A\) is defined then the diagonals of this matrix are replaced. Otherwise a matrix of \(m\) by \(n\) is created with the diagonals given by the columns of \(v\).
Negative values of \(d\) represent diagonals below the main diagonal, and positive values of \(d\) diagonals above the main diagonal.
For example:
```

spdiags (reshape (1:12, 4, 3), [-1 0 1], 5, 4)
=> 5 10 0 0
1 6 11 0
0}27\quad71
0}0030
0

```

See also: [diag], page 525 .
\(s=\) speye \((m, n)\)
\(s=\) speye ( \(m\) )
\(\mathrm{s}=\) speye ( sz )
Return a sparse identity matrix of size mxn.
The implementation is significantly more efficient than sparse (eye (m)) as the full matrix is not constructed.
Called with a single argument a square matrix of size \(m\)-by-m is created. If called with a single vector argument \(s z\), this argument is taken to be the size of the matrix to create.

See also: [sparse], page 664, [spdiags], page 661, [eye], page 525.
\(r=\) spones \((S)\)
Replace the nonzero entries of \(S\) with ones.
This creates a sparse matrix with the same structure as \(S\).
See also: [sparse], page 664, [sprand], page 662, [sprandn], page 662, [sprandsym], page 663, [spfun], page 632, [spy], page 668.
sprand (m, n, d)
sprand (m, n, d, rc)
sprand (s)
Generate a sparse matrix with uniformly distributed random values.
The size of the matrix is mxn with a density of values \(d\). \(d\) must be between 0 and 1 . Values will be uniformly distributed on the interval \((0,1)\).
If called with a single matrix argument, a sparse matrix is generated with random values wherever the matrix \(s\) is nonzero.
If called with a scalar fourth argument rc, a random sparse matrix with reciprocal condition number \(r c\) is generated. If \(r c\) is a vector, then it specifies the first singular values of the generated matrix (length \((r c)<=\min (m, n)\) ).
See also: [sprandn], page 662, [sprandsym], page 663, [rand], page 530.
sprandn (m, n, d)
sprandn ( \(m, n, d, r c\) )
sprandn (s)
Generate a sparse matrix with normally distributed random values.
The size of the matrix is \(m x n\) with a density of values \(d\). \(d\) must be between 0 and 1 . Values will be normally distributed with a mean of 0 and a variance of 1 .
If called with a single matrix argument, a sparse matrix is generated with random values wherever the matrix \(s\) is nonzero.
If called with a scalar fourth argument rc, a random sparse matrix with reciprocal condition number \(r c\) is generated. If \(r c\) is a vector, then it specifies the first singular values of the generated matrix (length ( \(r c\) ) <= \(\min (m, n)\) ).
See also: [sprand], page 662, [sprandsym], page 663, [randn], page 532.
sprandsym ( \(n, d\) )
sprandsym (s)
Generate a symmetric random sparse matrix.
The size of the matrix will be nxn, with a density of values given by \(d\). \(d\) must be between 0 and 1 inclusive. Values will be normally distributed with a mean of zero and a variance of 1 .

If called with a single matrix argument, a random sparse matrix is generated wherever the matrix \(s\) is nonzero in its lower triangular part.

See also: [sprand], page 662, [sprandn], page 662, [spones], page 662, [sparse], page 664 .

The recommended way for the user to create a sparse matrix, is to create two vectors containing the row and column index of the data and a third vector of the same size containing the data to be stored. For example,
```

ri = ci = d = [];
for j = 1:c
ri = [ri; randperm(r,n)'];
ci = [ci; j*ones(n,1)];
d = [d; rand(n,1)];
endfor
s = sparse (ri, ci, d, r, c);

```
creates an \(r\)-by-c sparse matrix with a random distribution of \(n(<r)\) elements per column. The elements of the vectors do not need to be sorted in any particular order as Octave will sort them prior to storing the data. However, pre-sorting the data will make the creation of the sparse matrix faster.

The function spconvert takes a three or four column real matrix. The first two columns represent the row and column index respectively and the third and four columns, the real and imaginary parts of the sparse matrix. The matrix can contain zero elements and the elements can be sorted in any order. Adding zero elements is a convenient way to define the size of the sparse matrix. For example:
```

s = spconvert ([1 2 3 4; 1 3 4 4; 1 2 3 0]')
C Compressed Column Sparse (rows=4, cols=4, nnz=3)
(1 , 1) -> 1
(2 , 3) -> 2
(3 , 4) -> 3

```

An example of creating and filling a matrix might be
```

k = 5;
nz = r * k;
s = spalloc (r, c, nz)
for j = 1:c
idx = randperm (r);
s (:, j) = [zeros(r - k, 1); ...
rand(k, 1)] (idx);
endfor

```

It should be noted, that due to the way that the Octave assignment functions are written that the assignment will reallocate the memory used by the sparse matrix at each iteration of the above loop. Therefore the spalloc function ignores the \(n z\) argument and does not pre-assign the memory for the matrix. Therefore, it is vitally important that code using to above structure should be vectorized as much as possible to minimize the number of assignments and reduce the number of memory allocations.
\(F M=\) full (SM)
Return a full storage matrix from a sparse, diagonal, or permutation matrix, or a range.

See also: [sparse], page 664, [issparse], page 666.
\(s=s p a l l o c(m, n, n z)\)
Create an m-by-n sparse matrix with pre-allocated space for at most \(n z\) nonzero elements.

This is useful for building a matrix incrementally by a sequence of indexed assignments. Subsequent indexed assignments after spalloc will reuse the pre-allocated memory, provided they are of one of the simple forms
- \(s(I: J)=x\)
- \(s(:, I: J)=x\)
- \(s(K: L, I: J)=x\)
and that the following conditions are met:
- the assignment does not decrease nnz (S).
- after the assignment, nnz ( \(S\) ) does not exceed nz.
- no index is out of bounds.

Partial movement of data may still occur, but in general the assignment will be more memory and time efficient under these circumstances. In particular, it is possible to efficiently build a pre-allocated sparse matrix from a contiguous block of columns.
The amount of pre-allocated memory for a given matrix may be queried using the function nzmax.
Programming Note: Octave always reserves memory for at least one value, even if \(n z\) is 0 .

See also: [nzmax], page 666, [sparse], page 664.
```

s = sparse (a)
s = sparse (i, j, sv,m,n)
s = sparse (i, j, sv)
s = sparse (m,n)
s = sparse (i, j, s, m, n, "unique")
s = sparse (i, j, sv, m, n, nzmax)

```

Create a sparse matrix from a full matrix, or row, column, value triplets.
If \(a\) is a full matrix, convert it to a sparse matrix representation, removing all zero values in the process.

Given the integer index vectors \(i\) and \(j\), and a 1-by-nnz vector of real or complex values \(s v\), construct the sparse matrix \(S(i(k), j(k))=s v(k)\) with overall dimensions \(m\) and \(n\). If any of \(s v\), \(i\) or \(j\) are scalars, they are expanded to have a common size.
If \(m\) or \(n\) are not specified their values are derived from the maximum index in the vectors \(i\) and \(j\) as given by \(m=\max\) ( \(i\) ), \(n=\max\) ( \(j\) ).
Note: if multiple values are specified with the same \(i, j\) indices, the corresponding value in \(s\) will be the sum of the values at the repeated location. See [accumarray], page 636, for an example of how to produce different behavior such as taking the minimum instead.
If the option "unique" is given, and more than one value is specified at the same \(i, j\) indices, then the last specified value will be used.
sparse ( \(m, n\) ) will create an empty \(m x n\) sparse matrix and is equivalent to sparse ([], [], [], m, n)

The argument nzmax is ignored but accepted for compatibility with matlab.
Example 1 (sum at repeated indices):
```

i = [11 1 2]; j = [11 1 2]; sv = [3 4 5 5];
sparse (i, j, sv, 3, 4)

# 

    Compressed Column Sparse (rows = 3, cols = 4, nnz = 2 [17%])
        (1, 1) -> 7
        (2, 2) -> 5
    ```

Example 2 ("unique" option):
```

i = [1 1 1 2]; j = [1 1 2]; sv = [3 4 5];
sparse (i, j, sv, 3, 4, "unique")

# 

    Compressed Column Sparse (rows = 3, cols = 4, nnz = 2 [17%])
        (1, 1) -> 4
        (2, 2) -> 5
    ```

See also: [full], page 664, [accumarray], page 636, [spalloc], page 664, [spdiags], page 661, [speye], page 662, [spones], page 662, [sprand], page 662, [sprandn], page 662, [sprandsym], page 663, [spconvert], page 665, [spfun], page 632.
x = spconvert (m)
Convert a simple sparse matrix format easily generated by other programs into Octave's internal sparse format.
The input \(m\) is either a 3 or 4 column real matrix, containing the row, column, real, and imaginary parts of the elements of the sparse matrix. An element with a zero real and imaginary part can be used to force a particular matrix size.

See also: [sparse], page 664.
The above problem of memory reallocation can be avoided in oct-files. However, the construction of a sparse matrix from an oct-file is more complex than can be discussed
here. See Appendix A [External Code Interface], page 1009, for a full description of the techniques involved.

\subsection*{22.1.3 Finding Information about Sparse Matrices}

There are a number of functions that allow information concerning sparse matrices to be obtained. The most basic of these is issparse that identifies whether a particular Octave object is in fact a sparse matrix.

Another very basic function is nnz that returns the number of nonzero entries there are in a sparse matrix, while the function nzmax returns the amount of storage allocated to the sparse matrix. Note that Octave tends to crop unused memory at the first opportunity for sparse objects. There are some cases of user created sparse objects where the value returned by nzmax will not be the same as nnz, but in general they will give the same result. The function spstats returns some basic statistics on the columns of a sparse matrix including the number of elements, the mean and the variance of each column.
issparse (x)
Return true if \(x\) is a sparse matrix.
See also: [ismatrix], page 68.
\(n=\operatorname{nnz}\) (a)
Return the number of nonzero elements in a.
See also: [nzmax], page 666, [nonzeros], page 666, [find], page 513.
\(v=\) nonzeros \((A)\)
Return a column vector of the nonzero values of the matrix \(A\).
See also: [find], page 513, [nnz], page 666.
\(n=\operatorname{nzmax}(S M)\)
Return the amount of storage allocated to the sparse matrix \(S M\).
Programming Note: Octave tends to crop unused memory at the first opportunity for sparse objects. Thus, in general the value of nzmax will be the same as nnz, except for some cases of user-created sparse objects.
Also, note that Octave always reserves storage for at least one value. Thus, for empty matrices \(n n z\) will report 0 , but nzmax will report 1 .

See also: [nnz], page 666, [spalloc], page 664, [sparse], page 664.
[count, mean, var] = spstats \((S)\)
[count, mean, var] \(=\operatorname{spstats}(S, j)\)
Return the stats for the nonzero elements of the sparse matrix \(S\).
count is the number of nonzeros in each column, mean is the mean of the nonzeros in each column, and var is the variance of the nonzeros in each column.
Called with two input arguments, if \(S\) is the data and \(j\) is the bin number for the data, compute the stats for each bin. In this case, bins can contain data values of zero, whereas with spstats ( \(S\) ) the zeros may disappear.

When solving linear equations involving sparse matrices Octave determines the means to solve the equation based on the type of the matrix (see Section 22.2 [Sparse Linear Algebra], page 680). Octave probes the matrix type when the div (/) or ldiv ( \(\backslash\) ) operator is first used with the matrix and then caches the type. However the matrix_type function can be used to determine the type of the sparse matrix prior to use of the div or ldiv operators. For example,
```

a = tril (sprandn (1024, 1024, 0.02), -1) ...
+ speye (1024);
matrix_type (a);
ans = Lower

```
shows that Octave correctly determines the matrix type for lower triangular matrices. matrix_type can also be used to force the type of a matrix to be a particular type. For example:
```

a = matrix_type (tril (sprandn (1024, ...
1024, 0.02), -1) + speye (1024), "Lower");

```

This allows the cost of determining the matrix type to be avoided. However, incorrectly defining the matrix type will result in incorrect results from solutions of linear equations, and so it is entirely the responsibility of the user to correctly identify the matrix type

There are several graphical means of finding out information about sparse matrices. The first is the spy command, which displays the structure of the nonzero elements of the matrix. See Figure 22.1, for an example of the use of spy. More advanced graphical information can be obtained with the treeplot, etreeplot and gplot commands.


Figure 22.1: Structure of simple sparse matrix.
One use of sparse matrices is in graph theory, where the interconnections between nodes are represented as an adjacency matrix. That is, if the i-th node in a graph is connected to the j-th node. Then the ij -th node (and in the case of undirected graphs the ji-th node) of the sparse adjacency matrix is nonzero. If each node is then associated with a set of coordinates,
then the gplot command can be used to graphically display the interconnections between nodes.

As a trivial example of the use of gplot consider the example,
```

A = sparse ([2,6,1,3,2,4,3,5,4,6,1,5],
[1,1,2,2,3,3,4,4,5,5,6,6],1,6,6);
xy = [0,4,8,6,4,2;5,0,5,7,5,7]';
gplot (A,xy)

```
which creates an adjacency matrix A where node 1 is connected to nodes 2 and 6 , node 2 with nodes 1 and 3, etc. The coordinates of the nodes are given in the n-by- 2 matrix xy . See Figure 22.2.


Figure 22.2: Simple use of the gplot command.
The dependencies between the nodes of a Cholesky factorization can be calculated in linear time without explicitly needing to calculate the Cholesky factorization by the etree command. This command returns the elimination tree of the matrix and can be displayed graphically by the command treeplot (etree (A)) if A is symmetric or treeplot (etree ( \(\left.A+A^{\prime}\right)\) ) otherwise.
```

spy (x)
spy (..., markersize)
spy (..., line_spec)

```

Plot the sparsity pattern of the sparse matrix \(x\).
If the optional numeric argument markersize is given, it determines the size of the markers used in the plot.
If the optional string line_spec is given it is passed to plot and determines the appearance of the plot.
See also: [plot], page 316, [gplot], page 669.
\(p=\) etree \((S)\)
\(p=\) etree (S, typ)
[p, q] = etree (S, typ)
Return the elimination tree for the matrix \(S\).
By default \(S\) is assumed to be symmetric and the symmetric elimination tree is returned. The argument typ controls whether a symmetric or column elimination tree is returned. Valid values of typ are "sym" or "col", for symmetric or column elimination tree respectively.
Called with a second argument, etree also returns the postorder permutations on the tree.
etreeplot (A)
etreeplot (A, node_style, edge_style)
Plot the elimination tree of the matrix \(A\) or \(A+A^{\prime}\) if \(A\) in not symmetric.
The optional parameters node_style and edge_style define the output style.
See also: [treeplot], page 669, [gplot], page 669.
gplot ( \(A, x y\) )
gplot (A, xy, line_style)
\([x, y]=\operatorname{gplot}(A, x y)\)
Plot a graph defined by \(A\) and xy in the graph theory sense.
\(A\) is the adjacency matrix of the array to be plotted and \(x y\) is an \(n\)-by- 2 matrix containing the coordinates of the nodes of the graph.
The optional parameter line_style defines the output style for the plot. Called with no output arguments the graph is plotted directly. Otherwise, return the coordinates of the plot in \(x\) and \(y\).
See also: [treeplot], page 669, [etreeplot], page 669, [spy], page 668.
treeplot (tree)
treeplot (tree, node_style, edge_style)
Produce a graph of tree or forest.
The first argument is vector of predecessors.
The optional parameters node_style and edge_style define the output plot style.
The complexity of the algorithm is \(\mathrm{O}(\mathrm{n})\) in terms of is time and memory requirements.
See also: [etreeplot], page 669, [gplot], page 669.
treelayout (tree)
treelayout (tree, permutation)
treelayout lays out a tree or a forest.
The first argument tree is a vector of predecessors.
The parameter permutation is an optional postorder permutation.
The complexity of the algorithm is \(\mathrm{O}(\mathrm{n})\) in terms of time and memory requirements.
See also: [etreeplot], page 669, [gplot], page 669, [treeplot], page 669.

\subsection*{22.1.4 Basic Operators and Functions on Sparse Matrices}

\subsection*{22.1.4.1 Sparse Functions}

Many Octave functions have been overloaded to work with either sparse or full matrices. There is no difference in calling convention when using an overloaded function with a sparse matrix, however, there is also no access to potentially sparse-specific features. At any time the sparse matrix specific version of a function can be used by explicitly calling its function name.

The table below lists all of the sparse functions of Octave. Note that the names of the specific sparse forms of the functions are typically the same as the general versions with a \(s p\) prefix. In the table below, and in the rest of this article, the specific sparse versions of functions are used.

Generate sparse matrices:
spalloc, spdiags, speye, sprand, sprandn, sprandsym
Sparse matrix conversion:
full, sparse, spconvert
Manipulate sparse matrices
issparse, nnz, nonzeros, nzmax, spfun, spones, spy
Graph Theory:
etree, etreeplot, gplot, treeplot
Sparse matrix reordering:
amd, ccolamd, colamd, colperm, csymamd, dmperm, symamd, randperm, symrcm

Linear algebra:
condest, eigs, matrix_type, normest, normest1, sprank, spaugment, svds
Iterative techniques:
ichol, ilu, pcg, pcr
Miscellaneous:
spparms, symbfact, spstats
In addition all of the standard Octave mapper functions (i.e., basic math functions that take a single argument) such as abs, etc. can accept sparse matrices. The reader is referred to the documentation supplied with these functions within Octave itself for further details.

\subsection*{22.1.4.2 Return Types of Operators and Functions}

The two basic reasons to use sparse matrices are to reduce the memory usage and to not have to do calculations on zero elements. The two are closely related in that the computation time on a sparse matrix operator or function is roughly linear with the number of nonzero elements.

Therefore, there is a certain density of nonzero elements of a matrix where it no longer makes sense to store it as a sparse matrix, but rather as a full matrix. For this reason operators and functions that have a high probability of returning a full matrix will always
return one. For example adding a scalar constant to a sparse matrix will almost always make it a full matrix, and so the example,
```

speye (3) +0
$\Rightarrow \quad 1 \quad 0 \quad 0$
010
$0 \quad 0 \quad 1$

```
returns a full matrix as can be seen.
Additionally, if sparse_auto_mutate is true, all sparse functions test the amount of memory occupied by the sparse matrix to see if the amount of storage used is larger than the amount used by the full equivalent. Therefore speye (2) * 1 will return a full matrix as the memory used is smaller for the full version than the sparse version.

As all of the mixed operators and functions between full and sparse matrices exist, in general this does not cause any problems. However, one area where it does cause a problem is where a sparse matrix is promoted to a full matrix, where subsequent operations would resparsify the matrix. Such cases are rare, but can be artificially created, for example (fliplr (speye (3)) + speye (3)) - speye (3) gives a full matrix when it should give a sparse one. In general, where such cases occur, they impose only a small memory penalty.

There is however one known case where this behavior of Octave's sparse matrices will cause a problem. That is in the handling of the diag function. Whether diag returns a sparse or full matrix depending on the type of its input arguments. So
\[
a=\operatorname{diag}(\text { sparse }([1,2,3]),-1) \text {; }
\]
should return a sparse matrix. To ensure this actually happens, the sparse function, and other functions based on it like speye, always returns a sparse matrix, even if the memory used will be larger than its full representation.
```

val = sparse_auto_mutate ()
old_val = sparse_auto_mutate (new_val)
sparse_auto_mutate (new_val, "local")

```

Query or set the internal variable that controls whether Octave will automatically mutate sparse matrices to full matrices to save memory.
For example:
```

s = speye (3);
sparse_auto_mutate (false);
s(:, 1) = 1;
typeinfo (s)
=> sparse matrix
sparse_auto_mutate (true);
s(1, :) = 1;
typeinfo (s)
m matrix

```

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

Note that the sparse_auto_mutate option is incompatible with MATLAB, and so it is off by default.

\subsection*{22.1.4.3 Mathematical Considerations}

The attempt has been made to make sparse matrices behave in exactly the same manner as there full counterparts. However, there are certain differences and especially differences with other products sparse implementations.

First, the ". /" and " . " operators must be used with care. Consider what the examples \(\mathrm{s}=\) speye (4);
a1 = s.^ 2;
a2 = s.^s;
a3 = s.^-2;
a4 = s ./ 2;
\(\mathrm{a} 5=2\)./ s;
\(\mathrm{a} 6=\mathrm{s} . / \mathrm{s} ;\)
will give. The first example of \(s\) raised to the power of 2 causes no problems. However \(s\) raised element-wise to itself involves a large number of terms \(0 .{ }^{\wedge} 0\) which is 1 . There \(s .{ }^{\text {- }}\) \(s\) is a full matrix.

Likewise s.^-2 involves terms like \(0 . \wedge-2\) which is infinity, and so s.^-2 is equally a full matrix.

For the "./" operator s./ 2 has no problems, but 2 ./s involves a large number of infinity terms as well and is equally a full matrix. The case of \(s . / s\) involves terms like 0 .\(/ 0\) which is a NaN and so this is equally a full matrix with the zero elements of \(s\) filled with NaN values.

The above behavior is consistent with full matrices, but is not consistent with sparse implementations in other products.

A particular problem of sparse matrices comes about due to the fact that as the zeros are not stored, the sign-bit of these zeros is equally not stored. In certain cases the sign-bit of zero is important. For example:
\[
\begin{array}{ll}
\mathrm{a}=0 . /[-1,1 ; 1, & -1] ; \\
\mathrm{b}=1 . / \mathrm{a} \\
\Rightarrow & -\operatorname{Inf} \\
& \operatorname{Inf} \\
\mathrm{c}= & 1 . / \text { sparse }(\mathrm{Inf} \\
\Rightarrow & \operatorname{Inf} \\
& \text { Inf }
\end{array}
\]

To correct this behavior would mean that zero elements with a negative sign-bit would need to be stored in the matrix to ensure that their sign-bit was respected. This is not done at this time, for reasons of efficiency, and so the user is warned that calculations where the sign-bit of zero is important must not be done using sparse matrices.

In general any function or operator used on a sparse matrix will result in a sparse matrix with the same or a larger number of nonzero elements than the original matrix. This is particularly true for the important case of sparse matrix factorizations. The usual way to address this is to reorder the matrix, such that its factorization is sparser than the factorization of the original matrix. That is the factorization of \(\mathrm{L} * \mathrm{U}=\mathrm{P} * \mathrm{~S} * \mathrm{Q}\) has sparser terms \(L\) and \(U\) than the equivalent factorization \(L * U=S\).

Several functions are available to reorder depending on the type of the matrix to be factorized. If the matrix is symmetric positive-definite, then symamd or csymamd should be used. Otherwise amd, colamd or ccolamd should be used. For completeness the reordering functions colperm and randperm are also available.

See Figure 22.3, for an example of the structure of a simple positive definite matrix.


Figure 22.3: Structure of simple sparse matrix.

The standard Cholesky factorization of this matrix can be obtained by the same command that would be used for a full matrix. This can be visualized with the command \(\mathrm{r}=\) chol (A) ; spy (r);. See Figure 22.4. The original matrix had 598 nonzero terms, while this Cholesky factorization has 10200 , with only half of the symmetric matrix being stored. This is a significant level of fill in, and although not an issue for such a small test case, can represents a large overhead in working with other sparse matrices.

The appropriate sparsity preserving permutation of the original matrix is given by symamd and the factorization using this reordering can be visualized using the command \(\mathrm{q}=\operatorname{symamd}(\mathrm{A}) ; \mathrm{r}=\operatorname{chol}(\mathrm{A}(\mathrm{q}, \mathrm{q}))\); spy ( r ). This gives 399 nonzero terms which is a significant improvement.

The Cholesky factorization itself can be used to determine the appropriate sparsity preserving reordering of the matrix during the factorization, In that case this might be obtained with three return arguments as \([r, p, q]=\operatorname{chol}(A) ; s p y(r)\).


Figure 22.4: Structure of the unpermuted Cholesky factorization of the above matrix.


Figure 22.5: Structure of the permuted Cholesky factorization of the above matrix.

In the case of an asymmetric matrix, the appropriate sparsity preserving permutation is colamd and the factorization using this reordering can be visualized using the command q \(=\operatorname{colamd}(A) ;[1, u, p]=l u(A(:, q)) ; s p y(l+u)\).

Finally, Octave implicitly reorders the matrix when using the div (/) and ldiv ( \(\backslash\) ) operators, and so no the user does not need to explicitly reorder the matrix to maximize performance.
\(p=\) amd (S)
\(p=\operatorname{amd}(S, o p t s)\)
Return the approximate minimum degree permutation of a matrix.
This is a permutation such that the Cholesky factorization of \(S(p, p)\) tends to be sparser than the Cholesky factorization of \(S\) itself. amd is typically faster than symamd but serves a similar purpose.
The optional parameter opts is a structure that controls the behavior of amd. The fields of the structure are
opts.dense Determines what amd considers to be a dense row or column of the input matrix. Rows or columns with more than max (16, (dense * sqrt (n))) entries, where \(n\) is the order of the matrix \(S\), are ignored by amd during the calculation of the permutation. The value of dense must be a positive scalar and the default value is 10.0
opts.aggressive
If this value is a nonzero scalar, then amd performs aggressive absorption. The default is not to perform aggressive absorption.
The author of the code itself is Timothy A. Davis (see http://faculty.cse.tamu. edu/davis/suitesparse.html).
See also: [symamd], page 679, [colamd], page 676.
\(p=\mathrm{ccolamd}(S)\)
\(p=\) ccolamd (S, knobs)
\(p=\) ccolamd (S, knobs, cmember)
[p, stats] = ccolamd (...)
Constrained column approximate minimum degree permutation.
\(p=\) ccolamd ( \(S\) ) returns the column approximate minimum degree permutation vector for the sparse matrix \(S\). For a non-symmetric matrix \(S, S(:, p)\) tends to have sparser LU factors than \(S\). chol \(\left(S(:, p)^{\prime} * S(:, p)\right)\) also tends to be sparser than chol \(\left(S^{\prime} * S\right) . p=\operatorname{ccolamd}(S, 1)\) optimizes the ordering for \(l u(S(:, p))\). The ordering is followed by a column elimination tree post-ordering.
knobs is an optional 1-element to 5 -element input vector, with a default value of [0 \(101010]\) if not present or empty. Entries not present are set to their defaults.
knobs(1) if nonzero, the ordering is optimized for \(l \mathrm{l}(\mathrm{S}(:, \mathrm{p})\) ). It will be a poor ordering for chol \(\left(S(:, p)^{\prime} * S(:, p)\right)\). This is the most important knob for ccolamd.
knobs(2) if \(S\) is m-by-n, rows with more than max (16, knobs(2) * sqrt (n)) entries are ignored.
knobs(3) columns with more than max (16, knobs(3) * sqrt (min (m, n))) entries are ignored and ordered last in the output permutation (subject to the cmember constraints).
knobs(4) if nonzero, aggressive absorption is performed.
knobs(5) if nonzero, statistics and knobs are printed.
cmember is an optional vector of length \(n\). It defines the constraints on the column ordering. If cmember \((j)=c\), then column \(j\) is in constraint set \(c\) ( \(c\) must be in the range 1 to \(n\) ). In the output permutation \(p\), all columns in set 1 appear first, followed by all columns in set 2 , and so on. cmember \(=\) ones ( \(1, n\) ) if not present or empty. ccolamd ( \(S,[], 1: n\) ) returns \(1: \mathrm{n}\)
\(p=\operatorname{ccolamd}(S)\) is about the same as \(p=\operatorname{colamd}(S)\). knobs and its default values differ. colamd always does aggressive absorption, and it finds an ordering suitable for both lu \((S(:, p))\) and chol \(\left(S(:, p)^{\prime} * S(:, p)\right)\); it cannot optimize its ordering for \(l u(S(:, p))\) to the extent that ccolamd ( \(S, 1\) ) can.
stats is an optional 20-element output vector that provides data about the ordering and the validity of the input matrix \(S\). Ordering statistics are in stats (1:3). stats (1) and stats(2) are the number of dense or empty rows and columns ignored by CCOLAMD and stats(3) is the number of garbage collections performed on the internal data structure used by CCOLAMD (roughly of size \(2.2 * \mathrm{nnz}(S)+4 * m+7\) * \(n\) integers).
stats ( \(4: 7\) ) provide information if CCOLAMD was able to continue. The matrix is OK if stats (4) is zero, or 1 if invalid. stats (5) is the rightmost column index that is unsorted or contains duplicate entries, or zero if no such column exists. stats (6) is the last seen duplicate or out-of-order row index in the column index given by stats (5), or zero if no such row index exists. stats (7) is the number of duplicate or out-of-order row indices. stats ( \(8: 20\) ) is always zero in the current version of CCOLAMD (reserved for future use).
The authors of the code itself are S. Larimore, T. Davis and S. Rajamanickam in collaboration with J. Bilbert and E. Ng. Supported by the National Science Foundation (DMS-9504974, DMS-9803599, CCR-0203270), and a grant from Sandia National Lab. See http://faculty.cse.tamu.edu/davis/suitesparse.html for ccolamd, csymamd, amd, colamd, symamd, and other related orderings.

See also: [colamd], page 676, [csymamd], page 677 .
```

p = colamd (S)
p = colamd (S, knobs)
[p, stats] = colamd (S)
[p, stats] = colamd (S, knobs)

```

Compute the column approximate minimum degree permutation.
\(p=\operatorname{col}\) amd ( \(S\) ) returns the column approximate minimum degree permutation vector for the sparse matrix \(S\). For a non-symmetric matrix \(S, S(:, p)\) tends to have sparser LU factors than \(S\). The Cholesky factorization of \(S(:, p)^{\prime} * S(:, p)\) also tends to be sparser than that of \(S^{\prime} * S\).
knobs is an optional one- to three-element input vector. If \(S\) is m-by-n, then rows with more than \(\max (16, \operatorname{knobs}(1) * \operatorname{sqrt}(\mathrm{n}))\) entries are ignored. Columns with more than max \((16, \operatorname{knobs}(2) * \operatorname{sqrt}(\min (m, n)))\) entries are removed prior to ordering, and ordered last in the output permutation \(p\). Only completely dense rows or columns are removed if knobs(1) and knobs(2) are < 0, respectively. If knobs(3) is nonzero, stats and knobs are printed. The default is knobs \(=\left[\begin{array}{lll}10 & 10 & 0\end{array}\right]\). Note that knobs differs from earlier versions of colamd.
stats is an optional 20 -element output vector that provides data about the ordering and the validity of the input matrix \(S\). Ordering statistics are in stats(1:3). stats(1) and stats(2) are the number of dense or empty rows and columns ignored by COLAMD and stats(3) is the number of garbage collections performed on the internal data structure used by COLAMD (roughly of size \(2.2 * \mathrm{nnz}(S)+4 * m+7 *\) \(n\) integers).
Octave built-in functions are intended to generate valid sparse matrices, with no duplicate entries, with ascending row indices of the nonzeros in each column, with a non-negative number of entries in each column (!) and so on. If a matrix is invalid, then COLAMD may or may not be able to continue. If there are duplicate entries (a row index appears two or more times in the same column) or if the row indices in a column are out of order, then ColamD can correct these errors by ignoring the duplicate entries and sorting each column of its internal copy of the matrix \(S\) (the input matrix \(S\) is not repaired, however). If a matrix is invalid in other ways then COLAMD cannot continue, an error message is printed, and no output arguments ( \(p\) or stats) are returned. COLAMD is thus a simple way to check a sparse matrix to see if it's valid.
stats(4:7) provide information if colamd was able to continue. The matrix is OK if stats(4) is zero, or 1 if invalid. stats(5) is the rightmost column index that is unsorted or contains duplicate entries, or zero if no such column exists. stats(6) is the last seen duplicate or out-of-order row index in the column index given by stats(5), or zero if no such row index exists. stats(7) is the number of duplicate or out-of-order row indices. stats( \(8: 20\) ) is always zero in the current version of COLAMD (reserved for future use).
The ordering is followed by a column elimination tree post-ordering.
The authors of the code itself are Stefan I. Larimore and Timothy A. Davis. The algorithm was developed in collaboration with John Gilbert, Xerox PARC, and Esmond Ng, Oak Ridge National Laboratory. (see http://faculty.cse.tamu.edu/davis/ suitesparse.html)
See also: [colperm], page 677, [symamd], page 679, [ccolamd], page 675.
\(p=\) colperm (s)
Return the column permutations such that the columns of \(s(:, p)\) are ordered in terms of increasing number of nonzero elements.
If \(s\) is symmetric, then \(p\) is chosen such that \(s(p, p)\) orders the rows and columns with increasing number of nonzero elements.
```

p = csymamd (S)
p = csymamd (S, knobs)
p = csymamd (S, knobs, cmember)
[p, stats] = csymamd (...)

```

For a symmetric positive definite matrix \(S\), return the permutation vector \(p\) such that \(S(p, p)\) tends to have a sparser Cholesky factor than \(S\).
Sometimes csymamd works well for symmetric indefinite matrices too. The matrix \(S\) is assumed to be symmetric; only the strictly lower triangular part is referenced. \(S\) must be square. The ordering is followed by an elimination tree post-ordering.
knobs is an optional 1-element to 3 -element input vector, with a default value of [10 \(10]\). Entries not present are set to their defaults.
knobs(1) If \(S\) is n-by-n, then rows and columns with more than \(\max (16\), knobs(1)*sqrt(n)) entries are ignored, and ordered last in the output permutation (subject to the cmember constraints).
knobs(2) If nonzero, aggressive absorption is performed.
knobs(3) If nonzero, statistics and knobs are printed.
cmember is an optional vector of length n. It defines the constraints on the ordering. If cmember \((\mathrm{j})=S\), then row/column j is in constraint set \(c\) ( \(c\) must be in the range 1 to n ). In the output permutation \(p\), rows/columns in set 1 appear first, followed by all rows/columns in set 2 , and so on. cmember \(=\) ones \((1, n)\) if not present or empty. csymamd ( \(\mathrm{S},[\mathrm{l}, 1: \mathrm{n}\) ) returns \(1: \mathrm{n}\).
\(p=\operatorname{csymamd}(S)\) is about the same as \(p=\operatorname{symamd}(S)\). knobs and its default values differ.
stats( \(4: 7\) ) provide information if CCOLAMD was able to continue. The matrix is OK if stats(4) is zero, or 1 if invalid. stats(5) is the rightmost column index that is unsorted or contains duplicate entries, or zero if no such column exists. stats(6) is the last seen duplicate or out-of-order row index in the column index given by stats(5), or zero if no such row index exists. stats(7) is the number of duplicate or out-of-order row indices. stats( \(8: 20\) ) is always zero in the current version of CCOLAMD (reserved for future use).
The authors of the code itself are S. Larimore, T. Davis and S. Rajamanickam in collaboration with J. Bilbert and E. Ng. Supported by the National Science Foundation (DMS-9504974, DMS-9803599, CCR-0203270), and a grant from Sandia National Lab. See http://faculty.cse.tamu.edu/davis/suitesparse.html for ccolamd, colamd, csymamd, amd, colamd, symamd, and other related orderings.

See also: [symamd], page 679, [ccolamd], page 675.
\(p=\) dmperm (A)
\([p, q, r, s, c c, r r]=\operatorname{dmperm}(A)\)
Perform a Dulmage-Mendelsohn permutation of the sparse matrix \(A\).
With a single output argument dmperm, return a maximum matching \(p\) such that \(\mathrm{p}(\mathrm{j})=\mathrm{i}\) if column \(j\) is matched to row \(i\), or 0 if column \(j\) is unmatched. If \(A\) is square and full structural rank, \(p\) is a row permutation and \(\mathrm{A}(\mathrm{p},:\) ) has a zero-free diagonal. The structural rank of \(A\) is \(\operatorname{sprank}(\mathrm{A})=\operatorname{sum}(\mathrm{p}>0)\).
Called with two or more output arguments, return the Dulmage-Mendelsohn decomposition of A. p and \(q\) are permutation vectors. \(c c\) and \(r r\) are vectors of length 5. c \(=\mathrm{A}(\mathrm{p}, q)\) is split into a 4 -by- 4 set of coarse blocks:
\begin{tabular}{cccc} 
A11 & A12 & A13 & A14 \\
0 & 0 & A23 & A24 \\
0 & 0 & 0 & A34 \\
0 & 0 & 0 & A44
\end{tabular}
where A12, A23, and A34 are square with zero-free diagonals. The columns of A11 are the unmatched columns, and the rows of A44 are the unmatched rows. Any
of these blocks can be empty. In the "coarse" decomposition, the ( \(\mathrm{i}, \mathrm{j}\) )-th block is \(C(r r(i): r r(i+1)-1, c c(j): c c(j+1)-1)\). In terms of a linear system, [A11A12] is the underdetermined part of the system (it is always rectangular and with more columns and rows, or 0 -by- 0 ), A23 is the well-determined part of the system (it is always square), and [A34 ; A44] is the over-determined part of the system (it is always rectangular with more rows than columns, or 0 -by- 0 ).
The structural rank of \(A\) is sprank \((A)=r r(4)-1\), which is an upper bound on the numerical \(\operatorname{rank}\) of \(A . \operatorname{sprank}(A)=\operatorname{rank}(f u l l(\operatorname{sprand}(A)))\) with probability 1 in exact arithmetic.
The A23 submatrix is further subdivided into block upper triangular form via the "fine" decomposition (the strongly-connected components of A23). If \(A\) is square and structurally non-singular, A23 is the entire matrix.
\(C(r(i): r(i+1)-1, s(j): s(j+1)-1)\) is the \((i, j)\)-th block of the fine decomposition. The ( 1,1 ) block is the rectangular block [A11 A12], unless this block is 0 -by- 0 . The ( \(\mathrm{b}, \mathrm{b}\) ) block is the rectangular block [A34 ; A44], unless this block is 0 -by- 0 , where b \(=\) length \((r)-1\). All other blocks of the form \(C(r(i): r(i+1)-1, s(i): s(i+1)-1)\) are diagonal blocks of A23, and are square with a zero-free diagonal.
The method used is described in: A. Pothen \& C.-J. Fan. Computing the Block Triangular Form of a Sparse Matrix. ACM Trans. Math. Software, 16(4):303-324, 1990.

See also: [colamd], page 676, [ccolamd], page 675.
```

p = symamd (S)
p = symamd (S, knobs)
[p, stats] = symamd (S)
[p, stats] = symamd (S, knobs)

```

For a symmetric positive definite matrix \(S\), returns the permutation vector p such that \(S(p, p)\) tends to have a sparser Cholesky factor than \(S\).
Sometimes symamd works well for symmetric indefinite matrices too. The matrix \(S\) is assumed to be symmetric; only the strictly lower triangular part is referenced. \(S\) must be square.
knobs is an optional one- to two-element input vector. If \(S\) is \(n\)-by-n, then rows and columns with more than max \((16, \operatorname{knobs}(1) * s q r t(n))\) entries are removed prior to ordering, and ordered last in the output permutation \(p\). No rows/columns are removed if knobs(1) < 0. If knobs(2) is nonzero, stats and knobs are printed. The default is knobs \(=\left[\begin{array}{ll}10 & 0\end{array}\right]\). Note that knobs differs from earlier versions of symamd.
stats is an optional 20 -element output vector that provides data about the ordering and the validity of the input matrix \(S\). Ordering statistics are in stats(1:3). \(\operatorname{stats}(1)=\operatorname{stats}(2)\) is the number of dense or empty rows and columns ignored by SYMAMD and stats(3) is the number of garbage collections performed on the internal data structure used by SYMAMD (roughly of size \(8.4 * \mathrm{nnz}\) (tril ( \(S,-1\) ) ) \(+9 * n\) integers).
Octave built-in functions are intended to generate valid sparse matrices, with no duplicate entries, with ascending row indices of the nonzeros in each column, with a non-negative number of entries in each column (!) and so on. If a matrix is invalid,
then SYMAMD may or may not be able to continue. If there are duplicate entries (a row index appears two or more times in the same column) or if the row indices in a column are out of order, then SYMAMD can correct these errors by ignoring the duplicate entries and sorting each column of its internal copy of the matrix \(S\) (the input matrix \(S\) is not repaired, however). If a matrix is invalid in other ways then SYMAMD cannot continue, an error message is printed, and no output arguments ( \(p\) or stats) are returned. SYMAMD is thus a simple way to check a sparse matrix to see if it's valid.
stats(4:7) provide information if SYMAMD was able to continue. The matrix is OK if stats (4) is zero, or 1 if invalid. stats(5) is the rightmost column index that is unsorted or contains duplicate entries, or zero if no such column exists. stats(6) is the last seen duplicate or out-of-order row index in the column index given by stats(5), or zero if no such row index exists. stats(7) is the number of duplicate or out-of-order row indices. stats ( \(8: 20\) ) is always zero in the current version of SYMAMD (reserved for future use).
The ordering is followed by a column elimination tree post-ordering.
The authors of the code itself are Stefan I. Larimore and Timothy A. Davis. The algorithm was developed in collaboration with John Gilbert, Xerox PARC, and Esmond Ng, Oak Ridge National Laboratory. (see http://faculty.cse.tamu.edu/davis/ suitesparse.html)
See also: [colperm], page 677, [colamd], page 676.
\(p=\operatorname{symrcm}(S)\)
Return the symmetric reverse Cuthill-McKee permutation of \(S\).
\(p\) is a permutation vector such that \(S(p, p)\) tends to have its diagonal elements closer to the diagonal than \(S\). This is a good preordering for LU or Cholesky factorization of matrices that come from "long, skinny" problems. It works for both symmetric and asymmetric \(S\).
The algorithm represents a heuristic approach to the NP-complete bandwidth minimization problem. The implementation is based in the descriptions found in
E. Cuthill, J. McKee. Reducing the Bandwidth of Sparse Symmetric Matrices. Proceedings of the 24th ACM National Conference, 157-172 1969, Brandon Press, New Jersey.
A. George, J.W.H. Liu. Computer Solution of Large Sparse Positive Definite Systems, Prentice Hall Series in Computational Mathematics, ISBN 0-13-165274-5, 1981.
See also: [colperm], page 677, [colamd], page 676, [symamd], page 679.

\subsection*{22.2 Linear Algebra on Sparse Matrices}

Octave includes a polymorphic solver for sparse matrices, where the exact solver used to factorize the matrix, depends on the properties of the sparse matrix itself. Generally, the cost of determining the matrix type is small relative to the cost of factorizing the matrix itself, but in any case the matrix type is cached once it is calculated, so that it is not re-determined each time it is used in a linear equation.

The selection tree for how the linear equation is solve is
1. If the matrix is diagonal, solve directly and goto 8
2. If the matrix is a permuted diagonal, solve directly taking into account the permutations. Goto 8
3. If the matrix is square, banded and if the band density is less than that given by spparms ("bandden") continue, else goto 4.
a. If the matrix is tridiagonal and the right-hand side is not sparse continue, else goto 3b.
1. If the matrix is Hermitian, with a positive real diagonal, attempt Cholesky factorization using LAPACK xPTSV.
2. If the above failed or the matrix is not Hermitian with a positive real diagonal use Gaussian elimination with pivoting using Lapack xGTSV, and goto 8.
b. If the matrix is Hermitian with a positive real diagonal, attempt Cholesky factorization using LAPACK xPBTRF.
c. if the above failed or the matrix is not Hermitian with a positive real diagonal use Gaussian elimination with pivoting using LAPACK xGBTRF, and goto 8 .
4. If the matrix is upper or lower triangular perform a sparse forward or backward substitution, and goto 8
5. If the matrix is an upper triangular matrix with column permutations or lower triangular matrix with row permutations, perform a sparse forward or backward substitution, and goto 8
6. If the matrix is square, Hermitian with a real positive diagonal, attempt sparse Cholesky factorization using Cholmod.
7. If the sparse Cholesky factorization failed or the matrix is not Hermitian with a real positive diagonal, and the matrix is square, factorize, solve, and perform one refinement iteration using UMFPACK.
8. If the matrix is not square, or any of the previous solvers flags a singular or near singular matrix, find a minimum norm solution using CXSPARSE \({ }^{3}\).

The band density is defined as the number of nonzero values in the band divided by the total number of values in the full band. The banded matrix solvers can be entirely disabled by using spparms to set bandden to 1 (i.e., spparms ("bandden", 1)).

The QR solver factorizes the problem with a Dulmage-Mendelsohn decomposition, to separate the problem into blocks that can be treated as over-determined, multiple well determined blocks, and a final over-determined block. For matrices with blocks of strongly connected nodes this is a big win as LU decomposition can be used for many blocks. It also significantly improves the chance of finding a solution to over-determined problems rather than just returning a vector of NaN 's.

All of the solvers above, can calculate an estimate of the condition number. This can be used to detect numerical stability problems in the solution and force a minimum norm solution to be used. However, for narrow banded, triangular or diagonal matrices, the cost of calculating the condition number is significant, and can in fact exceed the cost of factoring the matrix. Therefore the condition number is not calculated in these cases, and Octave

\footnotetext{
3 The ChOLMOD, UMFPACK and CXSPARSE packages were written by Tim Davis and are available at http: // faculty.cse.tamu.edu/davis/suitesparse.html
}
relies on simpler techniques to detect singular matrices or the underlying LAPACK code in the case of banded matrices.

The user can force the type of the matrix with the matrix_type function. This overcomes the cost of discovering the type of the matrix. However, it should be noted that identifying the type of the matrix incorrectly will lead to unpredictable results, and so matrix_type should be used with care.
```

nest = normest (A)
nest = normest (A, tol)
[nest, iter] = normest (...)

```

Estimate the 2 -norm of the matrix \(A\) using a power series analysis.
This is typically used for large matrices, where the cost of calculating norm (A) is prohibitive and an approximation to the 2-norm is acceptable.
tol is the tolerance to which the 2-norm is calculated. By default tol is \(1 \mathrm{e}-6\).
The optional output iter returns the number of iterations that were required for normest to converge.
See also: [normest1], page 682, [norm], page 592, [cond], page 586, [condest], page 683.
```

nest = normest1 (A)
nest = normest1 (A, t)
nest = normest1 (A, t, x0)
nest = normest1 (Afun, t, x0, p1, p2,...)
[nest, v] = normest1 (A,...)
[nest, v, w] = normest1 (A,...)
[nest, v, w, iter] = normest1 (A,...)

```

Estimate the 1-norm of the matrix \(A\) using a block algorithm.
normest1 is best for large sparse matrices where only an estimate of the norm is required. For small to medium sized matrices, consider using norm ( \(A, 1\) ). In addition, normest1 can be used for the estimate of the 1-norm of a linear operator \(A\) when matrix-vector products \(A * x\) and \(A^{\prime} * x\) can be cheaply computed. In this case, instead of the matrix \(A\), a function Afun (flag, x) is used; it must return:
- the dimension \(n\) of \(A\), if flag is "dim"
- true if \(A\) is a real operator, if flag is "real"
- the result \(A * x\), if flag is "notransp"
- the result \(A^{\prime} * x\), if flag is "transp"

A typical case is \(A\) defined by \(b^{\wedge} m\), in which the result \(A * x\) can be computed without even forming explicitly \(b^{\wedge} m\) by:
```

y = x;
for i = 1:m
y = b * y;
endfor

```

The parameters \(p 1, p 2, \ldots\) are arguments of Afun (flag, x, p1, p2, ...).
The default value for \(t\) is 2 . The algorithm requires matrix-matrix products with sizes \(n \times n\) and \(n \times t\).

The initial matrix \(x 0\) should have columns of unit 1-norm. The default initial matrix \(x 0\) has the first column ones \((n, 1) / n\) and, if \(t>1\), the remaining columns with random elements \(-1 / n, 1 / n\), divided by \(n\).
On output, nest is the desired estimate, \(v\) and \(w\) are vectors such that \(w=A * v\), with norm ( \(\mathrm{w}, 1\) ) \(=c *\) norm ( \(v, 1\) ). iter contains in iter (1) the number of iterations (the maximum is hardcoded to 5) and in iter(2) the total number of products \(A *\) x or \(A^{\prime} * \mathrm{x}\) performed by the algorithm.
Algorithm Note: normest1 uses random numbers during evaluation. Therefore, if consistent results are required, the "state" of the random generator should be fixed before invoking normest1.
Reference: N. J. Higham and F. Tisseur, A block algorithm for matrix 1-norm estimation, with and application to 1-norm pseudospectra, SIAM J. Matrix Anal. Appl., pp. 1185-1201, Vol 21, No. 4, 2000.
See also: [normest], page 682, [norm], page 592, [cond], page 586, [condest], page 683.
```

cest = condest (A)
cest = condest (A, t)
cest = condest (A, Ainvfcn)
cest = condest (A,Ainvfcn, t)
cest = condest (A, Ainvfcn, t, p1, p2,...)
cest = condest (Afcn, Ainvfcn)
cest = condest (Afcn, Ainvfcn, t)
cest = condest (Afcn, Ainvfcn, t, p1, p2,...)
[cest, v] = condest (...)

```

Estimate the 1-norm condition number of a square matrix \(A\) using \(t\) test vectors and a randomized 1-norm estimator.
The optional input \(t\) specifies the number of test vectors (default 5).
The input may be a matrix \(A\) (the algorithm is particularly appropriate for large, sparse matrices). Alternatively, the behavior of the matrix can be defined implicitly by functions. When using an implicit definition, condest requires the following functions:
- Afcn (flag, x) which must return
- the dimension \(n\) of \(A\), if flag is "dim"
- true if \(A\) is a real operator, if flag is "real"
- the result \(A * x\), if flag is "notransp"
- the result \(A^{\prime} * x\), if flag is "transp"
- Ainvfen (flag, x) which must return
- the dimension \(n\) of inv ( \(A\) ), if flag is "dim"
- true if inv (A) is a real operator, if flag is "real"
- the result inv (A) * x, if flag is "notransp"
- the result inv (A) ' \(* x\) x, if flag is "transp"

Any parameters \(p 1, p 2, \ldots\) are additional arguments of Afcn (flag, \(x, p 1, p 2\), ...) and Ainvfcn (flag, x, p1, p2,...).

The principal output is the 1-norm condition number estimate cest.
The optional second output \(v\) is an approximate null vector; it satisfies the equation norm ( \(A * V, 1\) ) \(==\operatorname{norm}(A, 1) * \operatorname{norm}(v, 1) /\) cest.
Algorithm Note: condest uses a randomized algorithm to approximate the 1-norms. Therefore, if consistent results are required, the "state" of the random generator should be fixed before invoking condest.
References:
- N.J. Higham and F. Tisseur, A Block Algorithm for Matrix 1-Norm Estimation, with an Application to 1-Norm Pseudospectra. SIMAX vol 21, no 4, pp 1185-1201. https://dx.doi.org/10.1137/S0895479899356080
- N.J. Higham and F. Tisseur, A Block Algorithm for Matrix 1-Norm Estimation, with an Application to 1-Norm Pseudospectra. https://citeseer.ist.psu. edu/223007.html

See also: [cond], page 586, [rcond], page 594, [norm], page 592, [normest1], page 682, [normest], page 682.
spparms ()
vals = spparms ()
[keys, vals] = spparms ()
val = spparms (key)
spparms (vals)
spparms ("default")
spparms ("tight")
spparms (key, val)
Query or set the parameters used by the sparse solvers and factorization functions.
The first four calls above get information about the current settings, while the others change the current settings. The parameters are stored as pairs of keys and values, where the values are all floats and the keys are one of the following strings:
'spumoni' Printing level of debugging information of the solvers (default 0)
'ths_rel' Included for compatibility. Not used. (default 1)
'ths_abs' Included for compatibility. Not used. (default 1)
'exact_d' Included for compatibility. Not used. (default 0)
'supernd' Included for compatibility. Not used. (default 3)
'rreduce' Included for compatibility. Not used. (default 3)
'wh_frac' Included for compatibility. Not used. (default 0.5)
'autommd' Flag whether the LU/QR and the ' ' and '/' operators will automatically use the sparsity preserving mmd functions (default 1)
'autoamd' Flag whether the LU and the ' \(\backslash\) ' and '/' operators will automatically use the sparsity preserving amd functions (default 1)
'piv_tol' The pivot tolerance of the UMFPACK solvers (default 0.1)
'sym_tol' The pivot tolerance of the UMFPACK symmetric solvers (default 0.001)
'bandden' The density of nonzero elements in a banded matrix before it is treated by the lapack banded solvers (default 0.5)
'umfpack' Flag whether the UMFPACK or mmd solvers are used for the LU , ' X ' and '/' operations (default 1)
The value of individual keys can be set with spparms (key, val). The default values can be restored with the special keyword "default". The special keyword "tight" can be used to set the mmd solvers to attempt a sparser solution at the potential cost of longer running time.
See also: [chol], page 595, [colamd], page 676, [lu], page 597, [qr], page 599, [symamd], page 679 .
\(p=\operatorname{sprank}(S)\)
Calculate the structural rank of the sparse matrix \(S\).
Note that only the structure of the matrix is used in this calculation based on a Dulmage-Mendelsohn permutation to block triangular form. As such the numerical rank of the matrix \(S\) is bounded by sprank ( \(S\) ) >= rank ( \(S\) ). Ignoring floating point errors sprank ( \(S\) ) == rank ( \(S\) ).
See also: [dmperm], page 678.
[count, h, parent, post, R] = symbfact (S)
[...] = symbfact (S, typ)
[...] = symbfact (S, typ, mode)
Perform a symbolic factorization analysis of the sparse matrix \(S\).
The input variables are
\(S \quad S\) is a real or complex sparse matrix.
typ Is the type of the factorization and can be one of "sym" (default)

Factorize \(S\). Assumes \(S\) is symmetric and uses the upper triangular portion of the matrix.
"col" Factorize \(S^{\prime} * S\).
"row" Factorize \(S * S^{\prime}\).
"lo" Factorize \(S^{\prime}\). Assumes \(S\) is symmetric and uses the lower triangular portion of the matrix.
mode When mode is unspecified return the Cholesky factorization for \(R\). If mode is "lower" or "L" then return the conjugate transpose \(R^{\prime}\) which is a lower triangular factor. The conjugate transpose version is faster and uses less memory, but still returns the same values for all other outputs: count, h, parent, and post.
The output variables are:
count The row counts of the Cholesky factorization as determined by typ. The computational difficulty of performing the true factorization using chol is sum (count .^ 2).
\(h \quad\) The height of the elimination tree.
parent The elimination tree itself.
post A sparse boolean matrix whose structure is that of the Cholesky factorization as determined by typ.

See also: [chol], page 595, [etree], page 669, [treelayout], page 669.
For non square matrices, the user can also utilize the spaugment function to find a least squares solution to a linear equation.
s = spaugment ( \(A, c\) )
Create the augmented matrix of \(A\).
This is given by
```

[c * eye(m, m), A;
A', zeros(n, n)]

```

This is related to the least squares solution of \(A \backslash b\), by
```

s * [ r / c; x] = [ b, zeros(n, columns(b)) ]

```
where \(r\) is the residual error
\[
r=b-A * x
\]

As the matrix \(s\) is symmetric indefinite it can be factorized with \(l u\), and the minimum norm solution can therefore be found without the need for a qr factorization. As the residual error will be zeros ( \(m, m\) ) for underdetermined problems, and example can be
```

m = 11; n = 10; mn = max (m, n);
A = spdiags ([ones(mn,1), 10*ones(mn,1), -ones(mn,1)],
[-1, 0, 1], m, n);
x0 = A \ ones (m,1);
s = spaugment (A);
[L, U, P, Q] = lu (s);
x1 = Q * (U \ (L \ (P * [ones(m,1); zeros(n,1)])));
x1 = x1(end - n + 1 : end);

```

To find the solution of an overdetermined problem needs an estimate of the residual error \(r\) and so it is more complex to formulate a minimum norm solution using the spaugment function.
In general the left division operator is more stable and faster than using the spaugment function.

See also: [mldivide], page 162.
Finally, the function eigs can be used to calculate a limited number of eigenvalues and eigenvectors based on a selection criteria and likewise for svds which calculates a limited number of singular values and vectors.
```

d = eigs (A)
d = eigs (A, k)
d = eigs (A, k, sigma)

```
```

d = eigs (A, k, sigma, opts)
d = eigs (A,B)
d = eigs (A, B, k)
d = eigs (A, B, k, sigma)
d = eigs (A,B, k, sigma, opts)
d = eigs (Af, n)
d = eigs (Af, n, k)
d = eigs (Af, n, k, sigma)
d = eigs (Af, n, k, sigma, opts)
d = eigs (Af, n, B)
d = eigs (Af, n, B, k)
d = eigs (Af, n, B, k, sigma)
d = eigs (Af, n, B, k, sigma, opts)
[V, D] = eigs (...)
[V, D, flag] = eigs (...)

```

Calculate a limited number of eigenvalues and eigenvectors based on a selection criteria.
By default, eigs solve the equation \(A \nu=\lambda \nu\), where \(\lambda\) is a scalar representing one of the eigenvalues, and \(\nu\) is the corresponding eigenvector. If given the positive definite matrix \(B\) then eigs solves the general eigenvalue equation \(A \nu=\lambda B \nu\).
The input \(A\) is a square matrix of dimension \(n\)-by-n. Typically, \(A\) is also large and sparse.

The input \(B\) for the generalized eigenvalue problem is a square matrix with the same size as \(A\) ( \(n\)-by- \(n\) ). Typically, \(B\) is also large and sparse.
The number of eigenvalues and eigenvectors to calculate is given by \(k\) and defaults to 6.

The argument sigma determines which eigenvalues are returned. sigma can be either a scalar or a string. When sigma is a scalar, the k eigenvalues closest to sigma are returned. If sigma is a string, it must be one of the following values.
"lm" Largest Magnitude (default).
"sm" Smallest Magnitude.
"la" Largest Algebraic (valid only for real symmetric problems).
"sa" Smallest Algebraic (valid only for real symmetric problems).
"be" Both Ends, with one more from the high-end if \(k\) is odd (valid only for real symmetric problems).
"lr" Largest Real part (valid only for complex or unsymmetric problems).
"sr" Smallest Real part (valid only for complex or unsymmetric problems).
"li" Largest Imaginary part (valid only for complex or unsymmetric problems).
"si" Smallest Imaginary part (valid only for complex or unsymmetric problems).

If opts is given, it is a structure defining possible options that eigs should use. The fields of the opts structure are:
issym If \(A f\) is given then this flag (true/false) determines whether the function \(A f\) defines a symmetric problem. It is ignored if a matrix \(A\) is given. The default is false.
isreal If \(A f\) is given then this flag (true/false) determines whether the function \(A f\) defines a real problem. It is ignored if a matrix \(A\) is given. The default is true.
tol Defines the required convergence tolerance, calculated as tol \(*\) norm (A). The default is eps.
maxit The maximum number of iterations. The default is 300 .
p The number of Lanczos basis vectors to use. More vectors will result in faster convergence, but a greater use of memory. The optimal value of \(p\) is problem dependent and should be in the range \(k+1\) to \(n\). The default value is \(2 * k\).
v0 The starting vector for the algorithm. An initial vector close to the final vector will speed up convergence. The default is for ARPack to randomly generate a starting vector. If specified, v0 must be an \(n\)-by- 1 vector where \(n=\) rows ( \(A\) ).
disp The level of diagnostic printout ( \(0|1| 2\) ). If disp is 0 then diagnostics are disabled. The default value is 0 .
cholB If the generalized eigenvalue problem is being calculated, this flag (true/false) specifies whether the \(B\) input represents chol ( \(B\) ) or simply the matrix \(B\). The default is false.
permB The permutation vector of the Cholesky factorization for \(B\) if cholB is true. It is obtained by \([R, \sim\), permB] \(=\operatorname{chol}(B\), vector"). The default is \(1: n\).

It is also possible to represent \(A\) by a function denoted Af. Af must be followed by a scalar argument \(n\) defining the length of the vector argument accepted by Af. Af can be a function handle, an inline function, or a string. When \(A f\) is a string it holds the name of the function to use.
Af is a function of the form \(\mathrm{y}=\mathrm{Af}\) ( x ) where the required return value of \(A f\) is determined by the value of sigma. The four possible forms are
\(A * x \quad\) if sigma is not given or is a string other than "sm".
\(\mathrm{A} \backslash \mathrm{x} \quad\) if sigma is 0 or " sm ".
( \(\mathrm{A}-\operatorname{sigma} * \mathrm{I}) \backslash \mathrm{x}\)
if sigma is a scalar not equal to 0 ; \(I\) is the identity matrix of the same size as \(A\).
\((\mathrm{A}-\operatorname{sigma} * \mathrm{~B}) \backslash \mathrm{x}\) for the general eigenvalue problem.

The return arguments and their form depend on the number of return arguments requested. For a single return argument, a column vector \(d\) of length \(k\) is returned containing the \(k\) eigenvalues that have been found. For two return arguments, \(V\) is an \(n\)-by- \(k\) matrix whose columns are the \(k\) eigenvectors corresponding to the returned eigenvalues. The eigenvalues themselves are returned in \(D\) in the form of a \(k\)-by- \(k\) matrix, where the elements on the diagonal are the eigenvalues.
The third return argument flag returns the status of the convergence. If flag is 0 then all eigenvalues have converged. Any other value indicates a failure to converge.
Programming Notes: For small problems, \(n<500\), consider using eig (full (A)).
If ARPACK fails to converge consider increasing the number of Lanczos vectors (opt.p), increasing the number of iterations (opt.maxiter), or decreasing the tolerance (opt.tol).
Reference: This function is based on the Arpack package, written by R. Lehoucq, K. Maschhoff, D. Sorensen, and C. Yang. For more information see http://www.caam. rice.edu/software/ARPACK/.
See also: [eig], page 587, [svds], page 689.
```

s = svds (A)
s = svds (A,k)
s = svds (A, k, sigma)
s = svds (A, k, sigma, opts)
[u, s, v] = svds (...)
[u, s, v, flag] = svds (...)

```

Find a few singular values of the matrix \(A\).
The singular values are calculated using
```

[m, n] = size (A);
s = eigs ([sparse(m, m), A;
A', sparse(n, n)])

```

The eigenvalues returned by eigs correspond to the singular values of \(A\). The number of singular values to calculate is given by \(k\) and defaults to 6 .
The argument sigma specifies which singular values to find. When sigma is the string 'L', the default, the largest singular values of \(A\) are found. Otherwise, sigma must be a real scalar and the singular values closest to sigma are found. As a corollary, sigma \(=0\) finds the smallest singular values. Note that for relatively small values of sigma, there is a chance that the requested number of singular values will not be found. In that case sigma should be increased.
opts is a structure defining options that svds will pass to eigs. The possible fields of this structure are documented in eigs. By default, svds sets the following three fields:
tol The required convergence tolerance for the singular values. The default value is \(1 \mathrm{e}-10\). eigs is passed tol / sqrt (2).
maxit The maximum number of iterations. The default is 300 .
disp \(\quad\) The level of diagnostic printout ( \(0|1| 2\) ). If disp is 0 then diagnostics are disabled. The default value is 0 .

If more than one output is requested then svds will return an approximation of the singular value decomposition of \(A\)
```

A_approx = u*S*v'

```
where \(A_{-}\)approx is a matrix of size \(A\) but only rank \(k\).
flag returns 0 if the algorithm has successfully converged, and 1 otherwise. The test for convergence is
```

norm (A*v - u*s, 1) <= tol * norm (A, 1)

```
svds is best for finding only a few singular values from a large sparse matrix. Otherwise, svd (full (A)) will likely be more efficient.
See also: [svd], page 607, [eigs], page 686.

\subsection*{22.3 Iterative Techniques Applied to Sparse Matrices}

The left division \and right division / operators, discussed in the previous section, use direct solvers to resolve a linear equation of the form \(x=A \backslash b\) or \(x=b / A\). Octave also includes a number of functions to solve sparse linear equations using iterative techniques.
```

x = pcg (A, b, tol, maxit, m1, m2, x0, ...)
x = pcg (A, b, tol, maxit, M, [], x0, ...)
[x, flag, relres, iter, resvec, eigest] = pcg (A, b,···)

```

Solve the linear system of equations \(A * x=b\) by means of the Preconditioned Conjugate Gradient iterative method.
The input arguments are:
- \(A\) is the matrix of the linear system and it must be square. \(A\) can be passed as a matrix, function handle, or inline function Afun such that Afun ( x ) \(=\mathrm{A} * \mathrm{x}\). Additional parameters to Afun may be passed after x0.
\(A\) has to be Hermitian and Positive Definite (HPD). If pcg detects \(A\) not to be positive definite, a warning is printed and the flag output is set.
- \(b\) is the right-hand side vector.
- tol is the required relative tolerance for the residual error, \(b-A * x\). The iteration stops if norm \((b-A * x) \leq t o l *\) norm (b). If tol is omitted or empty, then a tolerance of 1e-6 is used.
- maxit is the maximum allowed number of iterations; if maxit is omitted or empty then a value of 20 is used.
- \(m\) is a HPD preconditioning matrix. For any decomposition \(m=p 1 * p 2\) such that inv ( \(p 1\) ) \(* A *\) inv ( \(p 2\) ) is HPD, the conjugate gradient method is formally applied to the linear system inv (p1) *A * inv (p2) \(* y=i n v(p 1) * b\), with \(x\) \(=\operatorname{inv}(p 2) * y\) (split preconditioning). In practice, at each iteration of the conjugate gradient method a linear system with matrix \(m\) is solved with mldivide. If a particular factorization \(m=m 1 * m 2\) is available (for instance, an incomplete Cholesky factorization of a), the two matrices \(m 1\) and \(m 2\) can be passed and the relative linear systems are solved with the mldivide operator. Note that a proper choice of the preconditioner may dramatically improve the overall performance of the method. Instead of matrices \(m 1\) and \(m 2\), the user may pass two functions
which return the results of applying the inverse of \(m 1\) and \(m 2\) to a vector. If \(m 1\) is omitted or empty [], then no preconditioning is applied. If no factorization of \(m\) is available, \(m 2\) can be omitted or left [], and the input variable \(m 1\) can be used to pass the preconditioner \(m\).
- \(x 0\) is the initial guess. If \(x 0\) is omitted or empty then the function sets \(x 0\) to a zero vector by default.

The arguments which follow \(x 0\) are treated as parameters, and passed in an appropriate manner to any of the functions ( \(A\) or m 1 or m 2 ) that have been given to pcg . See the examples below for further details.

The output arguments are:
- \(x\) is the computed approximation to the solution of \(A * x=b\). If the algorithm did not converge, then \(x\) is the iteration which has the minimum residual.
- flag reports on the convergence:
- 0: The algorithm converged to within the prescribed tolerance.
- 1: The algorithm did not converge and it reached the maximum number of iterations.
- 2: The preconditioner matrix is singular.
- 3: The algorithm stagnated, i.e., the absolute value of the difference between the current iteration \(x\) and the previous is less than eps \(*\) norm ( \(x, 2\) ).
- 4: The algorithm detects that the input (preconditioned) matrix is not HPD.
- relres is the ratio of the final residual to its initial value, measured in the Euclidean norm.
- iter indicates the iteration of \(x\) which it was computed. Since the output \(x\) corresponds to the minimal residual solution, the total number of iterations that the method performed is given by length(resvec) - 1 .
- resvec describes the convergence history of the method. resvec ( \(i, 1\) ) is the Euclidean norm of the residual, and resvec ( \(i, 2\) ) is the preconditioned residual norm, after the ( \(i-1\) )-th iteration, \(i=1,2, \ldots\), iter +1 . The preconditioned residual norm is defined as \(r^{\prime} *(m \backslash r)\) where \(r=b-A * x\), see also the description of \(m\). If eigest is not required, only resvec (:,1) is returned.
- eigest returns the estimate for the smallest eigest(1) and largest eigest(2) eigenvalues of the preconditioned matrix \(P=m \backslash A\). In particular, if no preconditioning is used, the estimates for the extreme eigenvalues of \(A\) are returned. eigest(1) is an overestimate and eigest(2) is an underestimate, so that eigest(2) / eigest(1) is a lower bound for cond ( \(P, 2\) ), which nevertheless in the limit should theoretically be equal to the actual value of the condition number.

Let us consider a trivial problem with a tridiagonal matrix
```

n = 10;
A = toeplitz (sparse ([1, 1], [1, 2], [2, 1], 1, n));
b = A * ones (n, 1);
M1 = ichol (A); \# in this tridiagonal case it corresponds to chol (A)'
M2 = M1';
M = M1 * M2;
Afun = @(x) A * x;
Mfun = @(x) M \ x;
M1fun = @(x) M1 \ x;
M2fun = @(x) M2 \ x;

```

Example 1: Simplest use of pcg
\(\mathrm{x}=\mathrm{pcg}(\mathrm{A}, \mathrm{b})\)
Example 2: pcg with a function which computes \(A * x\)
\(\mathrm{x}=\mathrm{pcg}\) (Afun, b)
Example 3: pcg with a preconditioner matrix \(M\)
\(\mathrm{x}=\operatorname{pcg}(\mathrm{A}, \mathrm{b}, 1 \mathrm{e}-06,100, \mathrm{M})\)
EXAMPLE 4: pcg with a function as preconditioner
\(x=p c g\) (Afun, b, 1e-6, 100, Mfun)
Example 5: pcg with preconditioner matrices M1 and M2
\[
\mathrm{x}=\mathrm{pcg}(\mathrm{~A}, \mathrm{~b}, 1 \mathrm{e}-6,100, \mathrm{M} 1, \mathrm{M} 2)
\]

EXAMPLE 6: pcg with functions as preconditioners
```

x = pcg (Afun, b, 1e-6, 100, M1fun, M2fun)

```

EXAMPLE 7: pcg with as input a function requiring an argument
```

        function y = Ap (A, x, p) # compute A^p * x
        y = x;
        for i = 1:p
            y = A * y;
        endfor
    endfunction
    Apfun = @(x, p) Ap (A, x, p);
x = pcg (Apfun, b, [], [], [], [], [], 2);

```

ExAmple 8: explicit example to show that pcg uses a split preconditioner
```

M1 = ichol (A + 0.1 * eye (n)); \# factorization of A perturbed
M2 = M1';
M = M1 * M2;

## reference solution computed by pcg after two iterations

[x_ref, fl] = pcg (A, b, [], 2, M)

## split preconditioning

[y, fl] = pcg ((M1 \ A) / M2, M1 \ b, [], 2)
x = M2 \ y \# compare x and x_ref

```

\section*{References:}
1. C.T. Kelley, Iterative Methods for Linear and Nonlinear Equations, SIAM, 1995. (the base PCG algorithm)
2. Y. Saad, Iterative Methods for Sparse Linear Systems, PWS 1996. (condition number estimate from PCG) Revised version of this book is available online at https://www-users.cs.umn.edu/~saad/books.html

See also: [sparse], page 664, [pcr], page 693, [gmres], page 617, [bicg], page 611, [bicgstab], page 613, [cgs], page 615.
\(x=\operatorname{pcr}(A, b, t o l, \operatorname{maxit}, m, x 0, \ldots)\)
[x, flag, relres, iter, resvec] = pcr (...)
Solve the linear system of equations \(A * x=b\) by means of the Preconditioned Conjugate Residuals iterative method.
The input arguments are
- A can be either a square (preferably sparse) matrix or a function handle, inline function or string containing the name of a function which computes \(A * x\). In principle \(A\) should be symmetric and non-singular; if per finds \(A\) to be numerically singular, you will get a warning message and the flag output parameter will be set.
- \(b\) is the right hand side vector.
- tol is the required relative tolerance for the residual error, \(b-A * x\). The iteration stops if norm \((b-A * x)<=\) tol \(*\) norm ( \(b-A * x 0\) ). If tol is empty or is omitted, the function sets tol \(=1 \mathrm{e}-6\) by default.
- maxit is the maximum allowable number of iterations; if [] is supplied for maxit, or pcr has less arguments, a default value equal to 20 is used.
- \(m\) is the (left) preconditioning matrix, so that the iteration is (theoretically) equivalent to solving by pcr \(P * \mathrm{x}=m \backslash \mathrm{~b}\), with \(P=m \backslash A\). Note that a proper choice of the preconditioner may dramatically improve the overall performance of the method. Instead of matrix \(m\), the user may pass a function which returns the results of applying the inverse of \(m\) to a vector (usually this is the preferred way of using the preconditioner). If [] is supplied for \(m\), or \(m\) is omitted, no preconditioning is applied.
- \(x 0\) is the initial guess. If \(x 0\) is empty or omitted, the function sets \(x 0\) to a zero vector by default.

The arguments which follow \(x 0\) are treated as parameters, and passed in a proper way to any of the functions ( \(A\) or \(m\) ) which are passed to pcr. See the examples below for further details.
The output arguments are
- \(x\) is the computed approximation to the solution of \(A * x=b\).
- flag reports on the convergence. flag \(=0\) means the solution converged and the tolerance criterion given by \(t o l\) is satisfied. \(f l a g=1\) means that the maxit limit for the iteration count was reached. flag \(=3\) reports a pcr breakdown, see [1] for details.
- relres is the ratio of the final residual to its initial value, measured in the Euclidean norm.
- iter is the actual number of iterations performed.
- resvec describes the convergence history of the method, so that resvec (i) contains the Euclidean norms of the residual after the ( \(i-1\) )-th iteration, \(i=1,2\), ..., itert.
Let us consider a trivial problem with a diagonal matrix (we exploit the sparsity of A)
```

n = 10;
A = sparse (diag (1:n));
b = rand (N, 1);

```

Example 1: Simplest use of pcr
\[
\mathrm{x}=\operatorname{pcr}(\mathrm{A}, \mathrm{~b})
\]

Example 2: pcr with a function which computes \(A * x\).
```

function y = apply_a (x)
y = [1:10]' .* x;
endfunction
x = pcr ("apply_a", b)

```

Example 3: Preconditioned iteration, with full diagnostics. The preconditioner (quite strange, because even the original matrix \(A\) is trivial) is defined as a function
```

function y = apply_m (x)
k = floor (length (x) - 2);
y = x;
y(1:k) = x(1:k) ./ [1:k]';
endfunction
[x, flag, relres, iter, resvec] = ...
pcr (A, b, [], [], "apply_m")
semilogy ([1:iter+1], resvec);

```

Example 4: Finally, a preconditioner which depends on a parameter \(k\).
```

function $y=$ apply_m (x, varargin)
$\mathrm{k}=\operatorname{varargin}\{1\} ;$
$\mathrm{y}=\mathrm{x}$;
$y(1: k)=x(1: k) . /[1: k] ' ;$
endfunction
[x, flag, relres, iter, resvec] = ...
pcr (A, b, [], [], "apply_m"', [], 3)

```

Reference:
W. Hackbusch, Iterative Solution of Large Sparse Systems of Equations, section 9.5.4; Springer, 1994
See also: [sparse], page 664, [pcg], page 690.

The speed with which an iterative solver converges to a solution can be accelerated with the use of a pre-conditioning matrix \(M\). In this case the linear equation \(M^{\wedge}-1 * x=M^{\wedge}-1 *\) \(A \backslash b\) is solved instead. Typical pre-conditioning matrices are partial factorizations of the original matrix.
```

L = ichol (A)
L = ichol (A, opts)

```

Compute the incomplete Cholesky factorization of the sparse square matrix \(A\).
By default, ichol uses only the lower triangle of \(A\) and produces a lower triangular factor \(L\) such that \(\mathrm{L} * \mathrm{~L}\) ' approximates \(A\).
The factor given by this routine may be useful as a preconditioner for a system of linear equations being solved by iterative methods such as PCG (Preconditioned Conjugate Gradient).
The factorization may be modified by passing options in a structure opts. The option name is a field of the structure and the setting is the value of field. Names and specifiers are case sensitive.
type Type of factorization.
"nofill" (default)
Incomplete Cholesky factorization with no fill-in (IC(0)).
"ict" Incomplete Cholesky factorization with threshold dropping (ICT).
diagcomp A non-negative scalar alpha for incomplete Cholesky factorization of \(A+\) alpha * diag (diag (A)) instead of \(A\). This can be useful when \(A\) is not positive definite. The default value is 0 .
droptol A non-negative scalar specifying the drop tolerance for factorization if performing ICT. The default value is 0 which produces the complete Cholesky factorization.
Non-diagonal entries of \(L\) are set to 0 unless
abs \((L(i, j))>=\) droptol \(* \operatorname{norm}(A(j:\) end,\(j), 1)\).
michol Modified incomplete Cholesky factorization:
"off" (default)
Row and column sums are not necessarily preserved.
"on" The diagonal of \(L\) is modified so that row (and column) sums are preserved even when elements have been dropped during the factorization. The relationship preserved is: \(A * e=L *\) \(L^{\prime} * e\), where e is a vector of ones.
shape
\[
\begin{aligned}
& \text { "lower" (default) } \\
& \quad \text { Use only the lower triangle of } A \text { and return a lower triangular } \\
& \text { factor } L \text { such that } \mathrm{L} * \mathrm{~L} \text { ' approximates } A \text {. } \\
& \text { "upper" } \quad \begin{array}{l}
\text { Use only the upper triangle of } A \text { and return an upper trian- } \\
\text { gular factor } U \text { such that } \mathrm{U} ' * \mathrm{U} \text { approximates } A .
\end{array}
\end{aligned}
\]

\section*{EXAMPLES}

The following problem demonstrates how to factorize a sample symmetric positive definite matrix with the full Cholesky decomposition and with the incomplete one.
```

A = [ 0.37, -0.05, -0.05, -0.07;
-0.05, 0.116, 0.0, -0.05;
-0.05, 0.0, 0.116, -0.05;
-0.07, -0.05, -0.05, 0.202];
A = sparse (A);
nnz (tril (A))
ans = 9
L = chol (A, "lower");
nnz (L)
ans = 10
norm (A - L * L', "fro") / norm (A, "fro")
ans = 1.1993e-16
opts.type = "nofill";
L = ichol (A, opts);
nnz (L)
ans = 9
norm (A - L * L', "fro") / norm (A, "fro")
ans = 0.019736

```

Another example for decomposition is a finite difference matrix used to solve a boundary value problem on the unit square.
```

nx = 400; ny = 200;
hx = 1 / (nx + 1); hy = 1 / (ny + 1);
Dxx = spdiags ([ones(nx, 1), -2*ones(nx, 1), ones(nx, 1)],
[-1 0 1 ], nx, nx) / (hx ` 2);
Dyy = spdiags ([ones(ny, 1), -2*ones(ny, 1), ones(ny, 1)],
[-1 0 1 ], ny, ny) / (hy ~ 2);
A = -kron (Dxx, speye (ny)) - kron (speye (nx), Dyy);
nnz (tril (A))
ans = 239400
opts.type = "nofill";
L = ichol (A, opts);
nnz (tril (A))
ans = 239400
norm (A - L * L', "fro") / norm (A, "fro")
ans = 0.062327

```

References for implemented algorithms:
[1] Y. Saad. "Preconditioning Techniques." Iterative Methods for Sparse Linear Systems, PWS Publishing Company, 1996.
[2] M. Jones, P. Plassmann: An Improved Incomplete Cholesky Factorization, 1992.
See also: [chol], page 595, [ilu], page 696, [pcg], page 690.
ilu (A)
ilu (A, opts)
```

[L,U] = ilu (...)
[L, U, P] = ilu (...)

```

Compute the incomplete LU factorization of the sparse square matrix \(A\).
ilu returns a unit lower triangular matrix \(L\), an upper triangular matrix \(U\), and optionally a permutation matrix \(P\), such that \(L * U\) approximates \(P * A\).
The factors given by this routine may be useful as preconditioners for a system of linear equations being solved by iterative methods such as BICG (BiConjugate Gradients) or GMRES (Generalized Minimum Residual Method).
The factorization may be modified by passing options in a structure opts. The option name is a field of the structure and the setting is the value of field. Names and specifiers are case sensitive.
type Type of factorization.
"nofill" (default)
ILU factorization with no fill-in (ILU(0)).
Additional supported options: milu.
"crout" Crout version of ILU factorization (ILUC).
Additional supported options: milu, droptol.
"ilutp" ILU factorization with threshold and pivoting.
Additional supported options: milu, droptol, udiag, thresh.
droptol A non-negative scalar specifying the drop tolerance for factorization. The default value is 0 which produces the complete LU factorization.
Non-diagonal entries of \(U\) are set to 0 unless
abs ( \(U(i, j)\) ) >= droptol \(* \operatorname{norm}(A(:, j))\).
Non-diagonal entries of \(L\) are set to 0 unless
abs \((L(i, j))>=d r o p t o l * \operatorname{norm}(A(:, j)) / U(j, j)\).
milu Modified incomplete LU factorization:
"row" Row-sum modified incomplete LU factorization. The factorization preserves row sums: \(A * e=L * U * e\), where \(e\) is a vector of ones.
"col" Column-sum modified incomplete LU factorization. The factorization preserves column sums: \(\mathrm{e}^{\prime} * A=\mathrm{e}^{\prime} * L * U\).
"off" (default)
Row and column sums are not necessarily preserved.
udiag If true, any zeros on the diagonal of the upper triangular factor are replaced by the local drop tolerance droptol \(* \operatorname{norm}(A(:, j)) / U(j, j)\). The default is false.
thresh Pivot threshold for factorization. It can range between 0 (diagonal pivoting) and 1 (default), where the maximum magnitude entry in the column is chosen to be the pivot.

If ilu is called with just one output, the returned matrix is \(L+U\) - speye (size (A)), where \(L\) is unit lower triangular and \(U\) is upper triangular.

With two outputs, ilu returns a unit lower triangular matrix \(L\) and an upper triangular matrix \(U\). For opts.type \(==\) "ilutp", one of the factors is permuted based on the value of opts.milu. When opts.milu \(==\) "row", \(U\) is a column permuted upper triangular factor. Otherwise, \(L\) is a row-permuted unit lower triangular factor.
If there are three named outputs and opts.milu \(!=\) "row", \(P\) is returned such that \(L\) and \(U\) are incomplete factors of \(P * A\). When opts.milu \(==\) "row", \(P\) is returned such that \(L\) and \(U\) are incomplete factors of \(A * P\).
EXAMPLES
```

A = gallery ("neumann", 1600) + speye (1600);
opts.type = "nofill";
nnz (A)
ans = 7840
nnz (lu (A))
ans = 126478
nnz (ilu (A, opts))
ans = 7840

```

This shows that \(A\) has 7,840 nonzeros, the complete LU factorization has 126,478 nonzeros, and the incomplete LU factorization, with 0 level of fill-in, has 7,840 nonzeros, the same amount as \(A\). Taken from: https://www.mathworks.com/help/ matlab/ref/ilu.html
```

A = gallery ("wathen", 10, 10);
b = sum (A, 2);
tol = 1e-8;
maxit = 50;
opts.type = "crout";
opts.droptol = 1e-4;
[L, U] = ilu (A, opts);
x = bicg (A, b, tol, maxit, L, U);
norm (A * x - b, inf)

```

This example uses ILU as preconditioner for a random FEM-Matrix, which has a large condition number. Without \(L\) and \(U\) BICG would not converge.
See also: [lu], page 597, [ichol], page 695, [bicg], page 611, [gmres], page 617.

\subsection*{22.4 Real Life Example using Sparse Matrices}

A common application for sparse matrices is in the solution of Finite Element Models. Finite element models allow numerical solution of partial differential equations that do not have closed form solutions, typically because of the complex shape of the domain.

In order to motivate this application, we consider the boundary value Laplace equation. This system can model scalar potential fields, such as heat or electrical potential. Given a medium \(\Omega\) with boundary \(\partial \Omega\). At all points on the \(\partial \Omega\) the boundary conditions are
known, and we wish to calculate the potential in \(\Omega\). Boundary conditions may specify the potential (Dirichlet boundary condition), its normal derivative across the boundary (Neumann boundary condition), or a weighted sum of the potential and its derivative (Cauchy boundary condition).

In a thermal model, we want to calculate the temperature in \(\Omega\) and know the boundary temperature (Dirichlet condition) or heat flux (from which we can calculate the Neumann condition by dividing by the thermal conductivity at the boundary). Similarly, in an electrical model, we want to calculate the voltage in \(\Omega\) and know the boundary voltage (Dirichlet) or current (Neumann condition after diving by the electrical conductivity). In an electrical model, it is common for much of the boundary to be electrically isolated; this is a Neumann boundary condition with the current equal to zero.

The simplest finite element models will divide \(\Omega\) into simplexes (triangles in 2D, pyramids in 3D). We take as a 3-D example a cylindrical liquid filled tank with a small non-conductive ball from the EIDORS project \({ }^{4}\). This is model is designed to reflect an application of electrical impedance tomography, where current patterns are applied to such a tank in order to image the internal conductivity distribution. In order to describe the FEM geometry, we have a matrix of vertices nodes and simplices elems.

The following example creates a simple rectangular 2-D electrically conductive medium with 10 V and 20 V imposed on opposite sides (Dirichlet boundary conditions). All other edges are electrically isolated.
```

node_y = [1;1.2;1.5;1.8;2]*ones(1,11);
node_x = ones(5,1)*[1,1.05,1.1,1.2, ...
1.3,1.5,1.7,1.8,1.9,1.95,2];
nodes = [node_x(:), node_y(:)];
[h,w] = size (node_x);
elems = [];
for idx = 1:w-1
widx = (idx-1)*h;
elems = [elems; ...
widx+[(1:h-1);(2:h);h+(1:h-1)]'; ...
widx+[(2:h);h+(2:h);h+(1:h-1)]' ];
endfor
E = size (elems,1); \# No. of simplices
N = size (nodes,1); \# No. of vertices
D = size (elems,2); \# dimensions+1

```

This creates a N-by-2 matrix nodes and a E-by-3 matrix elems with values, which define finite element triangles:

\footnotetext{
\({ }^{4}\) EIDORS - Electrical Impedance Tomography and Diffuse optical Tomography Reconstruction Software http://eidors3d.sourceforge.net
}
```

nodes(1:7,:)'
1.00 1.00 1.00 1.00 1.00 1.05 1.05 ...
1.00 1.20 1.50 1.80 2.00 1.00 1.20 ...

```
elems(1:7,:)'
    \(1 \begin{array}{lllllll} & 2 & 3 & 4 & 2 & 3\end{array}\)
    \(\begin{array}{lllllll}2 & 3 & 4 & 5 & 7 & 8 & 9\end{array}\)
    \(\begin{array}{lllllll}6 & 7 & 8 & 9 & 6 & 7 & 8\end{array}\)

Using a first order FEM, we approximate the electrical conductivity distribution in \(\Omega\) as constant on each simplex (represented by the vector conductivity). Based on the finite element geometry, we first calculate a system (or stiffness) matrix for each simplex (represented as 3 -by- 3 elements on the diagonal of the element-wise system matrix SE ). Based on SE and a N-by-DE connectivity matrix C, representing the connections between simplices and vertices, the global connectivity matrix \(S\) is calculated.
```


## Element conductivity

conductivity = [1*ones (1,16), ...
2*ones (1,48), 1*ones (1,16)];

## Connectivity matrix

C = sparse ((1:D*E), reshape (elems', ...
D*E, 1), 1, D*E, N);

## Calculate system matrix

Siidx = floor ([0:D*E-1]'/D) * D * ...
ones(1,D) + ones(D*E,1)*(1:D) ;
Sjidx = [1:D*E]'*ones (1,D);
Sdata = zeros (D*E,D);
dfact = factorial (D-1);
for j = 1:E
a = inv ([ones(D,1), ...
nodes(elems(j,:), :)]);
const = conductivity(j) * 2 / ...
dfact / abs (det (a));
Sdata(D*(j-1)+(1:D),:) = const * ...
a(2:D,:)' * a(2:D,:);
endfor

## Element-wise system matrix

SE = sparse(Siidx,Sjidx,Sdata);

## Global system matrix

S = C'* SE *C;

```

The system matrix acts like the conductivity \(S\) in Ohm's law \(S V=I\). Based on the Dirichlet and Neumann boundary conditions, we are able to solve for the voltages at each vertex V.
```


## Dirichlet boundary conditions

D_nodes = [1:5, 51:55];
D_value = [10*ones (1,5), 20*ones (1,5)];

```
```

V = zeros (N,1);
V(D_nodes) = D_value;
idx = 1:N; \# vertices without Dirichlet
\# boundary condns
idx(D_nodes) = [];

## Neumann boundary conditions. Note that

## N_value must be normalized by the

## boundary length and element conductivity

N_nodes = [];
N_value = [];
Q = zeros (N,1);
Q(N_nodes) = N_value;
V(idx) = S(idx,idx) \ ( Q(idx) - ...
S(idx,D_nodes) * V(D_nodes));

```

Finally, in order to display the solution, we show each solved voltage value in the \(z\)-axis for each simplex vertex. See Figure 22.6.
```

elemx = elems(:,[1,2,3,1])';
xelems = reshape (nodes(elemx, 1), 4, E);
yelems = reshape (nodes(elemx, 2), 4, E);
velems = reshape (V(elemx), 4, E);
plot3 (xelems,yelems,velems,"k");
print "grid.eps";

```


Figure 22.6: Example finite element model the showing triangular elements. The height of each vertex corresponds to the solution value.

\section*{23 Numerical Integration}

Octave comes with several built-in functions for computing the integral of a function numerically (termed quadrature). These functions all solve 1-dimensional integration problems.

\subsection*{23.1 Functions of One Variable}

Octave supports five different adaptive quadrature algorithms for computing the integral
\[
\int_{a}^{b} f(x) d x
\]
of a function \(f\) over the interval from \(a\) to \(b\). These are
quad Numerical integration based on Gaussian quadrature.
quadv Numerical integration using an adaptive vectorized Simpson's rule.
quadl Numerical integration using an adaptive Lobatto rule.
quadgk Numerical integration using an adaptive Gauss-Konrod rule.
quadcc Numerical integration using adaptive Clenshaw-Curtis rules.
In addition, the following functions are also provided:
integral A compatibility wrapper function that will choose between quadv and quadgk depending on the integrand and options chosen.
```

trapz, cumtrapz

```

Numerical integration of data using the trapezoidal method.
The best quadrature algorithm to use depends on the integrand. If you have empirical data, rather than a function, the choice is trapz or cumtrapz. If you are uncertain about the characteristics of the integrand, quadcc will be the most robust as it can handle discontinuities, singularities, oscillatory functions, and infinite intervals. When the integrand is smooth quadgk may be the fastest of the algorithms.

\section*{Function Characteristics}
quad Low accuracy with nonsmooth integrands
quadv Medium accuracy with smooth integrands
quadl Medium accuracy with smooth integrands. Slower than quadgk.
quadgk Medium accuracy (1e-6 - 1e-9) with smooth integrands.
Handles oscillatory functions and infinite bounds
quadcc Low to High accuracy with nonsmooth/smooth integrands Handles oscillatory functions, singularities, and infinite bounds
Here is an example of using quad to integrate the function
\[
f(x)=x \sin (1 / x) \sqrt{|1-x|}
\]
from \(x=0\) to \(x=3\).
This is a fairly difficult integration (plot the function over the range of integration to see why).

The first step is to define the function:
```

function y = f (x)
y = x .* sin (1./x) .* sqrt (abs (1 - x));
endfunction

```

Note the use of the 'dot' forms of the operators. This is not necessary for the quad integrator, but is required by the other integrators. In any case, it makes it much easier to generate a set of points for plotting because it is possible to call the function with a vector argument to produce a vector result.

The second step is to call quad with the limits of integration:
```

[q, ier, nfun, err] = quad ("f", 0, 3)
=> 1.9819
=>1
F 5061
\# 1.1522e-07

```

Although quad returns a nonzero value for ier, the result is reasonably accurate (to see why, examine what happens to the result if you move the lower bound to 0.1 , then 0.01 , then 0.001, etc.).

The function "f" can be the string name of a function or a function handle. These options make it quite easy to do integration without having to fully define a function in an m-file. For example:
```


# Verify gamma function = (n-1)! for n = 4

f = @(x) x.^3 .* exp (-x);
quadcc (f, 0, Inf)
=>6.0000

```
```

q= quad (f,a,b)
q= quad (f,a,b,tol)
q = quad (f, a, b, tol, sing)
[q, ier, nfun, err] = quad (...)

```

Numerically evaluate the integral of \(f\) from a to \(b\) using Fortran routines from QUADPACK.
\(f\) is a function handle, inline function, or a string containing the name of the function to evaluate. The function must have the form \(y=f(x)\) where \(y\) and \(x\) are scalars.
\(a\) and \(b\) are the lower and upper limits of integration. Either or both may be infinite.
The optional argument tol is a vector that specifies the desired accuracy of the result. The first element of the vector is the desired absolute tolerance, and the second element is the desired relative tolerance. To choose a relative test only, set the absolute tolerance to zero. To choose an absolute test only, set the relative tolerance to zero. Both tolerances default to sqrt (eps) or approximately \(1.5 \mathrm{e}-8\).

The optional argument sing is a vector of values at which the integrand is known to be singular.
The result of the integration is returned in \(q\).
ier contains an integer error code ( 0 indicates a successful integration).
nfun indicates the number of function evaluations that were made.
err contains an estimate of the error in the solution.
The function quad_options can set other optional parameters for quad.
Note: because quad is written in Fortran it cannot be called recursively. This prevents its use in integrating over more than one variable by routines dblquad and triplequad.
See also: [quad_options], page 705, [quadv], page 705, [quadl], page 706, [quadgk], page 707, [quadcc], page 708, [trapz], page 711, [dblquad], page 713, [triplequad], page 713 .
quad_options ()
val = quad_options (opt)
quad_options (opt, val)
Query or set options for the function quad.
When called with no arguments, the names of all available options and their current values are displayed.
Given one argument, return the value of the option opt.
When called with two arguments, quad_options sets the option opt to value val.
Options include
"absolute tolerance"
Absolute tolerance; may be zero for pure relative error test.
"relative tolerance"
Non-negative relative tolerance. If the absolute tolerance is zero, the relative tolerance must be greater than or equal to max ( \(50 * \mathrm{eps}, 0.5 \mathrm{e}-28\) ).
"single precision absolute tolerance"
Absolute tolerance for single precision; may be zero for pure relative error test.
"single precision relative tolerance"
Non-negative relative tolerance for single precision. If the absolute tolerance is zero, the relative tolerance must be greater than or equal to \(\max\) ( \(50 * e \mathrm{eps}, 0.5 \mathrm{e}-28\) ).
\(q=\operatorname{quadv}(f, a, b)\)
\(q=\) quadv ( \(f, a, b, t o l\) )
\(q=\) quadv ( \(f, a, b\), tol, trace)
\(q=\) quadv ( \(f, a, b\), tol, trace, \(p 1, p 2, \ldots\) )
[q, nfun] = quadv (...)
Numerically evaluate the integral of \(f\) from a to \(b\) using an adaptive Simpson's rule.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. quadv is a vectorized version of quad and the function defined by \(f\) must accept a scalar or vector as input and return a scalar, vector, or array as output.
\(a\) and \(b\) are the lower and upper limits of integration. Both limits must be finite.
The optional argument tol defines the absolute tolerance used to stop the adaptation procedure. The default value is \(1 \mathrm{e}-6\).

The algorithm used by quadv involves recursively subdividing the integration interval and applying Simpson's rule on each subinterval. If trace is true then after computing each of these partial integrals display: (1) the total number of function evaluations, (2) the left end of the subinterval, (3) the length of the subinterval, (4) the approximation of the integral over the subinterval.
Additional arguments \(p 1\), etc., are passed directly to the function \(f\). To use default values for tol and trace, one may pass empty matrices ([]).

The result of the integration is returned in \(q\).
The optional output nfun indicates the total number of function evaluations performed.
Note: quadv is written in Octave's scripting language and can be used recursively in dblquad and triplequad, unlike the quad function.

See also: [quad], page 704, [quadl], page 706, [quadgk], page 707, [quadcc], page 708, [trapz], page 711, [dblquad], page 713, [triplequad], page 713, [integral], page 709, [integral2], page 715, [integral3], page 717.
```

q = quadl (f,a,b)
q = quadl (f, a, b, tol)
q = quadl (f, a, b, tol, trace)
q = quadl (f, a, b, tol, trace, p1, p2,...)
[q, nfun] = quadl (...)

```

Numerically evaluate the integral of \(f\) from \(a\) to \(b\) using an adaptive Lobatto rule.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be vectorized and return a vector of output values when given a vector of input values.
\(a\) and \(b\) are the lower and upper limits of integration. Both limits must be finite.
The optional argument tol defines the absolute tolerance with which to perform the integration. The default value is \(1 \mathrm{e}-6\).
The algorithm used by quadl involves recursively subdividing the integration interval. If trace is defined then for each subinterval display: (1) the total number of function evaluations, (2) the left end of the subinterval, (3) the length of the subinterval, (4) the approximation of the integral over the subinterval.
Additional arguments \(p 1\), etc., are passed directly to the function \(f\). To use default values for tol and trace, one may pass empty matrices ([]).
The result of the integration is returned in \(q\).
The optional output nfun indicates the total number of function evaluations performed.
Reference: W. Gander and W. Gautschi, Adaptive Quadrature - Revisited, BIT Vol. 40, No. 1, March 2000, pp. 84-101. https://www.inf.ethz.ch/personal/gander/
See also: [quad], page 704, [quadv], page 705, [quadgk], page 707, [quadcc], page 708, [trapz], page 711, [dblquad], page 713, [triplequad], page 713, [integral], page 709, [integral2], page 715, [integral3], page 717.
```

q = quadgk (f, a, b)
q = quadgk (f, a, b, abstol)
q = quadgk (f, a, b, abstol, trace)
q = quadgk (f, a, b, prop, val, ...)
[q, err] = quadgk (...)

```

Numerically evaluate the integral of \(f\) from a to \(b\) using adaptive Gauss-Kronrod quadrature.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be vectorized and return a vector of output values when given a vector of input values.
\(a\) and \(b\) are the lower and upper limits of integration. Either or both limits may be infinite or contain weak end singularities. Variable transformation will be used to treat any infinite intervals and weaken the singularities. For example:
quadgk (@(x) 1./ (sqrt (x) .* (x + 1)), 0, Inf)
Note that the formulation of the integrand uses the element-by-element operator ./ and all user functions to quadgk should do the same.
The optional argument tol defines the absolute tolerance used to stop the integration procedure. The default value is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).
The algorithm used by quadgk involves subdividing the integration interval and evaluating each subinterval. If trace is true then after computing each of these partial integrals display: (1) the number of subintervals at this step, (2) the current estimate of the error err, (3) the current estimate for the integral \(q\).
The behavior of the algorithm can be configured by passing arguments to quadgk as pairs "prop", val. Valid properties are
AbsTol Define the absolute error tolerance for the quadrature. The default absolute tolerance is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).

RelTol Define the relative error tolerance for the quadrature. The default relative tolerance is 1e-6 (1e-4 for single).

\section*{MaxIntervalCount}
quadgk initially subdivides the interval on which to perform the quadrature into 10 intervals. Subintervals that have an unacceptable error are subdivided and re-evaluated. If the number of subintervals exceeds 650 subintervals at any point then a poor convergence is signaled and the current estimate of the integral is returned. The property "MaxIntervalCount" can be used to alter the number of subintervals that can exist before exiting.

\section*{WayPoints}

Discontinuities in the first derivative of the function to integrate can be flagged with the "WayPoints" property. This forces the ends of a subinterval to fall on the breakpoints of the function and can result in significantly improved estimation of the error in the integral, faster computation, or both. For example,
```

quadgk (@(x) abs (1 - x.^2), 0, 2, "Waypoints", 1)

```
signals the breakpoint in the integrand at \(x=1\).

Trace If logically true quadgk prints information on the convergence of the quadrature at each iteration.

If any of \(a, b\), or waypoints is complex then the quadrature is treated as a contour integral along a piecewise continuous path defined by [a, waypoints(1), waypoints(2) , ..., b]. In this case the integral is assumed to have no edge singularities. For example,
```

quadgk (@(z) log (z), 1+1i, 1+1i, "WayPoints",
[-1+1i, -1-1i, +1-1i])

```
integrates \(\log (z)\) along the square defined by \([1+1 i,-1+1 i,-1-1 i,+1-1 i]\).
The result of the integration is returned in \(q\).
err is an approximate bound on the error in the integral abs ( \(q-I\) ), where \(I\) is the exact value of the integral.
Reference: L.F. Shampine, " Vectorized adaptive quadrature in matlab", Journal of Computational and Applied Mathematics, pp. 131-140, Vol 211, Issue 2, Feb 2008.

See also: [quad], page 704, [quadv], page 705, [quadl], page 706, [quadcc], page 708, [trapz], page 711, [dblquad], page 713, [triplequad], page 713, [integral], page 709, [integral2], page 715, [integral3], page 717 .
\(q=\) quadcc \((f, a, b)\)
\(q=\) quadcc ( \(f, a, b, t o l\) )
\(q=\) quadcc ( \(f, a, b\), tol, sing)
[q, err, nr_points] = quadcc (...)
Numerically evaluate the integral of \(f\) from a to \(b\) using doubly-adaptive ClenshawCurtis quadrature.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be vectorized and must return a vector of output values if given a vector of input values. For example,
```

f = @(x) x .* sin (1./x) .* sqrt (abs (1 - x));

```
which uses the element-by-element "dot" form for all operators.
\(a\) and \(b\) are the lower and upper limits of integration. Either or both limits may be infinite. quadcc handles an infinite limit by substituting the variable of integration with \(\mathrm{x}=\tan (\mathrm{pi} / 2 * \mathrm{u})\).
The optional argument tol is a 1 - or 2-element vector that specifies the desired accuracy of the result. The first element of the vector is the desired absolute tolerance, and the second element is the desired relative tolerance. To choose a relative test only, set the absolute tolerance to zero. To choose an absolute test only, set the relative tolerance to zero. The default absolute tolerance is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single), and the default relative tolerance is \(1 \mathrm{e}-6\) ( \(1 \mathrm{e}-4\) for single).
The optional argument sing contains a list of points where the integrand has known singularities, or discontinuities in any of its derivatives, inside the integration interval. For the example above, which has a discontinuity at \(\mathrm{x}=1\), the call to quadcc would be as follows
```

int = quadcc (f, a, b, [], [ 1 ]);

```

The result of the integration is returned in \(q\).
err is an estimate of the absolute integration error.
\(n r_{-}\)points is the number of points at which the integrand was evaluated.
If the adaptive integration did not converge, the value of err will be larger than the requested tolerance. Therefore, it is recommended to verify this value for difficult integrands.
quadcc is capable of dealing with non-numeric values of the integrand such as NaN or Inf. If the integral diverges, and quadcc detects this, then a warning is issued and Inf or -Inf is returned.
Note: quadcc is a general purpose quadrature algorithm and, as such, may be less efficient for a smooth or otherwise well-behaved integrand than other methods such as quadgk.
The algorithm uses Clenshaw-Curtis quadrature rules of increasing degree in each interval and bisects the interval if either the function does not appear to be smooth or a rule of maximum degree has been reached. The error estimate is computed from the L2-norm of the difference between two successive interpolations of the integrand over the nodes of the respective quadrature rules.
Reference: P. Gonnet, Increasing the Reliability of Adaptive Quadrature Using Explicit Interpolants, ACM Transactions on Mathematical Software, Vol. 37, Issue 3, Article No. 3, 2010.

See also: [quad], page 704, [quadv], page 705, [quadl], page 706, [quadgk], page 707, [trapz], page 711, [dblquad], page 713, [triplequad], page 713.
\(q=\) integral (f, \(a, b)\)
\(q=\) integral ( \(f, a, b\), prop, val, ...)
Numerically evaluate the integral of \(f\) from a to \(b\) using adaptive quadrature.
integral is a wrapper for quadcc (general real-valued, scalar integrands and limits), quadgk (integrals with specified integration paths), and quadv (array-valued integrands) that is intended to provide matlab compatibility. More control of the numerical integration may be achievable by calling the various quadrature functions directly.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be vectorized and return a vector of output values when given a vector of input values.
\(a\) and \(b\) are the lower and upper limits of integration. Either or both limits may be infinite or contain weak end singularities. If either or both limits are complex, integral will perform a straight line path integral. Alternatively, a complex domain path can be specified using the "Waypoints" option (see below).
Additional optional parameters can be specified using "property", value pairs. Valid properties are:

\section*{Waypoints}

Specifies points to be used in defining subintervals of the quadrature algorithm, or if \(a, b\), or waypoints are complex then the quadrature is calculated as a contour integral along a piecewise continuous path. For more detail, see [quadgk], page 707.

\section*{ArrayValued}
integral expects \(f\) to return a scalar value unless arrayvalued is specified as true. This option will cause integral to perform the integration over the entire array and return \(q\) with the same dimensions as returned by \(f\). For more detail see [quadv], page 705 .
AbsTol Define the absolute error tolerance for the quadrature. The default absolute tolerance is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).
RelTol Define the relative error tolerance for the quadrature. The default relative tolerance is 1e-6 (1e-4 for single).
Adaptive quadrature is used to minimize the estimate of error until the following is satisfied:
\[
\text { error } \leq \max (\text { AbsTol, RelTol } \cdot|q|)
\]

Known matlab incompatibilities:
1. If tolerances are left unspecified, and any integration limits or waypoints are of type single, then Octave's integral functions automatically reduce the default absolute and relative error tolerances as specified above. If tighter tolerances are desired they must be specified. matlab leaves the tighter tolerances appropriate for double inputs in place regardless of the class of the integration limits.
2. As a consequence of using quadcc, quadgk, and quadv, certain option combinations are not supported. Currently, "ArrayValued" cannot be combined with "RelTol" or "Waypoints".

See also: [integral2], page 715, [integral3], page 717, [quad], page 704, [quadgk], page 707, [quadv], page 705, [quadl], page 706, [quadcc], page 708, [trapz], page 711, [dblquad], page 713, [triplequad], page 713.

Sometimes one does not have the function, but only the raw ( \(\mathrm{x}, \mathrm{y}\) ) points from which to perform an integration. This can occur when collecting data in an experiment. The trapz function can integrate these values as shown in the following example where "data" has been collected on the cosine function over the range \([0, \mathrm{pi} / 2)\).
```

x = 0:0.1:pi/2; \# Uniformly spaced points
y = cos (x);
trapz (x, y)
\# 0.99666

```

The answer is reasonably close to the exact value of 1 . Ordinary quadrature is sensitive to the characteristics of the integrand. Empirical integration depends not just on the integrand, but also on the particular points chosen to represent the function. Repeating the example above with the sine function over the range \([0, \mathrm{pi} / 2)\) produces far inferior results.
```

x = 0:0.1:pi/2; \# Uniformly spaced points
y = sin (x);
trapz (x, y)
=> 0.92849

```

However, a slightly different choice of data points can change the result significantly. The same integration, with the same number of points, but spaced differently produces a more accurate answer.
```

x = linspace (0, pi/2, 16); \# Uniformly spaced, but including endpoint
y = sin (x);
trapz (x, y)
=> 0.99909

```

In general there may be no way of knowing the best distribution of points ahead of time. Or the points may come from an experiment where there is no freedom to select the best distribution. In any case, one must remain aware of this issue when using trapz.
```

q = trapz (y)
q = trapz (x, y)
q = trapz (..., dim)

```

Numerically evaluate the integral of points \(y\) using the trapezoidal method.
trapz ( \(y\) ) computes the integral of \(y\) along the first non-singleton dimension. When the argument \(x\) is omitted an equally spaced \(x\) vector with unit spacing (1) is assumed. trapz ( \(x, y\) ) evaluates the integral with respect to the spacing in \(x\) and the values in \(y\). This is useful if the points in \(y\) have been sampled unevenly.
If the optional dim argument is given, operate along this dimension.
Application Note: If \(x\) is not specified then unit spacing will be used. To scale the integral to the correct value you must multiply by the actual spacing value (deltaX). As an example, the integral of \(x^{3}\) over the range \([0,1]\) is \(x^{4} / 4\) or 0.25 . The following code uses trapz to calculate the integral in three different ways.
```

x = 0:0.1:1;
y = x.^3;

## No scaling

q = trapz (y)
\#q=2.5250

## Approximation to integral by scaling

q * 0.1
=> 0.25250

## Same result by specifying x

trapz (x, y)
\# 0.25250

```

See also: [cumtrapz], page 711 .
\(q=\) cumtrapz ( \(y\) )
\(q=\) cumtrapz ( \(x, y\) )
\(q=\) cumtrapz (..., dim)
Cumulative numerical integration of points \(y\) using the trapezoidal method.
cumtrapz ( \(y\) ) computes the cumulative integral of \(y\) along the first non-singleton dimension. Where trapz reports only the overall integral sum, cumtrapz reports the current partial sum value at each point of \(y\).
When the argument \(x\) is omitted an equally spaced \(x\) vector with unit spacing (1) is assumed. cumtrapz ( \(x, y\) ) evaluates the integral with respect to the spacing in \(x\) and the values in \(y\). This is useful if the points in \(y\) have been sampled unevenly.
If the optional dim argument is given, operate along this dimension.

Application Note: If \(x\) is not specified then unit spacing will be used. To scale the integral to the correct value you must multiply by the actual spacing value (deltaX).

See also: [trapz], page 711, [cumsum], page 556.

\subsection*{23.2 Orthogonal Collocation}
[r, amat, bmat, q] = colloc (n, "left", "right")
Compute derivative and integral weight matrices for orthogonal collocation.
Reference: J. Villadsen, M. L. Michelsen, Solution of Differential Equation Models by Polynomial Approximation.

Here is an example of using colloc to generate weight matrices for solving the second order differential equation \(u^{\prime}-\alpha u^{\prime \prime}=0\) with the boundary conditions \(u(0)=0\) and \(u(1)=1\).

First, we can generate the weight matrices for \(n\) points (including the endpoints of the interval), and incorporate the boundary conditions in the right hand side (for a specific value of \(\alpha\) ).
```

n = 7;
alpha = 0.1;
[r, a, b] = colloc (n-2, "left", "right");
at = a(2:n-1,2:n-1);
bt = b(2:n-1,2:n-1);
rhs = alpha * b(2:n-1,n) - a(2:n-1,n);

```

Then the solution at the roots \(r\) is
```

u = [ 0; (at - alpha * bt) \ rhs; 1]
[ [0.00; 0.004; 0.01 0.00; 0.12; 0.62;1.00 ]

```

\subsection*{23.3 Functions of Multiple Variables}

Octave includes several functions for computing the integral of functions of multiple variables. This procedure can generally be performed by creating a function that integrates \(f\) with respect to \(x\), and then integrates that function with respect to \(y\). This procedure can be performed manually using the following example which integrates the function:
\[
f(x, y)=\sin (\pi x y) \sqrt{x y}
\]
for \(x\) and \(y\) between 0 and 1 .
Using quadgk in the example below, a double integration can be performed. (Note that any of the 1-D quadrature functions can be used in this fashion except for quad since it is written in Fortran and cannot be called recursively.)
```

function q = g(y)
q = ones (size (y));
for i = 1:length (y)
f = @(x) sin (pi*x.*y(i)) .* sqrt (x.*y(i));
q(i) = quadgk (f, 0, 1);
endfor
endfunction
I = quadgk ("g", 0, 1)
=> 0.30022

```

The algorithm above is implemented in the function dblquad for integrals over two variables. The 3-D equivalent of this process is implemented in triplequad for integrals over three variables. As an example, the result above can be replicated with a call to dblquad as shown below.
```

I = dblquad (@(x, y) sin (pi*x.*y) .* sqrt (x.*y), 0, 1, 0, 1)
=> 0.30022

```
dblquad (f, xa, xb, ya, yb)
dblquad (f, xa, xb, ya, yb, tol)
dblquad (f, xa, xb, ya, yb, tol, quadf)
dblquad (f, xa, xb, ya, yb, tol, quadf,...)

Numerically evaluate the double integral of \(f\).
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must have the form \(z=f(x, y)\) where x is a vector and \(y\) is a scalar. It should return a vector of the same length and orientation as \(x\).
\(x a, y a\) and \(x b, y b\) are the lower and upper limits of integration for x and y respectively. The underlying integrator determines whether infinite bounds are accepted.
The optional argument tol defines the absolute tolerance used to integrate each subintegral. The default value is \(1 \mathrm{e}-6\).
The optional argument quadf specifies which underlying integrator function to use. Any choice but quad is available and the default is quadcc.
Additional arguments, are passed directly to \(f\). To use the default value for tol or quadf one may pass ':' or an empty matrix ([]).

See also: [integral2], page 715, [integral3], page 717, [triplequad], page 713, [quad], page 704, [quadv], page 705, [quadl], page 706, [quadgk], page 707, [quadcc], page 708, [trapz], page 711.
triplequad (f, xa, xb, ya, yb, za, zb)
triplequad (f, xa, xb, ya, yb, za, zb, tol)
triplequad (f, xa, xb, ya, yb, za, zb, tol, quadf)
triplequad (f, xa, xb, ya, yb, za, zb, tol, quadf,...)
Numerically evaluate the triple integral of \(f\).
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must have the form \(w=f(x, y, z)\) where either x or y is a vector and the remaining inputs are scalars. It should return a vector of the same length and orientation as \(x\) or \(y\).
\(x a, y a, z a\) and \(x b, y b, z b\) are the lower and upper limits of integration for \(x, y\), and z respectively. The underlying integrator determines whether infinite bounds are accepted.
The optional argument tol defines the absolute tolerance used to integrate each subintegral. The default value is \(1 \mathrm{e}-6\).
The optional argument quadf specifies which underlying integrator function to use. Any choice but quad is available and the default is quadcc.
Additional arguments, are passed directly to \(f\). To use the default value for tol or quadf one may pass ':' or an empty matrix ([]).

See also: [integral3], page 717, [integral2], page 715, [dblquad], page 713, [quad], page 704, [quadv], page 705, [quadl], page 706, [quadgk], page 707, [quadcc], page 708, [trapz], page 711.

The recursive algorithm for quadrature presented above is referred to as "iterated". A separate 2-D integration method is implemented in the function quad2d. This function performs a "tiled" integration by subdividing the integration domain into rectangular regions and performing separate integrations over those domains. The domains are further subdivided in areas requiring refinement to reach the desired numerical accuracy. For certain functions this method can be faster than the 2-D iteration used in the other functions above.
```

q = quad2d (f, xa, xb, ya, yb)
q = quad2d (f, xa, xb, ya, yb, prop, val,...)
[q, err, iter] = quad2d (...)

```

Numerically evaluate the two-dimensional integral of \(f\) using adaptive quadrature over the two-dimensional domain defined by xa, xb, ya, yb using tiled integration. Additionally, ya and \(y b\) may be scalar functions of \(x\), allowing for the integration over non-rectangular domains.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be of the form \(z=f(x, y)\) where \(x\) is a vector and \(y\) is a scalar. It should return a vector of the same length and orientation as \(x\).
Additional optional parameters can be specified using "property", value pairs. Valid properties are:

AbsTol Define the absolute error tolerance for the quadrature. The default value is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).

RelTol Define the relative error tolerance for the quadrature. The default value is \(1 \mathrm{e}-6\) ( \(1 \mathrm{e}-4\) for single).

MaxFunEvals
The maximum number of function calls to the vectorized function \(f\). The default value is 5000 .

Singular Enable/disable transforms to weaken singularities on the edge of the integration domain. The default value is true.

\section*{Vectorized}

Option to disable vectorized integration, forcing Octave to use only scalar inputs when calling the integrand. The default value is false.

\section*{FailurePlot}

If quad2d fails to converge to the desired error tolerance before MaxFunEvals is reached, a plot of the areas that still need refinement is created. The default value is false.

Adaptive quadrature is used to minimize the estimate of error until the following is satisfied:
\[
\text { error } \leq \max (\text { AbsTol, RelTol } \cdot|q|)
\]

The optional output err is an approximate bound on the error in the integral abs ( \(q-I\) ), where \(I\) is the exact value of the integral. The optional output iter is the number of vectorized function calls to the function \(f\) that were used.
Example 1: integrate a rectangular region in \(x\) - \(y\) plane
```

f = @(x,y) 2*ones (size (x));
q = quad2d (f, 0, 1, 0, 1)
\#q=2

```

The result is a volume, which for this constant-value integrand, is just Length * Width * Height.

Example 2 : integrate a triangular region in x - y plane
```

f = @(x,y) 2*ones (size (x));
ymax = @(x) 1 - x;
q = quad2d (f, 0, 1, 0, ymax)
=>q=1

```

The result is a volume, which for this constant-value integrand, is the Triangle Area x Height or \(1 / 2 *\) Base \(*\) Width \(*\) Height.
Programming Notes: If there are singularities within the integration region it is best to split the integral and place the singularities on the boundary.
Known matlab incompatibility: If tolerances are left unspecified, and any integration limits are of type single, then Octave's integral functions automatically reduce the default absolute and relative error tolerances as specified above. If tighter tolerances are desired they must be specified. matlab leaves the tighter tolerances appropriate for double inputs in place regardless of the class of the integration limits.
Reference: L.F. Shampine, matlab program for quadrature in 2D, Applied Mathematics and Computation, pp. 266-274, Vol 1, 2008.
See also: [integral2], page 715, [dblquad], page 713, [integral], page 709, [quad], page 704, [quadgk], page 707, [quadv], page 705, [quadl], page 706, [quadcc], page 708, [trapz], page 711, [integral3], page 717, [triplequad], page 713.

Finally, the functions integral2 and integral3 are provided as general 2-D and 3-D integration functions. They will auto-select between iterated and tiled integration methods and, unlike dblquad and triplequad, will work with non-rectangular integration domains.
```

q = integral2 (f, xa, xb, ya, yb)
q = integral2 (f, xa, xb, ya, yb, prop, val, ...)
[q, err] = integral2 (...)

```

Numerically evaluate the two-dimensional integral of \(f\) using adaptive quadrature over the two-dimensional domain defined by xa, xb, ya, yb (scalars may be finite or
infinite). Additionally, ya and \(y b\) may be scalar functions of \(x\), allowing for integration over non-rectangular domains.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be of the form \(z=f(x, y)\) where x is a vector and \(y\) is a scalar. It should return a vector of the same length and orientation as \(x\).
Additional optional parameters can be specified using "property", value pairs. Valid properties are:

AbsTol Define the absolute error tolerance for the quadrature. The default value is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).
RelTol Define the relative error tolerance for the quadrature. The default value is \(1 \mathrm{e}-6\) ( \(1 \mathrm{e}-4\) for single).

Method Specify the two-dimensional integration method to be used, with valid options being "auto" (default), "tiled", or "iterated". When using "auto", Octave will choose the "tiled" method unless any of the integration limits are infinite.

\section*{Vectorized}

Enable or disable vectorized integration. A value of false forces Octave to use only scalar inputs when calling the integrand, which enables integrands \(f(x, y)\) that have not been vectorized and only accept x and y as scalars to be used. The default value is true.
Adaptive quadrature is used to minimize the estimate of error until the following is satisfied:
\[
\text { error } \leq \max (\text { AbsTol, RelTol } \cdot|q|)
\]
err is an approximate bound on the error in the integral abs ( \(q-I\) ), where \(I\) is the exact value of the integral.
Example 1 : integrate a rectangular region in \(\mathrm{x}-\mathrm{y}\) plane
```

$f=@(x, y) 2 *$ ones (size (x));
$q=i n t e g r a l 2(f, 0,1,0,1)$
$\Rightarrow q=2$

```

The result is a volume, which for this constant-value integrand, is just Length * Width * Height.

Example 2 : integrate a triangular region in x - y plane
```

f = @(x,y) 2*ones (size (x));
ymax = @(x) 1 - x;
q = integral2 (f, 0, 1, 0, ymax)
\#q=1

```

The result is a volume, which for this constant-value integrand, is the Triangle Area x Height or \(1 / 2 *\) Base \(*\) Width \(*\) Height.
Programming Notes: If there are singularities within the integration region it is best to split the integral and place the singularities on the boundary.
Known matlab incompatibility: If tolerances are left unspecified, and any integration limits are of type single, then Octave's integral functions automatically reduce the
default absolute and relative error tolerances as specified above. If tighter tolerances are desired they must be specified. matlab leaves the tighter tolerances appropriate for double inputs in place regardless of the class of the integration limits.
Reference: L.F. Shampine, matlab program for quadrature in 2D, Applied Mathematics and Computation, pp. 266-274, Vol 1, 2008.
See also: [quad2d], page 714, [dblquad], page 713, [integral], page 709, [quad], page 704, [quadgk], page 707, [quadv], page 705, [quadl], page 706, [quadcc], page 708, [trapz], page 711, [integral3], page 717, [triplequad], page 713.
\(q=\) integral3 (f, xa, xb, ya, yb, za, zb)
\(q=\) integral3 (f, xa, xb, ya, yb, za, zb, prop, val, ...)
Numerically evaluate the three-dimensional integral of \(f\) using adaptive quadrature over the three-dimensional domain defined by xa, xb, ya, yb, za, zb (scalars may be finite or infinite). Additionally, ya and yb may be scalar functions of \(x\) and \(z a\), and \(z b\) maybe be scalar functions of \(x\) and \(y\), allowing for integration over non-rectangular domains.
\(f\) is a function handle, inline function, or string containing the name of the function to evaluate. The function \(f\) must be of the form \(z=f(x, y)\) where x is a vector and \(y\) is a scalar. It should return a vector of the same length and orientation as \(x\).
Additional optional parameters can be specified using "property", value pairs. Valid properties are:
AbsTol Define the absolute error tolerance for the quadrature. The default value is \(1 \mathrm{e}-10\) ( \(1 \mathrm{e}-5\) for single).

RelTol Define the relative error tolerance for the quadrature. The default value is \(1 \mathrm{e}-6\) ( \(1 \mathrm{e}-4\) for single).
Method Specify the two-dimensional integration method to be used, with valid options being "auto" (default), "tiled", or "iterated". When using "auto", Octave will choose the "tiled" method unless any of the integration limits are infinite.

\section*{Vectorized}

Enable or disable vectorized integration. A value of false forces Octave to use only scalar inputs when calling the integrand, which enables integrands \(f(x, y)\) that have not been vectorized and only accept x and y as scalars to be used. The default value is true.

Adaptive quadrature is used to minimize the estimate of error until the following is satisfied:
\[
\text { error } \leq \max (\text { AbsTol, RelTol } \cdot|q|)
\]
err is an approximate bound on the error in the integral abs ( \(q-I\) ), where \(I\) is the exact value of the integral.
Example 1 : integrate over a rectangular volume
```

f = @(x,y,z) ones (size (x));
q = integral3 (f, 0, 1, 0, 1, 0, 1)
\#q=1.00000

```

For this constant-value integrand, the result is a volume which is just Length * Width * Height.

Example 2: integrate over a spherical volume
```

f=@(x,y) ones (size (x));
ymax = @(x) sqrt (1 - x.^2);
zmax = @(x,y) sqrt (1 - x.^2 - y.^2);
q = integral3 (f, 0, 1, 0, ymax, 0, zmax)
=>q = 0.52360

```

For this constant-value integrand, the result is a volume which is \(1 / 8\) th of a unit sphere or \(1 / 8 * 4 / 3 *\) pi.
Programming Notes: If there are singularities within the integration region it is best to split the integral and place the singularities on the boundary.
Known matlab incompatibility: If tolerances are left unspecified, and any integration limits are of type single, then Octave's integral functions automatically reduce the default absolute and relative error tolerances as specified above. If tighter tolerances are desired they must be specified. matlab leaves the tighter tolerances appropriate for double inputs in place regardless of the class of the integration limits.
Reference: L.F. Shampine, matlab program for quadrature in 2D, Applied Mathematics and Computation, pp. 266-274, Vol 1, 2008.
See also: [triplequad], page 713, [integral], page 709, [quad], page 704, [quadgk], page 707, [quadv], page 705, [quadl], page 706, [quadcc], page 708, [trapz], page 711, [integral2], page 715, [quad2d], page 714, [dblquad], page 713.

The above integrations can be fairly slow, and that problem increases exponentially with the dimensionality of the integral. Another possible solution for 2-D integration is to use Orthogonal Collocation as described in the previous section (see Section 23.2 [Orthogonal Collocation], page 712). The integral of a function \(f(x, y)\) for \(x\) and \(y\) between 0 and 1 can be approximated using \(n\) points by
\[
\int_{0}^{1} \int_{0}^{1} f(x, y) d x d y \approx \sum_{i=1}^{n} \sum_{j=1}^{n} q_{i} q_{j} f\left(r_{i}, r_{j}\right)
\]
where \(q\) and \(r\) is as returned by colloc ( n ). The generalization to more than two variables is straight forward. The following code computes the studied integral using \(n=8\) points.
```

f = @(x,y) sin (pi*x*y') .* sqrt (x*y');
n = 8;
[t, ~, ~, q] = colloc (n);
I = q'*f(t,t)*q;
=0.30022

```

It should be noted that the number of points determines the quality of the approximation. If the integration needs to be performed between \(a\) and \(b\), instead of 0 and 1 , then a change of variables is needed.

\section*{24 Differential Equations}

Octave has built-in functions for solving ordinary differential equations (ODEs), and differential-algebraic equations (DAEs).

\subsection*{24.1 Ordinary Differential Equations}

The function lsode can be used to solve ODEs of the form
\[
\frac{d x}{d t}=f(x, t)
\]
using Hindmarsh's ODE solver LSODE.
[x, istate, msg] = lsode (fcn, \(\left.x_{-} 0, t\right)\)
[x, istate, msg\(]=1\) sode (fcn, \(\left.x_{-} 0, t_{\text {, }} t_{-} c r i t\right)\)
Ordinary Differential Equation (ODE) solver.
The set of differential equations to solve is
\[
\frac{d x}{d t}=f(x, t)
\]
with
\[
x\left(t_{0}\right)=x_{0}
\]

The solution is returned in the matrix \(x\), with each row corresponding to an element of the vector \(t\). The first element of \(t\) should be \(t_{0}\) and should correspond to the initial state of the system \(x_{-} 0\), so that the first row of the output is \(x_{-} 0\).
The first argument, fcn, is a string, inline, or function handle that names the function \(f\) to call to compute the vector of right hand sides for the set of equations. The function must have the form
```

xdot = f (x, t)

```
in which \(x d o t\) and \(x\) are vectors and \(t\) is a scalar.
If \(f c n\) is a two-element string array or a two-element cell array of strings, inline functions, or function handles, the first element names the function \(f\) described above, and the second element names a function to compute the Jacobian of \(f\). The Jacobian function must have the form
\[
j a c=j(x, t)
\]
in which jac is the matrix of partial derivatives
\[
J=\frac{\partial f_{i}}{\partial x_{j}}=\left[\begin{array}{cccc}
\frac{\partial f_{1}}{\partial x_{1}} & \frac{\partial f_{1}}{\partial x_{2}} & \cdots & \frac{\partial f_{1}}{\partial x_{N}} \\
\frac{\partial f_{2}}{\partial x_{1}} & \frac{\partial f_{2}}{\partial x_{2}} & \cdots & \frac{\partial f_{2}}{\partial x_{N}} \\
\vdots & \vdots & \ddots & \vdots \\
\frac{\partial f_{M}}{\partial x_{1}} & \frac{\partial f_{M}}{\partial x_{2}} & \cdots & \frac{\partial f_{M}}{\partial x_{N}}
\end{array}\right]
\]

The second argument specifies the initial state of the system \(x_{0}\). The third argument is a vector, \(t\), specifying the time values for which a solution is sought.

The fourth argument is optional, and may be used to specify a set of times that the ODE solver should not integrate past. It is useful for avoiding difficulties with singularities and points where there is a discontinuity in the derivative.
After a successful computation, the value of istate will be 2 (consistent with the Fortran version of LSODE).

If the computation is not successful, istate will be something other than 2 and msg will contain additional information.
You can use the function 1sode_options to set optional parameters for 1sode.
See Alan C. Hindmarsh, ODEPACK, A Systematized Collection of ODE Solvers, in Scientific Computing, R. S. Stepleman, editor, (1983) or https://computing.llnl. gov/projects/odepack for more information about the inner workings of lsode.
Example: Solve the Van der Pol equation
```

fvdp = @(y,t) [y(2); (1 - y(1)~2) * y(2) - y(1)];
t = linspace (0, 20, 100);
y = lsode (fvdp, [2; 0], t);

```

See also: [daspk], page 722, [dassl], page 725, [dasrt], page 727.
```

lsode_options ()
val = lsode_options (opt)
lsode_options (opt, val)

```

Query or set options for the function lsode.
When called with no arguments, the names of all available options and their current values are displayed.
Given one argument, return the value of the option opt.
When called with two arguments, lsode_options sets the option opt to value val.
Options include

\section*{"absolute tolerance"}

Absolute tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector.
"relative tolerance"
Relative tolerance parameter. Unlike the absolute tolerance, this parameter may only be a scalar.
The local error test applied at each integration step is
```

abs (local error in x(i)) <= ...
rtol * abs (y(i)) + atol(i)

```
    "integration method"

A string specifying the method of integration to use to solve the ODE system. Valid values are
```

"adams"
"non-stiff"

```

No Jacobian used (even if it is available).
"bdf"
"stiff" Use stiff backward differentiation formula (BDF) method. If a function to compute the Jacobian is not supplied, lsode will compute a finite difference approximation of the Jacobian matrix.
```

"initial step size"

```

The step size to be attempted on the first step (default is determined automatically).
```

"maximum order"

```

Restrict the maximum order of the solution method. If using the Adams method, this option must be between 1 and 12 . Otherwise, it must be between 1 and 5, inclusive.
"maximum step size"
Setting the maximum stepsize will avoid passing over very large regions (default is not specified).
```

"minimum step size"

```

The minimum absolute step size allowed (default is 0 ).
"step limit"
Maximum number of steps allowed (default is 100000).
Here is an example of solving a set of three differential equations using lsode. Given the function
```


## oregonator differential equation

function xdot = f (x, t)
xdot = zeros (3,1);
xdot(1) = 77.27 * (x(2) - x(1)*x(2) + x(1) ...
- 8.375e-06*x(1) ^2);
xdot(2) = (x(3) - x(1)*x(2) - x(2)) / 77.27;
xdot(3) = 0.161*(x(1) - x(3));
endfunction

```
and the initial condition \(x 0=[4 ; 1.1 ; 4]\), the set of equations can be integrated using the command
```

t = linspace (0, 500, 1000);
y = lsode ("f", x0, t);

```

If you try this, you will see that the value of the result changes dramatically between \(t\) \(=0\) and 5 , and again around \(t=305\). A more efficient set of output points might be
```

t = [0, logspace(-1, log10(303), 150), ...
logspace(log10(304), log10(500), 150)];

```

An m-file for the differential equation used above is included with the Octave distribution in the examples directory under the name oregonator.m.

\subsection*{24.2 Differential-Algebraic Equations}

The function daspk can be used to solve DAEs of the form
\[
0=f(\dot{x}, x, t), \quad x(t=0)=x_{0}, \dot{x}(t=0)=\dot{x}_{0}
\]
where \(\dot{x}=\frac{d x}{d t}\) is the derivative of \(x\). The equation is solved using Petzold's DAE solver DASPK.
```

[x, xdot, istate, msg] = daspk (fcn, x_0, xdot_0, t, t_crit)

```

Solve a set of differential-algebraic equations.
daspk solves the set of equations
\[
0=f(x, \dot{x}, t)
\]
with
\[
x\left(t_{0}\right)=x_{0}, \dot{x}\left(t_{0}\right)=\dot{x}_{0}
\]

The solution is returned in the matrices \(x\) and \(x d o t\), with each row in the result matrices corresponding to one of the elements in the vector \(t\). The first element of \(t\) should be \(t_{0}\) and correspond to the initial state of the system x_0 and its derivative xdot_0, so that the first row of the output x is \(\mathrm{x}_{-} 0\) and the first row of the output xdot is xdot_0.
The first argument, fcn, is a string, inline, or function handle that names the function \(f\) to call to compute the vector of residuals for the set of equations. It must have the form
```

res = f (x, xdot, t)

```
in which \(x\), xdot, and res are vectors, and \(t\) is a scalar.
If fcn is a two-element string array or a two-element cell array of strings, inline functions, or function handles, the first element names the function \(f\) described above, and the second element names a function to compute the modified Jacobian
\[
J=\frac{\partial f}{\partial x}+c \frac{\partial f}{\partial \dot{x}}
\]

The modified Jacobian function must have the form
```

jac = j (x, xdot, t, c)

```

The second and third arguments to daspk specify the initial condition of the states and their derivatives, and the fourth argument specifies a vector of output times at which the solution is desired, including the time corresponding to the initial condition.
The set of initial states and derivatives are not strictly required to be consistent. If they are not consistent, you must use the daspk_options function to provide additional information so that daspk can compute a consistent starting point.
The fifth argument is optional, and may be used to specify a set of times that the DAE solver should not integrate past. It is useful for avoiding difficulties with singularities and points where there is a discontinuity in the derivative.

After a successful computation, the value of istate will be greater than zero (consistent with the Fortran version of DASPK).
If the computation is not successful, the value of istate will be less than zero and msg will contain additional information.
You can use the function daspk_options to set optional parameters for daspk.
See also: [dassl], page 725.
```

daspk_options ()

```
val = daspk_options (opt)
daspk_options (opt, val)

Query or set options for the function daspk.
When called with no arguments, the names of all available options and their current values are displayed.

Given one argument, return the value of the option opt.
When called with two arguments, daspk_options sets the option opt to value val.
Options include
"absolute tolerance"
Absolute tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the relative tolerance must also be a vector of the same length.
"relative tolerance"
Relative tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the absolute tolerance must also be a vector of the same length.
The local error test applied at each integration step is
```

abs (local error in x(i))
<= rtol(i) * abs (Y(i)) + atol(i)

```
"compute consistent initial condition"
Denoting the differential variables in the state vector by ' \(Y\) _ \(d\) ' and the algebraic variables by ' Y _a', ddaspk can solve one of two initialization problems:
1. Given Y_d, calculate Y_a and Y'_d
2. Given Y', calculate Y.

In either case, initial values for the given components are input, and initial guesses for the unknown components must also be provided as input. Set this option to 1 to solve the first problem, or 2 to solve the second (the default is 0 , so you must provide a set of initial conditions that are consistent).
If this option is set to a nonzero value, you must also set the "algebraic variables" option to declare which variables in the problem are algebraic.
"use initial condition heuristics"
Set to a nonzero value to use the initial condition heuristics options described below.
"initial condition heuristics"
A vector of the following parameters that can be used to control the initial condition calculation.

MXNIT Maximum number of Newton iterations (default is 5).
MXNJ Maximum number of Jacobian evaluations (default is 6).
MXNH Maximum number of values of the artificial stepsize parameter to be tried if the "compute consistent initial condition" option has been set to 1 (default is 5 ).
Note that the maximum total number of Newton iterations allowed is MXNIT*MXNJ*MXNH if the "compute consistent initial condition" option has been set to 1 and MXNIT*MXNJ if it is set to 2.

LSOFF Set to a nonzero value to disable the linesearch algorithm (default is 0 ).

STPTOL Minimum scaled step in linesearch algorithm (default is \(\left.\operatorname{eps}^{\wedge}(2 / 3)\right)\).

EPINIT Swing factor in the Newton iteration convergence test. The test is applied to the residual vector, premultiplied by the approximate Jacobian. For convergence, the weighted RMS norm of this vector (scaled by the error weights) must be less than EPINIT*EPCON, where EPCON \(=0.33\) is the analogous test constant used in the time steps. The default is EPINIT \(=0.01\).
"print initial condition info"
Set this option to a nonzero value to display detailed information about the initial condition calculation (default is 0 ).
"exclude algebraic variables from error test"
Set to a nonzero value to exclude algebraic variables from the error test. You must also set the "algebraic variables" option to declare which variables in the problem are algebraic (default is 0 ).
"algebraic variables"
A vector of the same length as the state vector. A nonzero element indicates that the corresponding element of the state vector is an algebraic variable (i.e., its derivative does not appear explicitly in the equation set).
This option is required by the "compute consistent initial condition" and "exclude algebraic variables from error test" options.
"enforce inequality constraints"
Set to one of the following values to enforce the inequality constraints specified by the "inequality constraint types" option (default is 0 ).
1. To have constraint checking only in the initial condition calculation.
2. To enforce constraint checking during the integration.
3. To enforce both options 1 and 2.
"inequality constraint types"
A vector of the same length as the state specifying the type of inequality constraint. Each element of the vector corresponds to an element of the state and should be assigned one of the following codes
-2 Less than zero.
-1 Less than or equal to zero.
\(0 \quad\) Not constrained.
1 Greater than or equal to zero.
2 Greater than zero.
This option only has an effect if the "enforce inequality constraints" option is nonzero.
"initial step size"
Differential-algebraic problems may occasionally suffer from severe scaling difficulties on the first step. If you know a great deal about the scaling of your problem, you can help to alleviate this problem by specifying an initial stepsize (default is computed automatically).
"maximum order"
Restrict the maximum order of the solution method. This option must be between 1 and 5 , inclusive (default is 5 ).

\section*{"maximum step size"}

Setting the maximum stepsize will avoid passing over very large regions (default is not specified).

Octave also includes DASSL, an earlier version of DASPK, and DASRT, which can be used to solve DAEs with constraints (stopping conditions).

Solve a set of differential-algebraic equations.
dassl solves the set of equations
\[
0=f(x, \dot{x}, t)
\]
with
\[
x\left(t_{0}\right)=x_{0}, \dot{x}\left(t_{0}\right)=\dot{x}_{0}
\]

The solution is returned in the matrices x and xdot , with each row in the result matrices corresponding to one of the elements in the vector \(t\). The first element of \(t\)
should be \(t_{0}\) and correspond to the initial state of the system x_0 and its derivative xdot_0, so that the first row of the output x is \(\mathrm{x}_{-} 0\) and the first row of the output xdot is xdot_0.

The first argument, fcn, is a string, inline, or function handle that names the function \(f\) to call to compute the vector of residuals for the set of equations. It must have the form
```

res = f (x, xdot, t)

```
in which \(x\), xdot, and res are vectors, and \(t\) is a scalar.
If fcn is a two-element string array or a two-element cell array of strings, inline functions, or function handles, the first element names the function \(f\) described above, and the second element names a function to compute the modified Jacobian
\[
J=\frac{\partial f}{\partial x}+c \frac{\partial f}{\partial \dot{x}}
\]

The modified Jacobian function must have the form
```

jac = j (x, xdot, t, c)

```

The second and third arguments to dassl specify the initial condition of the states and their derivatives, and the fourth argument specifies a vector of output times at which the solution is desired, including the time corresponding to the initial condition.
The set of initial states and derivatives are not strictly required to be consistent. In practice, however, DASSL is not very good at determining a consistent set for you, so it is best if you ensure that the initial values result in the function evaluating to zero. The fifth argument is optional, and may be used to specify a set of times that the DAE solver should not integrate past. It is useful for avoiding difficulties with singularities and points where there is a discontinuity in the derivative.

After a successful computation, the value of istate will be greater than zero (consistent with the Fortran version of DASSL).
If the computation is not successful, the value of istate will be less than zero and msg will contain additional information.
You can use the function dassl_options to set optional parameters for dassl.
See also: [daspk], page 722, [dasrt], page 727, [lsode], page 719.
```

dassl_options ()
val = dassl_options (opt)
dassl_options (opt, val)

```

Query or set options for the function dassl.
When called with no arguments, the names of all available options and their current values are displayed.
Given one argument, return the value of the option opt.
When called with two arguments, dassl_options sets the option opt to value val.

Options include
"absolute tolerance"
Absolute tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the relative tolerance must also be a vector of the same length.
"relative tolerance"
Relative tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the absolute tolerance must also be a vector of the same length.
The local error test applied at each integration step is
```

abs (local error in x(i))
<= rtol(i) * abs (Y(i)) + atol(i)

```
"compute consistent initial condition"
If nonzero, dassl will attempt to compute a consistent set of initial conditions. This is generally not reliable, so it is best to provide a consistent set and leave this option set to zero.
"enforce nonnegativity constraints"
If you know that the solutions to your equations will always be nonnegative, it may help to set this parameter to a nonzero value. However, it is probably best to try leaving this option set to zero first, and only setting it to a nonzero value if that doesn't work very well.
"initial step size"
Differential-algebraic problems may occasionally suffer from severe scaling difficulties on the first step. If you know a great deal about the scaling of your problem, you can help to alleviate this problem by specifying an initial stepsize.
"maximum order"
Restrict the maximum order of the solution method. This option must be between 1 and 5 , inclusive.
"maximum step size"
Setting the maximum stepsize will avoid passing over very large regions (default is not specified).
"step limit"
Maximum number of integration steps to attempt on a single call to the underlying Fortran code.
```

[x, xdot, t_out, istat, msg] = dasrt (fcn, g, $\left.x_{-} 0, x d o t_{-} 0, t\right)$
$\ldots=\operatorname{dasrt}\left(f c n, g, x_{-} 0, x_{d o t} 0, t, t_{-} c r i t\right)$
$\ldots=\operatorname{dasrt}\left(f c n, x_{-} 0, x_{d o t} 0, t\right)$
$\ldots=\operatorname{dasrt}\left(f c n, x_{-} 0, x d o t_{-} 0, t, t_{-} c r i t\right)$

```

Solve a set of differential-algebraic equations.
dasrt solves the set of equations
\[
0=f(x, \dot{x}, t)
\]
with
\[
x\left(t_{0}\right)=x_{0}, \dot{x}\left(t_{0}\right)=\dot{x}_{0}
\]
with functional stopping criteria (root solving).
The solution is returned in the matrices \(x\) and \(x d o t\), with each row in the result matrices corresponding to one of the elements in the vector \(t_{-}\)out. The first element of \(t\) should be \(t_{0}\) and correspond to the initial state of the system x_0 and its derivative xdot_0, so that the first row of the output x is \(\mathrm{x}_{-} 0\) and the first row of the output \(x d o t\) is xdot_0.

The vector \(t\) provides an upper limit on the length of the integration. If the stopping condition is met, the vector \(t_{-}\)out will be shorter than \(t\), and the final element of \(t_{-}\)out will be the point at which the stopping condition was met, and may not correspond to any element of the vector \(t\).
The first argument, fcn, is a string, inline, or function handle that names the function \(f\) to call to compute the vector of residuals for the set of equations. It must have the form
```

res = f (x, xdot, t)

```
in which \(x, x d o t\), and res are vectors, and \(t\) is a scalar.
If fcn is a two-element string array or a two-element cell array of strings, inline functions, or function handles, the first element names the function \(f\) described above, and the second element names a function to compute the modified Jacobian
\[
J=\frac{\partial f}{\partial x}+c \frac{\partial f}{\partial \dot{x}}
\]

The modified Jacobian function must have the form
```

jac = j (x, xdot, t, c)

```

The optional second argument names a function that defines the constraint functions whose roots are desired during the integration. This function must have the form
g_out \(=g(x, t)\)
and return a vector of the constraint function values. If the value of any of the constraint functions changes sign, DASRT will attempt to stop the integration at the point of the sign change.
If the name of the constraint function is omitted, dasrt solves the same problem as daspk or dassl.
Note that because of numerical errors in the constraint functions due to round-off and integration error, DASRT may return false roots, or return the same root at two or more nearly equal values of \(T\). If such false roots are suspected, the user should consider smaller error tolerances or higher precision in the evaluation of the constraint functions.
If a root of some constraint function defines the end of the problem, the input to DASRT should nevertheless allow integration to a point slightly past that root, so that DASRT can locate the root by interpolation.

The third and fourth arguments to dasrt specify the initial condition of the states and their derivatives, and the fourth argument specifies a vector of output times at which the solution is desired, including the time corresponding to the initial condition. The set of initial states and derivatives are not strictly required to be consistent. In practice, however, DASSL is not very good at determining a consistent set for you, so it is best if you ensure that the initial values result in the function evaluating to zero.
The sixth argument is optional, and may be used to specify a set of times that the DAE solver should not integrate past. It is useful for avoiding difficulties with singularities and points where there is a discontinuity in the derivative.
After a successful computation, the value of istate will be greater than zero (consistent with the Fortran version of DASSL).
If the computation is not successful, the value of istate will be less than zero and msg will contain additional information.
You can use the function dasrt_options to set optional parameters for dasrt.
See also: [dasrt_options], page 729, [daspk], page 722, [dasrt], page 727, [lsode], page 719 .
dasrt_options ()
val = dasrt_options (opt)
dasrt_options (opt, val)
Query or set options for the function dasrt.
When called with no arguments, the names of all available options and their current values are displayed.
Given one argument, return the value of the option opt.
When called with two arguments, dasrt_options sets the option opt to value val.
Options include
"absolute tolerance"
Absolute tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the relative tolerance must also be a vector of the same length.
"relative tolerance"
Relative tolerance. May be either vector or scalar. If a vector, it must match the dimension of the state vector, and the absolute tolerance must also be a vector of the same length.
The local error test applied at each integration step is
```

abs (local error in x(i)) <= ...
rtol(i) * abs (Y(i)) + atol(i)

```
    "initial step size"

Differential-algebraic problems may occasionally suffer from severe scaling difficulties on the first step. If you know a great deal about the scaling of your problem, you can help to alleviate this problem by specifying an initial stepsize.
"maximum order"
Restrict the maximum order of the solution method. This option must be between 1 and 5 , inclusive.
"maximum step size"
Setting the maximum stepsize will avoid passing over very large regions.
"step limit"
Maximum number of integration steps to attempt on a single call to the underlying Fortran code.

See K. E. Brenan, et al., Numerical Solution of Initial-Value Problems in DifferentialAlgebraic Equations, North-Holland (1989), DOI: https://doi.org/10.1137/1. 9781611971224 , for more information about the implementation of DASSL.

\subsection*{24.3 Matlab-compatible solvers}

Octave also provides a set of solvers for initial value problems for ordinary differential equations (ODEs) that have a matLab-compatible interface. The options for this class of methods are set using the functions.
- [odeset], page 737,
- [odeget], page 739,

Currently implemented solvers are:
- Runge-Kutta methods
- [ode45], page 730, integrates a system of non-stiff ODEs or index-1 differentialalgebraic equations (DAEs) using the high-order, variable-step Dormand-Prince method. It requires six function evaluations per integration step, but may take larger steps on smooth problems than ode23: potentially offering improved efficiency at smaller tolerances.
- [ode23], page 731, integrates a system of non-stiff ODEs or (or index-1 DAEs). It uses the third-order Bogacki-Shampine method and adapts the local step size in order to satisfy a user-specified tolerance. The solver requires three function evaluations per integration step.
- [ode23s], page 732, integrates a system of stiff ODEs (or index-1 DAEs) using a modified second-order Rosenbrock method.
- Linear multistep methods
- [ode15s], page 733, integrates a system of stiff ODEs (or index-1 DAEs) using a variable step, variable order method based on Backward Difference Formulas (BDF).
- [ode15i], page 735, integrates a system of fully-implicit ODEs (or index-1 DAEs) using the same variable step, variable order method as ode15s. [decic], page 736, can be used to compute consistent initial conditions for ode15i.

Detailed information on the solvers are given in L. F. Shampine and M. W. Reichelt, The MATLAB ODE Suite, SIAM Journal on Scientific Computing, Vol. 18, 1997, pp. 1-22, DOI: https://doi.org/10.1137/S1064827594276424.
[t, y] = ode45 (fun, trange, init)
[t, y] = ode45 (fun, trange, init, ode_opt)
[t, y, te, ye, ie] = ode45 (...)
solution \(=\) ode45 (...)
ode45 (...)
Solve a set of non-stiff Ordinary Differential Equations (non-stiff ODEs) with the well known explicit Dormand-Prince method of order 4.
fun is a function handle, inline function, or string containing the name of the function that defines the ODE: \(y^{\prime}=f(t, y)\). The function must accept two inputs where the first is time \(t\) and the second is a column vector of unknowns \(y\).
trange specifies the time interval over which the ODE will be evaluated. Typically, it is a two-element vector specifying the initial and final times ([tinit, tfinal]). If there are more than two elements then the solution will also be evaluated at these intermediate time instances.

By default, ode45 uses an adaptive timestep with the integrate_adaptive algorithm. The tolerance for the timestep computation may be changed by using the options "RelTol" and "AbsTol".
init contains the initial value for the unknowns. If it is a row vector then the solution \(y\) will be a matrix in which each column is the solution for the corresponding initial value in init.
The optional fourth argument ode_opt specifies non-default options to the ODE solver. It is a structure generated by odeset.
The function typically returns two outputs. Variable \(t\) is a column vector and contains the times where the solution was found. The output \(y\) is a matrix in which each column refers to a different unknown of the problem and each row corresponds to a time in \(t\).
The output can also be returned as a structure solution which has a field \(x\) containing a row vector of times where the solution was evaluated and a field \(y\) containing the solution matrix such that each column corresponds to a time in \(x\). Use fieldnames (solution) to see the other fields and additional information returned.
If no output arguments are requested, and no "OutputFcn" is specified in ode_opt, then the "OutputFcn" is set to odeplot and the results of the solver are plotted immediately.
If using the "Events" option then three additional outputs may be returned. te holds the time when an Event function returned a zero. ye holds the value of the solution at time te. ie contains an index indicating which Event function was triggered in the case of multiple Event functions.
Example: Solve the Van der Pol equation
```

fvdp = @(t,y) [y(2); (1 - y(1)^2) * y(2) - y(1)];
[t,y] = ode45 (fvdp, [0, 20], [2, 0]);

```

See also: [odeset], page 737, [odeget], page 739, [ode23], page 731, [ode15s], page 733.
\([t, y]=\) ode23 (fun, trange, init)
[t, y] = ode23 (fun, trange, init, ode_opt)
[t, y, te, ye, ie] = ode23 (...)
solution \(=\) ode23 (...)
ode23 (...)
Solve a set of non-stiff Ordinary Differential Equations (non-stiff ODEs) with the well known explicit Bogacki-Shampine method of order 3.
fun is a function handle, inline function, or string containing the name of the function that defines the ODE: \(y^{\prime}=f(t, y)\). The function must accept two inputs where the first is time \(t\) and the second is a column vector of unknowns \(y\).
trange specifies the time interval over which the ODE will be evaluated. Typically, it is a two-element vector specifying the initial and final times ([tinit, tfinal]). If there are more than two elements then the solution will also be evaluated at these intermediate time instances.
By default, ode23 uses an adaptive timestep with the integrate_adaptive algorithm. The tolerance for the timestep computation may be changed by using the options "RelTol" and "AbsTol".
init contains the initial value for the unknowns. If it is a row vector then the solution \(y\) will be a matrix in which each column is the solution for the corresponding initial value in init.
The optional fourth argument ode_opt specifies non-default options to the ODE solver. It is a structure generated by odeset.
The function typically returns two outputs. Variable \(t\) is a column vector and contains the times where the solution was found. The output \(y\) is a matrix in which each column refers to a different unknown of the problem and each row corresponds to a time in \(t\).

The output can also be returned as a structure solution which has a field \(x\) containing a row vector of times where the solution was evaluated and a field \(y\) containing the solution matrix such that each column corresponds to a time in \(x\). Use fieldnames (solution) to see the other fields and additional information returned.
If no output arguments are requested, and no "OutputFcn" is specified in ode_opt, then the "OutputFcn" is set to odeplot and the results of the solver are plotted immediately.

If using the "Events" option then three additional outputs may be returned. te holds the time when an Event function returned a zero. ye holds the value of the solution at time te. ie contains an index indicating which Event function was triggered in the case of multiple Event functions.
Example: Solve the Van der Pol equation
```

fvdp = @(t,y) [y(2); (1 - y(1) ^2) * y(2) - y(1)];
[t,y] = ode23 (fvdp, [0, 20], [2, 0]);

```

Reference: For the definition of this method see https://en.wikipedia.org/wiki/ List_of_Runge\%E2\%80\%93Kutta_methods.

See also: [odeset], page 737, [odeget], page 739, [ode45], page 730, [ode15s], page 733.
\([t, y]=\) ode23s (fun, trange, init)
\([t, y]=\) ode23s (fun, trange, init, ode_opt)
\([t, y]=\) ode23s (..., par1, par2, ...)
[t, y, te, ye, ie] = ode23s (...)
solution = ode23s (...)
Solve a set of stiff Ordinary Differential Equations (stiff ODEs) with a Rosenbrock method of order \((2,3)\).
fun is a function handle, inline function, or string containing the name of the function that defines the ODE: \(M y^{\prime}=f(t, y)\). The function must accept two inputs where the first is time \(t\) and the second is a column vector of unknowns \(y . M\) is a constant mass matrix, non-singular and possibly sparse. Set the field "Mass" in odeopts using odeset to specify a mass matrix.
trange specifies the time interval over which the ODE will be evaluated. Typically, it is a two-element vector specifying the initial and final times ([tinit, tfinal]). If there are more than two elements then the solution will also be evaluated at these intermediate time instances using an interpolation procedure of the same order as the one of the solver.
By default, ode23s uses an adaptive timestep with the integrate_adaptive algorithm. The tolerance for the timestep computation may be changed by using the options "RelTol" and "AbsTol".
init contains the initial value for the unknowns. If it is a row vector then the solution \(y\) will be a matrix in which each column is the solution for the corresponding initial value in init.
The optional fourth argument ode_opt specifies non-default options to the ODE solver. It is a structure generated by odeset. ode23s will ignore the following options: "BDF", "InitialSlope", "MassSingular", "MStateDependence", "MvPattern", "MaxOrder", "Non-negative".
The function typically returns two outputs. Variable \(t\) is a column vector and contains the times where the solution was found. The output \(y\) is a matrix in which each column refers to a different unknown of the problem and each row corresponds to a time in \(t\). If trange specifies intermediate time steps, only those will be returned.
The output can also be returned as a structure solution which has a field x containing a row vector of times where the solution was evaluated and a field \(y\) containing the solution matrix such that each column corresponds to a time in x . Use fieldnames (solution) to see the other fields and additional information returned.
If using the "Events" option then three additional outputs may be returned. te holds the time when an Event function returned a zero. ye holds the value of the solution at time te. ie contains an index indicating which Event function was triggered in the case of multiple Event functions.
Example: Solve the stiff Van der Pol equation
```

f = @(t,y) [y(2); 1000*(1 - y(1)^2) * y(2) - y(1)];
opt = odeset ('Mass', [1 0; 0 1], 'MaxStep', 1e-1);
[vt, vy] = ode23s (f, [0 2000], [2 0], opt);

```

See also: [odeset], page 737, [daspk], page 722, [dassl], page 725 .
\([t, y]=\) ode15s (fun, trange, \(y 0\) )
\([t, y]=\) ode15s (fun, trange, \(y 0\), ode_opt)
[t, y, te, ye, ie] = ode15s (...)
solution \(=\) ode15s (...)
ode15s (...)
Solve a set of stiff Ordinary Differential Equations (ODEs) or stiff semi-explicit index 1 Differential Algebraic Equations (DAEs).
ode15s uses a variable step, variable order BDF (Backward Differentiation Formula) method that ranges from order 1 to 5 .
fun is a function handle, inline function, or string containing the name of the function that defines the ODE: \(y^{\prime}=f(t, y)\). The function must accept two inputs where the first is time \(t\) and the second is a column vector of unknowns \(y\).
trange specifies the time interval over which the ODE will be evaluated. Typically, it is a two-element vector specifying the initial and final times ([tinit, tfinal]). If there are more than two elements then the solution will also be evaluated at these intermediate time instances.
init contains the initial value for the unknowns. If it is a row vector then the solution \(y\) will be a matrix in which each column is the solution for the corresponding initial value in init.
The optional fourth argument ode_opt specifies non-default options to the ODE solver. It is a structure generated by odeset.
The function typically returns two outputs. Variable \(t\) is a column vector and contains the times where the solution was found. The output \(y\) is a matrix in which each column refers to a different unknown of the problem and each row corresponds to a time in \(t\).
The output can also be returned as a structure solution which has a field \(x\) containing a row vector of times where the solution was evaluated and a field \(y\) containing the solution matrix such that each column corresponds to a time in \(x\). Use fieldnames (solution) to see the other fields and additional information returned.

If no output arguments are requested, and no "OutputFcn" is specified in ode_opt, then the "OutputFcn" is set to odeplot and the results of the solver are plotted immediately.

If using the "Events" option then three additional outputs may be returned. te holds the time when an Event function returned a zero. ye holds the value of the solution at time te. ie contains an index indicating which Event function was triggered in the case of multiple Event functions.

Example: Solve Robertson's equations:
```

function r = robertson_dae ( }t,y\mathrm{ )
r = [ -0.04*y(1) + 1e4*y(2)*y(3)
0.04*y(1) - 1e4*y(2)*y(3) - 3e7*y(2) ^2
y(1) + y(2) + y(3) - 1 ];
endfunction
opt = odeset ("Mass", [1 0 0; 0 1 0; 0 0 0], "MStateDependence", "none");
[t,y] = ode15s (@robertson_dae, [0, 1e3], [1; 0; 0], opt);

```

See also: [decic], page 736, [odeset], page 737, [odeget], page 739, [ode23], page 731, [ode45], page 730 .
[t, y] = ode15i (fun, trange, y0, yp0)
[t, y] = ode15i (fun, trange, y0, yp0, ode_opt)
[t, y, te, ye, ie] = ode15i (...)
solution = ode15i (...)
ode15i (...)
Solve a set of fully-implicit Ordinary Differential Equations (ODEs) or index 1 Differential Algebraic Equations (DAEs).
ode15i uses a variable step, variable order BDF (Backward Differentiation Formula) method that ranges from order 1 to 5 .
fun is a function handle, inline function, or string containing the name of the function that defines the ODE: \(0=f(t, y, y p)\). The function must accept three inputs where the first is time \(t\), the second is the function value \(y\) (a column vector), and the third is the derivative value yp (a column vector).
trange specifies the time interval over which the ODE will be evaluated. Typically, it is a two-element vector specifying the initial and final times ([tinit, tfinal]). If there are more than two elements then the solution will also be evaluated at these intermediate time instances.
\(y 0\) and yp0 contain the initial values for the unknowns \(y\) and \(y p\). If they are row vectors then the solution \(y\) will be a matrix in which each column is the solution for the corresponding initial value in y0 and yp0.
\(y 0\) and yp0 must be consistent initial conditions, meaning that \(f(t, y 0, y p 0)=0\) is satisfied. The function decic may be used to compute consistent initial conditions given initial guesses.

The optional fifth argument ode_opt specifies non-default options to the ODE solver. It is a structure generated by odeset.

The function typically returns two outputs. Variable \(t\) is a column vector and contains the times where the solution was found. The output \(y\) is a matrix in which each column refers to a different unknown of the problem and each row corresponds to a time in \(t\).

The output can also be returned as a structure solution which has a field x containing a row vector of times where the solution was evaluated and a field \(y\) containing the solution matrix such that each column corresponds to a time in \(x\). Use fieldnames (solution) to see the other fields and additional information returned.

If no output arguments are requested, and no "OutputFcn" is specified in ode_opt, then the "OutputFcn" is set to odeplot and the results of the solver are plotted immediately.

If using the "Events" option then three additional outputs may be returned. te holds the time when an Event function returned a zero. ye holds the value of the solution at time te. ie contains an index indicating which Event function was triggered in the case of multiple Event functions.

Example: Solve Robertson's equations:
```

function r = robertson_dae (t, y, yp)
r = [ -(yp(1) + 0.04*y(1) - 1e4*y(2)*y(3))
-(yp(2) - 0.04*y(1) + 1e4*y(2)*y(3) + 3e7*y(2)^2)
y(1) + y(2) + y(3) - 1];
endfunction
[t,y] = ode15i (@robertson_dae, [0, 1e3], [1; 0; 0], [-1e-4; 1e-4; 0]);

```

See also: [decic], page 736, [odeset], page 737, [odeget], page 739.
[y0_new, yp0_new] = decic (fun, t0, y0, fixed_y0, yp0, fixed_yp0)
[y0_new, yp0_new] = decic (fun, t0, y0, fixed_y0, yp0, fixed_yp0,
options)
[y0_new, yp0_new, resnorm] = decic (...)
Compute consistent implicit ODE initial conditions y0_new and yp0_new given initial guesses y0 and yp0.
A maximum of length (y0) components between fixed_y0 and fixed_yp0 may be chosen as fixed values.
fun is a function handle. The function must accept three inputs where the first is time \(t\), the second is a column vector of unknowns \(y\), and the third is a column vector of unknowns \(y p\).
\(t 0\) is the initial time such that \(f u n\left(t 0, y 0 \_n e w, y p 0_{n} n e w\right)=0\), specified as a scalar.
\(y 0\) is a vector used as the initial guess for \(y\).
fixed_y0 is a vector which specifies the components of y0 to hold fixed. Choose a maximum of length ( \(y 0\) ) components between fixed_y0 and fixed_yp0 as fixed values. Set fixed_y0(i) component to 1 if you want to fix the value of y0(i). Set fixed_y0(i) component to 0 if you want to allow the value of \(y 0(\mathrm{i})\) to change.
\(y p 0\) is a vector used as the initial guess for \(y p\).
fixed_yp0 is a vector which specifies the components of yp0 to hold fixed. Choose a maximum of length (yp0) components between fixed_y0 and fixed_yp0 as fixed values. Set fixed_yp0(i) component to 1 if you want to fix the value of yp0(i). Set fixed_yp \(0(\mathrm{i})\) component to 0 if you want to allow the value of \(y p 0(\mathrm{i})\) to change.
The optional seventh argument options is a structure array. Use odeset to generate this structure. The relevant options are RelTol and AbsTol which specify the error thresholds used to compute the initial conditions.
The function typically returns two outputs. Variable y0_new is a column vector and contains the consistent initial value of \(y\). The output yp0_new is a column vector and contains the consistent initial value of \(y p\).
The optional third output resnorm is the norm of the vector of residuals. If resnorm is small, decic has successfully computed the initial conditions. If the value of resnorm is large, use RelTol and AbsTol to adjust it.
Example: Compute initial conditions for Robertson's equations:
```

function r = robertson_dae (t, y, yp)
r = [ - (yp(1) + 0.04*y(1) - 1e4*y(2)*y(3))
-(yp(2) - 0.04*y(1) + 1e4*y(2)*y(3) + 3e7*y(2)^2)
y(1) + y(2) + y(3) - 1 ];
endfunction
[y0_new,ypO_new] = decic (@robertson_dae, 0, [1; 0; 0], [1; 1; 0],

```
[-1e-4; 1; 0], [0; 0; 0]);
See also: [ode15i], page 735, [odeset], page 737.
```

odestruct = odeset ()

```
odestruct \(=\) odeset ("field1", value1, "field2", value2, ...)
odestruct = odeset (oldstruct, "field1", value1, "field2", value2, ...)
odestruct \(=\) odeset (oldstruct, newstruct)
odeset ()

Create or modify an ODE options structure.
When called with no input argument and one output argument, return a new ODE options structure that contains all possible fields initialized to their default values. If no output argument is requested, display a list of the common ODE solver options along with their default value.
If called with name-value input argument pairs "field1", " value1", "field2", " value2", ... return a new ODE options structure with all the most common option fields initialized, and set the values of the fields "field1", "field2", . . . to the values value1, value2, ....
If called with an input structure oldstruct then overwrite the values of the options "field1", "field2", ... with new values value1, value2, ... and return the modified structure.
When called with two input ODE options structures oldstruct and newstruct overwrite all values from the structure oldstruct with new values from the structure newstruct. Empty values in newstruct will not overwrite values in oldstruct.
The most commonly used ODE options, which are always assigned a value by odeset, are the following:
AbsTol: positive scalar | vector, def. 1e-6
Absolute error tolerance.
BDF: \{"off"\} | "on"
Use BDF formulas in implicit multistep methods. Note: This option is not yet implemented.
Events: function_handle
Event function. An event function must have the form [value, isterminal, direction] = my_events_f ( \(t, y\) )

InitialSlope: vector
Consistent initial slope vector for DAE solvers.
InitialStep: positive scalar
Initial time step size.
Jacobian: matrix | function_handle
Jacobian matrix, specified as a constant matrix or a function of time and state.
JConstant: \{"off"\} | "on"
Specify whether the Jacobian is a constant matrix or depends on the state.

JPattern: sparse matrix
If the Jacobian matrix is sparse and non-constant but maintains a constant sparsity pattern, specify the sparsity pattern.

Mass: matrix | function_handle
Mass matrix, specified as a constant matrix or a function of time and state.

MassSingular: \{"maybe"\} | "yes" | "on"
Specify whether the mass matrix is singular.
MaxOrder: \(\{5\}|4| 3|2| 1\)
Maximum order of formula.
MaxStep: positive scalar
Maximum time step value.
MStateDependence: \{"weak"\} | "none" | "strong"
Specify whether the mass matrix depends on the state or only on time.
MvPattern: sparse matrix
If the mass matrix is sparse and non-constant but maintains a constant sparsity pattern, specify the sparsity pattern. Note: This option is not yet implemented.

NonNegative: scalar | vector
Specify elements of the state vector that are expected to remain nonnegative during the simulation.

NormControl: \{"off"\} | "on"
Control error relative to the 2 -norm of the solution, rather than its absolute value.

OutputFcn: function_handle
Function to monitor the state during the simulation. For the form of the function to use see [odeplot], page 739.

OutputSel: scalar | vector
Indices of elements of the state vector to be passed to the output monitoring function.

Refine: positive scalar
Specify whether output should be returned only at the end of each time step or also at intermediate time instances. The value should be a scalar indicating the number of equally spaced time points to use within each timestep at which to return output. Note: This option is not yet implemented.

RelTol: positive scalar
Relative error tolerance.
Stats: \{"off"\} | "on"
Print solver statistics after simulation.

Vectorized: \{"off"\} | "on"
Specify whether odefun can be passed multiple values of the state at once.

Field names that are not in the above list are also accepted and added to the result structure.
See also: [odeget], page 739 .
val = odeget (ode_opt, field)
val = odeget (ode_opt, field, default)
Query the value of the property field in the ODE options structure ode_opt.
If called with two input arguments and the first input argument ode_opt is an ODE option structure and the second input argument field is a string specifying an option name, then return the option value val corresponding to field from ode_opt.
If called with an optional third input argument, and field is not set in the structure ode_opt, then return the default value default instead.

See also: [odeset], page 737.
```

stop_solve = odeplot (t, y, flag)

```

Open a new figure window and plot the solution of an ode problem at each time step during the integration.
The types and values of the input parameters \(t\) and \(y\) depend on the input flag that is of type string. Valid values of flag are:
"init" The input \(t\) must be a column vector of length 2 with the first and last time step ([tfirst tlast]. The input y contains the initial conditions for the ode problem (y0).
"" The input \(t\) must be a scalar double specifying the time for which the solution in input \(y\) was calculated.
"done" The inputs should be empty, but are ignored if they are present.
odeplot always returns false, i.e., don't stop the ode solver.
Example: solve an anonymous implementation of the "Van der Pol" equation and display the results while solving.
```

fvdp = @(t,y) [y(2); (1 - y(1)^2) * y(2) - y(1)];
opt = odeset ("OutputFcn", @odeplot, "RelTol", 1e-6);
sol = ode45 (fvdp, [0 20], [2 0], opt);

```

Background Information: This function is called by an ode solver function if it was specified in the "OutputFcn" property of an options structure created with odeset. The ode solver will initially call the function with the syntax odeplot ([tfirst, tlast], y0, "init"). The function initializes internal variables, creates a new figure window, and sets the x limits of the plot. Subsequently, at each time step during the integration the ode solver calls odeplot ( \(t, y\), []). At the end of the solution the ode solver calls odeplot ([] , [], "done") so that odeplot can perform any clean-up actions required.
See also: [odeset], page 737, [odeget], page 739, [ode23], page 731, [ode45], page 730.

\section*{25 Optimization}

Octave comes with support for solving various kinds of optimization problems. Specifically Octave can solve problems in Linear Programming, Quadratic Programming, Nonlinear Programming, and Linear Least Squares Minimization.

\subsection*{25.1 Linear Programming}

Octave can solve Linear Programming problems using the glpk function. That is, Octave can solve
\[
\min _{x} c^{T} x
\]
subject to the linear constraints \(A x=b\) where \(x \geq 0\).
The glpk function also supports variations of this problem.
[xopt, fmin, errnum, extra] = glpk (c, A, b, lb, ub, ctype, vartype, sense, param)
Solve a linear program using the GNU GLPK library.
Given three arguments, glpk solves the following standard LP:
\[
\min _{x} C^{T} x
\]
subject to
\[
A x=b \quad x \geq 0
\]
but may also solve problems of the form
\[
\left[\min _{x} \mid \max _{x}\right] C^{T} x
\]
subject to
\[
A x[=|\leq| \geq] b \quad L B \leq x \leq U B
\]

Input arguments:
c A column array containing the objective function coefficients.
A A matrix containing the constraints coefficients.
\(b \quad\) A column array containing the right-hand side value for each constraint in the constraint matrix.
\(l b \quad\) An array containing the lower bound on each of the variables. If \(l b\) is not supplied, the default lower bound for the variables is zero.
\(u b \quad\) An array containing the upper bound on each of the variables. If \(u b\) is not supplied, the default upper bound is assumed to be infinite.
ctype An array of characters containing the sense of each constraint in the constraint matrix. Each element of the array may be one of the following values
"F" A free (unbounded) constraint (the constraint is ignored).
\[
\begin{array}{ll}
\text { "U" } & \begin{array}{l}
\text { An inequality constraint with an upper bound }(\mathrm{A}(\mathrm{i},::) * \mathrm{x}<= \\
\mathrm{b}(\mathrm{i})) .
\end{array} \\
\text { "S" } & \text { An equality constraint }(\mathrm{A}(\mathrm{i},::) * \mathrm{x}=\mathrm{b}(\mathrm{i})) . \\
\text { "L" } & \text { An inequality with a lower bound }(\mathrm{A}(\mathrm{i},::) * \mathrm{x}>=\mathrm{b}(\mathrm{i})) . \\
\text { "D" } & \begin{array}{l}
\text { An inequality constraint with both upper and lower bounds } \\
(\mathrm{A}(\mathrm{i},:) * \mathrm{x}>=-\mathrm{b}(\mathrm{i})) \text { and }(\mathrm{A}(\mathrm{i},: \mathrm{O}) * \mathrm{x}<=\mathrm{b}(\mathrm{i})) .
\end{array}
\end{array}
\]
vartype A column array containing the types of the variables.
"C" A continuous variable.
"I" An integer variable.
sense If sense is 1 , the problem is a minimization. If sense is -1 , the problem is a maximization. The default value is 1 .
param A structure containing the following parameters used to define the behavior of solver. Missing elements in the structure take on default values, so you only need to set the elements that you wish to change from the default.

Integer parameters:
msglev (default: 1)
Level of messages output by solver routines:
0 (GLP_MSG_OFF)
No output.
1 (GLP_MSG_ERR)
Error and warning messages only.
2 (GLP_MSG_ON)
Normal output.
3 (GLP_MSG_ALL)
Full output (includes informational messages).
scale (default: 16)
Scaling option. The values can be combined with the bitwise OR operator and may be the following:

1 (GLP_SF_GM)
Geometric mean scaling.
16 (GLP_SF_EQ)
Equilibration scaling.
32 (GLP_SF_2N)
Round scale factors to power of two.
64 (GLP_SF_SKIP)
Skip if problem is well scaled.

Alternatively, a value of 128 (GLP_SF_AUTO) may be also specified, in which case the routine chooses the scaling options automatically.
```

dual (default: 1)

```

Simplex method option:
1 (GLP_PRIMAL)
Use two-phase primal simplex.
2 (GLP_DUALP)
Use two-phase dual simplex, and if it fails, switch to the primal simplex.

3 (GLP_DUAL)
Use two-phase dual simplex.
price (default: 34)
Pricing option (for both primal and dual simplex):
17 (GLP_PT_STD)
Textbook pricing.
34 (GLP_PT_PSE)
Steepest edge pricing.
itlim (default: intmax)
Simplex iterations limit. It is decreased by one each time when one simplex iteration has been performed, and reaching zero value signals the solver to stop the search.
outfrq (default: 200)
Output frequency, in iterations. This parameter specifies how frequently the solver sends information about the solution to the standard output.
branch (default: 4)
Branching technique option (for MIP only):
1 (GLP_BR_FFV)
First fractional variable.
2 (GLP_BR_LFV)
Last fractional variable.
3 (GLP_BR_MFV)
Most fractional variable.
4 (GLP_BR_DTH)
Heuristic by Driebeck and Tomlin.
5 (GLP_BR_PCH)
Hybrid pseudocost heuristic.
```

btrack (default: 4)
Backtracking technique option (for MIP only):
1 (GLP_BT_DFS)
Depth first search.
2 (GLP_BT_BFS)
Breadth first search.
3 (GLP_BT_BLB)
Best local bound.
4(GLP_BT_BPH)
Best projection heuristic.
presol (default: 1)

```
    If this flag is set, the simplex solver uses the built-in LP
    presolver. Otherwise the LP presolver is not used.
lpsolver (default: 1)
    Select which solver to use. If the problem is a MIP problem
    this flag will be ignored.
    1 Revised simplex method.
    2 Interior point method.
rtest (default: 34)
    Ratio test technique:
    17 (GLP_RT_STD)
            Standard ("textbook").
    34 (GLP_RT_HAR)
            Harris' two-pass ratio test.
tmlim (default: intmax)
    Searching time limit, in milliseconds.
outdly (default: 0)

Output delay, in seconds. This parameter specifies how long the solver should delay sending information about the solution to the standard output.
save (default: 0)
If this parameter is nonzero, save a copy of the problem in CPLEX LP format to the file "outpb.lp". There is currently no way to change the name of the output file.

Real parameters:
tolbnd (default: 1e-7)
Relative tolerance used to check if the current basic solution is primal feasible. It is not recommended that you change this parameter unless you have a detailed understanding of its purpose.
toldj (default: 1e-7)
Absolute tolerance used to check if the current basic solution is dual feasible. It is not recommended that you change this parameter unless you have a detailed understanding of its purpose.
tolpiv (default: 1e-10)
Relative tolerance used to choose eligible pivotal elements of the simplex table. It is not recommended that you change this parameter unless you have a detailed understanding of its purpose.
objll (default: -DBL_MAX)
Lower limit of the objective function. If the objective function reaches this limit and continues decreasing, the solver stops the search. This parameter is used in the dual simplex method only.
objul (default: +DBL_MAX)
Upper limit of the objective function. If the objective function reaches this limit and continues increasing, the solver stops the search. This parameter is used in the dual simplex only.
tolint (default: 1e-5)
Relative tolerance used to check if the current basic solution is integer feasible. It is not recommended that you change this parameter unless you have a detailed understanding of its purpose.
tolobj (default: 1e-7)
Relative tolerance used to check if the value of the objective function is not better than in the best known integer feasible solution. It is not recommended that you change this parameter unless you have a detailed understanding of its purpose.
Output values:
xopt The optimizer (the value of the decision variables at the optimum).
fopt The optimum value of the objective function.
errnum Error code.
0 No error.
1 (GLP_EBADB)
Invalid basis.
2 (GLP_ESING)
Singular matrix.
3 (GLP_ECOND)
Ill-conditioned matrix.
```

        4(GLP_EBOUND)
            Invalid bounds.
    5 (GLP_EFAIL)
Solver failed.
6(GLP_EOBJLL)
Objective function lower limit reached.
7(GLP_EOBJUL)
Objective function upper limit reached.
8(GLP_EITLIM)
Iterations limit exhausted.
9(GLP_ETMLIM)
Time limit exhausted.
10 (GLP_ENOPFS)
No primal feasible solution.
11 (GLP_ENODFS)
No dual feasible solution.
12(GLP_EROOT)
Root LP optimum not provided.
13 (GLP_ESTOP)
Search terminated by application.
14(GLP_EMIPGAP)
Relative MIP gap tolerance reached.
15 (GLP_ENOFEAS)
No primal/dual feasible solution.
16 (GLP_ENOCVG)
No convergence.
17 (GLP_EINSTAB)
Numerical instability.
18 (GLP_EDATA)
Invalid data.
19 (GLP_ERANGE)
Result out of range.
extra A data structure containing the following fields:
lambda Dual variables.
redcosts Reduced Costs.
time Time (in seconds) used for solving LP/MIP problem.
status Status of the optimization.
1 (GLP_UNDEF)
Solution status is undefined.

```

2 (GLP_FEAS)
Solution is feasible.
3 (GLP_INFEAS)
Solution is infeasible.
4 (GLP_NOFEAS)
Problem has no feasible solution.
5 (GLP_OPT)
Solution is optimal.
6 (GLP_UNBND)
Problem has no unbounded solution.
Example:
```

c = [10, 6, 4]';
A = [ 1, 1, 1;
10, 4, 5;
2, 2, 6];
b = [100, 600, 300]';
lb = [0, 0, 0]';
ub = [];
ctype = "UUU";
vartype = "CCC";
s = -1;
param.msglev = 1;
param.itlim = 100;
[xmin, fmin, status, extra] = ...
glpk (c, A, b, lb, ub, ctype, vartype, s, param);

```

\subsection*{25.2 Quadratic Programming}

Octave can also solve Quadratic Programming problems, this is
\[
\min _{x} \frac{1}{2} x^{T} H x+x^{T} q
\]
subject to
\[
A x=b \quad l b \leq x \leq u b \quad A_{l b} \leq A_{i n} x \leq A_{u b}
\]
[x, obj, info, lambda] = qp (x0, H)
[x, obj, info, lambda] = qp (x0, H, q)
[x, obj, info, lambda] = qp (x0, H, q, A, b)
[x, obj, info, lambda] = qp (x0, H, q, A, b, lb, ub)
[x, obj, info, lambda] = qp (x0, H, q, A, b, lb, ub, \(\left.A_{-} l b, A_{-} i n, A_{-} u b\right)\)
[x, obj, info, lambda] = qp (..., options)
Solve a quadratic program (QP).

Solve the quadratic program defined by
\[
\min _{x} \frac{1}{2} x^{T} H x+x^{T} q
\]
subject to
\[
A x=b \quad l b \leq x \leq u b \quad A_{l b} \leq A_{i n} x \leq A_{u b}
\]
using a null-space active-set method.
Any bound ( \(\left.A, b, l b, u b, A_{-} i n, A_{-} l b, A_{-} u b\right)\) may be set to the empty matrix ([]) if not present. The constraints \(A\) and \(A_{-}\)in are matrices with each row representing a single constraint. The other bounds are scalars or vectors depending on the number of constraints. The algorithm is faster if the initial guess is feasible.
options is a structure specifying additional parameters which control the algorithm. Currently, qp recognizes these options: "MaxIter", "TolX".
"MaxIter" proscribes the maximum number of algorithm iterations before optimization is halted. The default value is 200 . The value must be a positive integer.
"TolX" specifies the termination tolerance for the unknown variables \(x\). The default is sqrt (eps) or approximately \(1 \mathrm{e}-8\).
On return, \(x\) is the location of the minimum and fval contains the value of the objective function at \(x\).
info Structure containing run-time information about the algorithm. The following fields are defined:
```

solveiter
The number of iterations required to find the solution.
info An integer indicating the status of the solution.
$0 \quad$ The problem is feasible and convex. Global solution found.
1 The problem is not convex. Local solution found.
2 The problem is not convex and unbounded.
3 Maximum number of iterations reached.
6 The problem is infeasible.

```

See also: [sqp], page 749.
\(x=\) pqpnonneg \((c, d)\)
\(x=\) pqpnonneg \((c, d, x 0)\)
\(x=\) pqpnonneg ( \(c, d, x 0\), options)
[x, minval] = pqpnonneg (...)
[x, minval, exitflag] = pqpnonneg (...)
[x, minval, exitflag, output] = pqpnonneg (...)
[x, minval, exitflag, output, lambda] = pqpnonneg (...)
Minimize ( \(1 / 2 * x^{\prime} * c * x+d^{\prime} * x\) ) subject to \(x>=0\).
\(c\) and \(d\) must be real matrices, and \(c\) must be symmetric and positive definite.
x 0 is an optional initial guess for the solution x .
options is an options structure to change the behavior of the algorithm (see [optimset], page 753). pqpnonneg recognizes one option: "MaxIter".
Outputs:
\(x \quad\) The solution matrix
minval The minimum attained model value, \(1 / 2 * x m i n ' * c * x m i n+d^{\prime} * x m i n\)
exitflag An indicator of convergence. 0 indicates that the iteration count was exceeded, and therefore convergence was not reached; >0 indicates that the algorithm converged. (The algorithm is stable and will converge given enough iterations.)
output A structure with two fields:
- "algorithm": The algorithm used ("nnls")
- "iterations": The number of iterations taken.
lambda Lagrange multipliers. If these are non-zero, the corresponding \(x\) values should be zero, indicating the solution is pressed up against a coordinate plane. The magnitude indicates how much the residual would improve if the \(x>=0\) constraints were relaxed in that direction.

See also: [lsqnonneg], page 752, [qp], page 747, [optimset], page 753.

\subsection*{25.3 Nonlinear Programming}

Octave can also perform general nonlinear minimization using a successive quadratic programming solver.
[x, obj, info, iter, nf, lambda] = sqp (x0, phi)
[...] = sqp (x0, phi, g)
[...] = sqp (x0, phi, g, h)
[...] \(=\operatorname{sqp}(x 0, p h i, g, h, l b, u b)\)
[...] = sqp (x0, phi, g, h, lb, ub, maxiter)
[...] = sqp (x0, phi, g, h, lb, ub, maxiter, tolerance)
Minimize an objective function using sequential quadratic programming (SQP).
Solve the nonlinear program
\[
\min _{x} \phi(x)
\]
subject to
\[
g(x)=0 \quad h(x) \geq 0 \quad l b \leq x \leq u b
\]
using a sequential quadratic programming method.
The first argument is the initial guess for the vector \(x 0\).
The second argument is a function handle pointing to the objective function phi. The objective function must accept one vector argument and return a scalar.
The second argument may also be a 2 - or 3 -element cell array of function handles. The first element should point to the objective function, the second should point to a function that computes the gradient of the objective function, and the third
should point to a function that computes the Hessian of the objective function. If the gradient function is not supplied, the gradient is computed by finite differences. If the Hessian function is not supplied, a BFGS update formula is used to approximate the Hessian.
When supplied, the gradient function phi\{2\} must accept one vector argument and return a vector. When supplied, the Hessian function phi\{3\} must accept one vector argument and return a matrix.
The third and fourth arguments \(g\) and \(h\) are function handles pointing to functions that compute the equality constraints and the inequality constraints, respectively. If the problem does not have equality (or inequality) constraints, then use an empty matrix ([]) for \(g\) (or \(h\) ). When supplied, these equality and inequality constraint functions must accept one vector argument and return a vector.
The third and fourth arguments may also be 2 -element cell arrays of function handles. The first element should point to the constraint function and the second should point to a function that computes the gradient of the constraint function:
\[
\left(\frac{\partial f(x)}{\partial x_{1}}, \frac{\partial f(x)}{\partial x_{2}}, \ldots, \frac{\partial f(x)}{\partial x_{N}}\right)^{T}
\]

The fifth and sixth arguments, \(l b\) and \(u b\), contain lower and upper bounds on \(x\). These must be consistent with the equality and inequality constraints \(g\) and \(h\). If the arguments are vectors then \(x(i)\) is bound by \(l b(i)\) and \(u b(i)\). A bound can also be a scalar in which case all elements of \(x\) will share the same bound. If only one bound ( \(\mathrm{lb}, \mathrm{ub}\) ) is specified then the other will default to (-realmax, +realmax).
The seventh argument maxiter specifies the maximum number of iterations. The default value is 100 .
The eighth argument tolerance specifies the tolerance for the stopping criteria. The default value is sqrt (eps).
The value returned in info may be one of the following:
101 The algorithm terminated normally. All constraints meet the specified tolerance.

102 The BFGS update failed.
103
The maximum number of iterations was reached.
104 The stepsize has become too small, i.e., \(\Delta x\), is less than tol \(*\) norm ( x ).
An example of calling sqp:
```

function r = g (x)
r = [ sumsq(x)-10;
x (2)*x(3)-5*x (4)*x (5);
x(1)^ }3+x(2)^3+1]
endfunction
function obj = phi (x)
obj = exp (prod (x)) - 0.5*(x(1)^3+x(2)^3+1)^2;

```
```

endfunction
x0 = [-1.8; 1.7; 1.9; -0.8; -0.8];
[x, obj, info, iter, nf, lambda] = sqp (x0, @phi, @g, [])
x =
-1.71714
1.59571
1.82725
-0.76364
-0.76364
obj = 0.053950
info = 101
iter = 8
nf = 10
lambda =
-0.0401627
0.0379578
-0.0052227

```

See also: [qp], page 747 .

\subsection*{25.4 Linear Least Squares}

Octave also supports linear least squares minimization. That is, Octave can find the parameter \(b\) such that the model \(y=x b\) fits data \((x, y)\) as well as possible, assuming zero-mean Gaussian noise. If the noise is assumed to be isotropic the problem can be solved using the ' \(\backslash\) ' or '/' operators, or the ols function. In the general case where the noise is assumed to be anisotropic the gls is needed.
[beta, sigma, r] = ols ( \(\mathrm{y}, \mathrm{x}\) )
Ordinary least squares (OLS) estimation.
OLS applies to the multivariate model \(y=x b+e\) where \(y\) is a \(t \times p\) matrix, \(x\) is a \(t \times k\) matrix, \(b\) is a \(k \times p\) matrix, and \(e\) is a \(t \times p\) matrix.
Each row of \(y\) is a \(p\)-variate observation in which each column represents a variable. Likewise, the rows of \(x\) represent \(k\)-variate observations or possibly designed values. Furthermore, the collection of observations \(x\) must be of adequate rank, \(k\), otherwise \(b\) cannot be uniquely estimated.
The observation errors, \(e\), are assumed to originate from an underlying \(p\)-variate distribution with zero mean and \(p\)-by- \(p\) covariance matrix \(S\), both constant conditioned on x. Furthermore, the matrix \(S\) is constant with respect to each observation such that \(\bar{e}=0\) and \(\operatorname{cov}(\operatorname{vec}(e))=\operatorname{kron}(s, I)\). (For cases that don't meet this criteria, such as autocorrelated errors, see generalized least squares, gls, for more efficient estimations.)

The return values beta, sigma, and \(r\) are defined as follows.
beta The OLS estimator for matrix \(b\). beta is calculated directly via \(\left(x^{T} x\right)^{-1} x^{T} y\) if the matrix \(x^{T} x\) is of full rank. Otherwise, beta \(=\) pinv \((x) * y\) where pinv ( \(x\) ) denotes the pseudoinverse of \(x\).
sigma The OLS estimator for the matrix \(s\),
\[
\text { sigma }=(y-x * \text { beta })^{\prime} *(y-x * \text { beta }) /(t-\operatorname{rank}(x))
\]
\(r \quad\) The matrix of OLS residuals, \(r=y-x *\) beta.
See also: [gls], page 752, [pinv], page 593.
[beta, v, r] = gls ( \(\mathrm{y}, \mathrm{x}, \mathrm{o}\) )
Generalized least squares (GLS) model.
Perform a generalized least squares estimation for the multivariate model \(y=x b+e\) where \(y\) is a \(t \times p\) matrix, x is a \(t \times k\) matrix, \(b\) is a \(k \times p\) matrix and \(e\) is a \(t \times p\) matrix.
Each row of \(y\) is a \(p\)-variate observation in which each column represents a variable. Likewise, the rows of x represent \(k\)-variate observations or possibly designed values. Furthermore, the collection of observations x must be of adequate rank, \(k\), otherwise \(b\) cannot be uniquely estimated.
The observation errors, \(e\), are assumed to originate from an underlying \(p\)-variate distribution with zero mean but possibly heteroscedastic observations. That is, in general, \(\bar{e}=0\) and \(\operatorname{cov}(\operatorname{vec}(e))=s^{2} o\) in which \(s\) is a scalar and \(o\) is a \(t p \times t p\) matrix. The return values beta, \(v\), and \(r\) are defined as follows.
beta The GLS estimator for matrix \(b\).
\(v \quad\) The GLS estimator for scalar \(s^{2}\).
\(r \quad\) The matrix of GLS residuals, \(r=y-x *\) beta.
See also: [ols], page 751.
```

x = lsqnonneg (c,d)

```
\(x=1\) sqnonneg ( \(c, d, x 0\) )
\(x=1\) sqnonneg ( \(c, d, x 0\), options)
[x, resnorm] = lsqnonneg (...)
[x, resnorm, residual] = lsqnonneg (...)
[x, resnorm, residual, exitflag] = lsqnonneg (...)
[x, resnorm, residual, exitflag, output] = lsqnonneg (...)
[x, resnorm, residual, exitflag, output, lambda] = lsqnonneg (...) Minimize norm ( \(c * x-d\) ) subject to \(x>=0\).
\(c\) and \(d\) must be real matrices.
x 0 is an optional initial guess for the solution x .
options is an options structure to change the behavior of the algorithm (see [optimset], page 753). 1sqnonneg recognizes these options: "MaxIter", "TolX".
Outputs:
resnorm The squared 2-norm of the residual: norm ( \(\left.c * x^{*}-d\right)^{\wedge} 2\)
residual The residual: \(d-c * x\)
exitflag An indicator of convergence. 0 indicates that the iteration count was exceeded, and therefore convergence was not reached; >0 indicates that the algorithm converged. (The algorithm is stable and will converge given enough iterations.)
output A structure with two fields:
- "algorithm": The algorithm used ("nnls")
- "iterations": The number of iterations taken.
lambda Lagrange multipliers. If these are non-zero, the corresponding \(x\) values should be zero, indicating the solution is pressed up against a coordinate plane. The magnitude indicates how much the residual would improve if the \(\mathrm{x}>=0\) constraints were relaxed in that direction.

See also: [pqpnonneg], page 748, [lscov], page 753, [optimset], page 753.
```

x = lscov (A, b)
x = lscov (A, b,V)
x = lscov (A, b, V, alg)
[x, stdx, mse, S] = lscov (...)

```

Compute a generalized linear least squares fit.
Estimate \(x\) under the model \(b=A x+w\), where the noise \(w\) is assumed to follow a normal distribution with covariance matrix \(\sigma^{2} V\).
If the size of the coefficient matrix \(A\) is n-by-p, the size of the vector/array of constant terms \(b\) must be n -by-k.
The optional input argument \(V\) may be an n-element vector of positive weights (inverse variances), or an n-by-n symmetric positive semi-definite matrix representing the covariance of \(b\). If \(V\) is not supplied, the ordinary least squares solution is returned.
The alg input argument, a guidance on solution method to use, is currently ignored.
Besides the least-squares estimate matrix x ( \(\mathrm{p}-\mathrm{by}-\mathrm{k}\) ), the function also returns stdx (p-by-k), the error standard deviation of estimated \(x\); mse ( \(k\)-by-1), the estimated data error covariance scale factors \(\left(\sigma^{2}\right)\); and \(S\) (p-by-p, or p-by-p-by-k if k > 1 ), the error covariance of \(x\).
Reference: Golub and Van Loan (1996), Matrix Computations (3rd Ed.), Johns Hopkins, Section 5.6.3
See also: [ols], page 751, [gls], page 752, [lsqnonneg], page 752.
optimset ()
options = optimset ()
options = optimset (par, val, ...)
options = optimset (old, par, val, ...)
options = optimset (old, new)
Create options structure for optimization functions.
When called without any input or output arguments, optimset prints a list of all valid optimization parameters.

When called with one output and no inputs, return an options structure with all valid option parameters initialized to [].
When called with a list of parameter/value pairs, return an options structure with only the named parameters initialized.
When the first input is an existing options structure old, the values are updated from either the par/val list or from the options structure new.
Valid parameters are:
AutoScaling
ComplexEqn
Display Request verbose display of results from optimizations. Values are:
```

"off" [default]

```

No display.
"iter" Display intermediate results for every loop iteration.
"final" Display the result of the final loop iteration.
"notify" Display the result of the final loop iteration if the function has failed to converge.
FinDiffType
FunValCheck
When enabled, display an error if the objective function returns an invalid value (a complex number, NaN, or Inf). Must be set to "on" or "off" [default]. Note: the functions fzero and fminbnd correctly handle Inf values and only complex values or NaN will cause an error in this case.

GradObj When set to "on", the function to be minimized must return a second argument which is the gradient, or first derivative, of the function at the point x. If set to "off" [default], the gradient is computed via finite differences.

Jacobian When set to "on", the function to be minimized must return a second argument which is the Jacobian, or first derivative, of the function at the point x. If set to "off" [default], the Jacobian is computed via finite differences.

MaxFunEvals
Maximum number of function evaluations before optimization stops. Must be a positive integer.

MaxIter Maximum number of algorithm iterations before optimization stops. Must be a positive integer.
OutputFen
A user-defined function executed once per algorithm iteration.
TolFun Termination criterion for the function output. If the difference in the calculated objective function between one algorithm iteration and the next is less than TolFun the optimization stops. Must be a positive scalar.

TolX Termination criterion for the function input. If the difference in \(x\), the current search point, between one algorithm iteration and the next is less than TolX the optimization stops. Must be a positive scalar.

\section*{TypicalX}

Updating
See also: [optimget], page 755 .
optimget (options, parname)
optimget (options, parname, default)
Return the specific option parname from the optimization options structure options created by optimset.
If parname is not defined then return default if supplied, otherwise return an empty matrix.

See also: [optimset], page 753.

\section*{26 Statistics}

Octave has support for various statistical methods. The emphasis is on basic descriptive statistics, but the Octave Forge statistics package includes probability distributions, statistical tests, random number generation, and much more.

The functions that analyze data all assume that multi-dimensional data is arranged in a matrix where each row is an observation, and each column is a variable. Thus, the matrix defined by
\[
\begin{aligned}
\mathrm{a}=\left[\begin{array}{ll}
0.9, & 0.7 \\
0.1, & 0.1 ; \\
0.5, & 0.4
\end{array}\right]
\end{aligned}
\]
contains three observations from a two-dimensional distribution. While this is the default data arrangement, most functions support different arrangements.

It should be noted that the statistics functions don't test for data containing NaN, NA, or Inf. These values need to be detected and dealt with explicitly. See [isnan], page 512, [isna], page 45, [isinf], page 512, [isfinite], page 513.

\subsection*{26.1 Descriptive Statistics}

One principal goal of descriptive statistics is to represent the essence of a large data set concisely. Octave provides the mean, median, and mode functions which all summarize a data set with just a single number corresponding to the central tendency of the data.
```

mean (x)
mean (x, dim)
mean (x,opt)
mean (x, dim, opt)
mean (..., outtype)

```

Compute the mean of the elements of the vector x .
The mean is defined as
\[
\operatorname{mean}(x)=\bar{x}=\frac{1}{N} \sum_{i=1}^{N} x_{i}
\]
where \(N\) is the number of elements of \(x\).
If \(x\) is a matrix, compute the mean for each column and return them in a row vector. If the optional argument dim is given, operate along this dimension.
The optional argument opt selects the type of mean to compute. The following options are recognized:
"a" Compute the (ordinary) arithmetic mean. [default]
"g" Compute the geometric mean.
"h" Compute the harmonic mean.

The optional argument outtype selects the data type of the output value. The following options are recognized:
"default"
Output will be of class double unless \(x\) is of class single, in which case the output will also be single.
"double" Output will be of class double.
"native" Output will be the same class as x unless x is of class logical in which case it returns of class double.

Both dim and opt are optional. If both are supplied, either may appear first.
See also: [median], page 758, [mode], page 758.
median (x)
median ( \(x\), dim)
Compute the median value of the elements of the vector \(x\).
When the elements of \(x\) are sorted, say \(s=\operatorname{sort}(x)\), the median is defined as
\[
\operatorname{median}(x)= \begin{cases}s(\lceil N / 2\rceil), & N \text { odd; } \\ (s(N / 2)+s(N / 2+1)) / 2, & N \text { even }\end{cases}
\]
where \(N\) is the number of elements of \(x\).
If \(x\) is of a discrete type such as integer or logical, then the case of even \(N\) rounds up (or toward true).
If \(x\) is a matrix, compute the median value for each column and return them in a row vector.
If the optional dim argument is given, operate along this dimension.
See also: [mean], page 757, [mode], page 758.
mode ( \(x\) )
mode ( \(x\), dim)
\([m, f, c]=\operatorname{mode}(. .\).
Compute the most frequently occurring value in a dataset (mode).
mode determines the frequency of values along the first non-singleton dimension and returns the value with the highest frequency. If two, or more, values have the same frequency mode returns the smallest.
If the optional argument dim is given, operate along this dimension.
The return variable \(f\) is the number of occurrences of the mode in the dataset.
The cell array \(c\) contains all of the elements with the maximum frequency.
See also: [mean], page 757, [median], page 758.
Using just one number, such as the mean, to represent an entire data set may not give an accurate picture of the data. One way to characterize the fit is to measure the dispersion of the data. Octave provides several functions for measuring dispersion.
[s, l] = bounds (x)
[s, l] = bounds (x, dim)
[s, l] = bounds (..., "nanflag")
Return the smallest and largest values of the input data \(x\).
If \(x\) is a vector, the bounds are calculated over the elements of \(x\). If \(x\) is a matrix, the bounds are calculated for each column. For a multi-dimensional array, the bounds are calculated over the first non-singleton dimension.
If the optional argument dim is given, operate along this dimension.
The optional argument "nanflag" defaults to "omitnan" which does not include NaN values in the result. If the argument "includenan" is given, and there is a NaN present, then the result for both smallest ( \(s\) ) and largest (1) elements will be NaN.
The bounds are a quickly computed measure of the dispersion of a data set, but are less accurate than iqr if there are outlying data points.

See also: [range], page 759, [iqr], page 759, [mad], page 760, [std], page 760.
range ( \(x\) )
range ( \(x\), dim)
Return the range, i.e., the difference between the maximum and the minimum of the input data.
If \(x\) is a vector, the range is calculated over the elements of \(x\). If \(x\) is a matrix, the range is calculated over each column of \(x\).
If the optional argument dim is given, operate along this dimension.
The range is a quickly computed measure of the dispersion of a data set, but is less accurate than iqr if there are outlying data points.
See also: [bounds], page 758, [iqr], page 759, [mad], page 760, [std], page 760.
\(Z=i q r(x)\)
\(Z=i q r(x, \operatorname{dim})\)
\(Z=i q r(x\), "ALL")
Return the interquartile range of \(x\), defined as the distance between the 25th and 75 th percentile values of \(x\) calculated using: quantile ( x , [0.25 0.75])
If \(x\) is a vector, iqr ( \(x\) ) will operate on the data in \(x\).
If \(x\) is a matrix, iqr ( \(x\) ) will operate independently on each column in \(x\) returning a row vector \(Z\).
If \(x\) is a n-dimensional array, iqr ( \(x\) ) will operate independently on the first nonsingleton dimension in \(x\), returning an array \(Z\) the same shape as \(x\) with the nonsingleton dimenion reduced to 1 .
The optional variable dim can be used to force iqr to operate over the specified dimension. dim can either be a scalar dimension or a vector of non-repeating dimensions over which to operate. In either case dim must be positive integers. A vector dim concatenates all specified dimensions for independent operation by iqr.
Specifying dimension "ALL" will force iqr to operate on all elements of \(x\), and is equivalent to iqr ( \(x(:)\) ). Similarly, specifying a vector dimension including all nonsingleton dimensions of \(x\) is equivalent to iqr ( \(x\), "ALL").

If \(x\) is a scalar, or only singleton dimensions are specified for dim, the output will be zeros (size (x)).
As a measure of dispersion, the interquartile range is less affected by outliers than either range or std.
See also: [bounds], page 758, [mad], page 760, [range], page 759, [std], page 760, [prctile], page 765, [quantile], page 764.
mad ( \(x\) )
mad ( \(x\), opt)
mad ( \(x\), opt, dim)
Compute the mean or median absolute deviation of the elements of \(x\).
The mean absolute deviation is defined as
```

mad = mean (abs (x - mean (x)))

```

The median absolute deviation is defined as
```

mad = median (abs (x - median (x)))

```

If \(x\) is a matrix, compute mad for each column and return results in a row vector. For a multi-dimensional array, the calculation is done over the first non-singleton dimension.
The optional argument opt determines whether mean or median absolute deviation is calculated. The default is 0 which corresponds to mean absolute deviation; A value of 1 corresponds to median absolute deviation.
If the optional argument dim is given, operate along this dimension.
As a measure of dispersion, mad is less affected by outliers than std.
See also: [bounds], page 758, [range], page 759, [iqr], page 759, [std], page 760, [mean], page 757, [median], page 758 .
meansq ( \(x\) )
meansq ( \(x\), dim)
Compute the mean square of the elements of the vector \(x\).
The mean square is defined as
\[
\operatorname{meansq}(x)=\frac{\sum_{i=1}^{N} x_{i}{ }^{2}}{N}
\]
where \(N\) is the number of elements of \(x\).
If \(x\) is a matrix, return a row vector containing the mean square of each column.
If the optional argument dim is given, operate along this dimension.
See also: [var], page 761, [std], page 760, [moment], page 763.
std (x)
std ( \(x\), w)
std ( \(x\), w, dim)
std (x, w, "ALL")
Compute the standard deviation of the elements of the vector x .

The standard deviation is defined as
\[
\operatorname{std}(x)=\sigma=\sqrt{\frac{\sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{2}}{N-1}}
\]
where \(\bar{x}\) is the mean value of x and \(N\) is the number of elements of x .
If \(x\) is an array, compute the standard deviation for each column and return them in a row vector (or for an n-D array, the result is returned as an array of dimension 1 x n x m x ...).
The optional argument \(w\) determines the weighting scheme to use. Valid values are: 0 [default]:

Normalize with \(N-1\). This provides the square root of the best unbiased estimator of the variance.

1: \(\quad\) Normalize with \(N\). This provides the square root of the second moment around the mean.
a vector: Compute the weighted standard deviation with nonnegative scalar weights. The length of \(w\) must be equal to the size of \(x\) along dimension dim.

If \(N\) is equal to 1 the value of \(W\) is ignored and normalization by \(N\) is used.
The optional variable dim forces std to operate over the specified dimension. dim can either be a scalar dimension or a vector of non-repeating dimensions over which to operate. Dimensions must be positive integers, and the standard deviation is calculated over the array slice defined by dim.
Specifying dimension "ALL" will force std to operate on all elements of \(x\), and is equivalent to \(\operatorname{std}(x(:))\).
When dim is a vector or "ALL", w must be either 0 or 1 .
See also: [var], page 761, [bounds], page 758, [mad], page 760, [range], page 759, [iqr], page 759, [mean], page 757, [median], page 758.

In addition to knowing the size of a dispersion it is useful to know the shape of the data set. For example, are data points massed to the left or right of the mean? Octave provides several common measures to describe the shape of the data set. Octave can also calculate moments allowing arbitrary shape measures to be developed.
\(\operatorname{var}(x)\)
\(\operatorname{var}(x, w)\)
var (x, w, dim)
var (x, w, "ALL")
Compute the variance of the elements of the vector \(x\).
The variance is defined as
\[
\operatorname{var}(x)=\sigma^{2}=\frac{\sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{2}}{N-1}
\]
where \(\bar{x}\) is the mean value of x and \(N\) is the number of elements of x .

If \(x\) is an array, compute the variance for each column and return them in a row vector (or for an n-D array, the result is returned as an array of dimension \(1 \times \mathrm{nxm}\) x ...).
The optional argument \(w\) determines the weighting scheme to use. Valid values are 0 [default]:

Normalize with \(N-1\). This provides the square root of the best unbiased estimator of the variance.

1: \(\quad\) Normalize with \(N\), this provides the square root of the second moment around the mean
a vector: Compute the weighted variance with nonnegative scalar weights. The length of \(w\) must be equal to the size of \(x\) along dimension dim.

If \(N\) is equal to 1 the value of \(W\) is ignored and normalization by \(N\) is used.
The optional variable dim forces var to operate over the specified dimension. dim can either be a scalar dimension or a vector of non-repeating dimensions over which to operate. Dimensions must be positive integers, and the variance is calculated over the array slice defined by dim.
Specifying dimension "ALL" will force var to operate on all elements of \(x\), and is equivalent to \(\operatorname{var}(x(:))\).
When dim is a vector or "ALL", w must be either 0 or 1 .
See also: [cov], page 786, [std], page 760, [skewness], page 762, [kurtosis], page 763, [moment], page 763.
```

skewness (x)
skewness (x, flag)
skewness (x, flag, dim)

```

Compute the sample skewness of the elements of \(x\).
The sample skewness is defined as
\[
\operatorname{skewness}(x)=\frac{\frac{1}{N} \sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{3}}{\sigma^{3}}
\]
where \(N\) is the length of \(x, \bar{x}\) its mean and \(\sigma\) its (uncorrected) standard deviation.
The optional argument flag controls which normalization is used. If flag is equal to 1 (default value, used when flag is omitted or empty), return the sample skewness as defined above. If flag is equal to 0 , return the adjusted skewness coefficient instead:
\[
\operatorname{skewness}(\mathrm{x})=\frac{\sqrt{N(N-1)}}{N-2} \times \frac{\frac{1}{N} \sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{3}}{\sigma^{3}}
\]

The adjusted skewness coefficient is obtained by replacing the sample second and third central moments by their bias-corrected versions.
If \(x\) is a matrix, or more generally a multi-dimensional array, return the skewness along the first non-singleton dimension. If the optional dim argument is given, operate along this dimension.
See also: [var], page 761, [kurtosis], page 763, [moment], page 763.
kurtosis (x)
kurtosis (x, flag)
kurtosis (x, flag, dim)
Compute the sample kurtosis of the elements of \(x\).
The sample kurtosis is defined as
\[
\kappa_{1}=\frac{\frac{1}{N} \sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{4}}{\sigma^{4}}
\]
where \(N\) is the length of \(x, \bar{x}\) its mean, and \(\sigma\) its (uncorrected) standard deviation.
The optional argument flag controls which normalization is used. If flag is equal to 1 (default value, used when flag is omitted or empty), return the sample kurtosis as defined above. If flag is equal to 0 , return the "bias-corrected" kurtosis coefficient instead:
\[
\kappa_{0}=3+\frac{N-1}{(N-2)(N-3)}\left((N+1) \kappa_{1}-3(N-1)\right)
\]

The bias-corrected kurtosis coefficient is obtained by replacing the sample second and fourth central moments by their unbiased versions. It is an unbiased estimate of the population kurtosis for normal populations.
If \(x\) is a matrix, or more generally a multi-dimensional array, return the kurtosis along the first non-singleton dimension. If the optional dim argument is given, operate along this dimension.
See also: [var], page 761, [skewness], page 762, [moment], page 763.
```

moment (x, p)
moment (x, p, type)
moment (x, p, dim)
moment (x, p, type, dim)
moment (x, p, dim, type)

```

Compute the \(p\)-th central moment of the vector \(x\).
The \(p\)-th central moment of \(x\) is defined as:
\[
\frac{\sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)^{p}}{N}
\]
where \(\bar{x}\) is the mean value of x and \(N\) is the number of elements of x .
If \(x\) is a matrix, return the row vector containing the \(p\)-th central moment of each column.
If the optional argument dim is given, operate along this dimension.
The optional string type specifies the type of moment to be computed. Valid options are:
"c" Central Moment (default).
"a"
"ac" Absolute Central Moment. The moment about the mean ignoring sign defined as
\[
\frac{\sum_{i=1}^{N}\left|x_{i}-\bar{x}\right|^{p}}{N}
\]
"r" Raw Moment. The moment about zero defined as
\[
\operatorname{moment}(x)=\frac{\sum_{i=1}^{N} x_{i}^{p}}{N}
\]
"ar" Absolute Raw Moment. The moment about zero ignoring sign defined as
\[
\frac{\sum_{i=1}^{N}\left|x_{i}\right|^{p}}{N}
\]

If both type and dim are given they may appear in any order.
See also: [var], page 761, [skewness], page 762, [kurtosis], page 763.
```

q = quantile (x)
q = quantile (x, p)
q = quantile (x, p, dim)
q = quantile (x, p, dim, method)

```

For a sample, \(x\), calculate the quantiles, \(q\), corresponding to the cumulative probability values in \(p\). All non-numeric values ( NaNs ) of \(x\) are ignored.
If \(x\) is a matrix, compute the quantiles for each column and return them in a matrix, such that the i-th row of \(q\) contains the \(p(\mathrm{i})\) th quantiles of each column of x .
If \(p\) is unspecified, return the quantiles for [0.00 0.250 .500 .751 .00 ]. The optional argument dim determines the dimension along which the quantiles are calculated. If dim is omitted it defaults to the first non-singleton dimension.
The methods available to calculate sample quantiles are the nine methods used by \(R\) (https://www.r-project.org/). The default value is method \(=5\).
Discontinuous sample quantile methods 1 , 2, and 3
1. Method 1: Inverse of empirical distribution function.
2. Method 2: Similar to method 1 but with averaging at discontinuities.
3. Method 3: SAS definition: nearest even order statistic.

Continuous sample quantile methods 4 through 9 , where \(p(k)\) is the linear interpolation function respecting each method's representative cdf.
4. Method 4: \(p(k)=k / N\). That is, linear interpolation of the empirical cdf, where \(N\) is the length of \(P\).
5. Method 5: \(p(k)=(k-0.5) / N\). That is, a piecewise linear function where the knots are the values midway through the steps of the empirical cdf.
6. Method 6: \(p(k)=k /(N+1)\).
7. Method 7: \(p(k)=(k-1) /(N-1)\).
8. Method 8: \(p(k)=(k-1 / 3) /(N+1 / 3)\). The resulting quantile estimates are approximately median-unbiased regardless of the distribution of \(x\).
9. Method 9: \(p(k)=(k-3 / 8) /(N+1 / 4)\). The resulting quantile estimates are approximately unbiased for the expected order statistics if \(x\) is normally distributed.

Hyndman and Fan (1996) recommend method 8. Maxima, S, and R (versions prior to 2.0.0) use 7 as their default. Minitab and SPSS use method 6. matlab uses method 5.

References:
- Becker, R. A., Chambers, J. M. and Wilks, A. R. (1988) The New S Language. Wadsworth \& Brooks/Cole.
- Hyndman, R. J. and Fan, Y. (1996) Sample quantiles in statistical packages, American Statistician, 50, 361-365.
- R: A Language and Environment for Statistical Computing; https://cran. r-project.org/doc/manuals/fullrefman.pdf.

Examples:
\[
\begin{aligned}
& \mathrm{x}=\operatorname{randi}(1000,[10,1]) ; \quad \text { \# Create empirical data in range } 1-1000 \\
& \mathrm{q}=\text { quantile }(\mathrm{x},[0,1]) ; \text { \# Return minimum, maximum of distribution } \\
& \mathrm{q}=\text { quantile }(\mathrm{x},[0.250 .50 .75]) ; \text { \# Return quartiles of distribution }
\end{aligned}
\]

See also: [prctile], page 765 .
```

q = prctile (x)
q = prctile (x, p)
q = prctile (x, p, dim)

```

For a sample \(x\), compute the quantiles, \(q\), corresponding to the cumulative probability values, \(p\), in percent.

If \(x\) is a matrix, compute the percentiles for each column and return them in a matrix, such that the i-th row of \(q\) contains the \(p(i)\) th percentiles of each column of \(x\).

If \(p\) is unspecified, return the quantiles for [0 255075 100].
The optional argument dim determines the dimension along which the percentiles are calculated. If dim is omitted it defaults to the first non-singleton dimension.
Programming Note: All non-numeric values (NaNs) of \(x\) are ignored.
See also: [quantile], page 764.
A summary view of a data set can be generated quickly with the statistics function.
```

statistics (x)
statistics (x, dim)

```

Return a vector with the minimum, first quartile, median, third quartile, maximum, mean, standard deviation, skewness, and kurtosis of the elements of the vector \(x\).
If \(x\) is a matrix, calculate statistics over the first non-singleton dimension.
If the optional argument dim is given, operate along this dimension.
See also: [min], page 559, [max], page 558, [median], page 758, [mean], page 757, [std], page 760, [skewness], page 762, [kurtosis], page 763.

\subsection*{26.2 Statistics on Sliding Windows of Data}

It is often useful to calculate descriptive statistics over a subsection (i.e., window) of a full dataset. Octave provides the function movfun which will call an arbitrary function handle with windows of data and accumulate the results. Many of the most commonly desired functions, such as the moving average over a window of data (movmean), are already provided.
```

y = movfun (fcn, x, wlen)
y = movfun (fcn, x, [nb, na])
y = movfun (..., "property", value)

```

Apply function fcn to a moving window of length wlen on data x .
If wlen is a scalar, the function \(f c n\) is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen-1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movfun uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movfun uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
During calculations the data input x is reshaped into a 2 -dimensional wlen-by- \(N\) matrix and \(f c n\) is called on this new matrix. Therefore, fcn must accept an array input argument and apply the computation along dimension 1, i.e., down the columns of the array.
When applied to an array (possibly multi-dimensional) with \(n\) columns, fcn may return a result in either of two formats: Format 1) an array of size 1-by-n-by-dim3-by-. . .-by-dimN. This is the typical output format from Octave core functions. Type demo ("movfun", 5) for an example of this use case. Format 2) a row vector of length \(n *\) numel_higher_dims where numel_higher_dims is prod (size ( \(x\) )(3:end)). The output of fcn for the i-th input column must be found in the output at indices i:n: (n*numel_higher_dims). This format is useful when concatenating functions into arrays, or when using nthargout. Type demo ("movfun", 6) for an example of this case.
The calculation can be controlled by specifying property/value pairs. Valid properties are
"dim" Operate along the dimension specified, rather than the default of the first non-singleton dimension.
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=f c n(x(1: 2))\), and \(y\) (end) \(=f c n(x(\) end -1 :end \()\) ).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) \(f c n([\mathrm{NaN}, x(1: 2)])\), and \(y(e n d)=f c n\) ([x(end-1:end), \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how fcn handles NaN, and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\) \(=f c n\left(\left[u s e r_{-} v a l u e, x(1: 2)\right]\right)\) and \(y(e n d)=f c n\) ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=f c n([x(1), x(1: 2)])\), and \(y\) (end) \(=f c n([x(\) end -1 :end \(), x(e n d)])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=f c n([x(e n d), x(1: 2)])\), and \(y(\) end \()=f c n([x(\) end -1 :end \(), x(1)])\).
Note that for some of these choices, the window size at the boundaries is not the same as for the central part, and fcn must work in these cases.
"nancond"
Controls whether NaN and NA values should be included (value: "includenan"), or excluded (value: "omitnan"), from the data passed to fcn. The default is "includenan". Caution: The "omitnan" option is not yet implemented.
"outdim" A row vector that selects which dimensions of the calculation will appear in the output \(y\). This is only useful when fcn returns an N -dimensional array in Format 1. The default is to return all output dimensions.

Programming Note: The property "outdim" can be used to save memory when the output of fcn has many dimensions, or when a wrapper to the base function that selects the desired outputs is too costly. When memory is not an issue, the easiest way to select output dimensions is to first calculate the complete result with movfun and then filter that result with indexing. If code complexity is not an issue then a wrapper can be created using anonymous functions. For example, if basefcn is a function returning a \(K\)-dimensional row output, and only dimension \(D\) is desired, then the following wrapper could be used.
```

fcn =@(x) basefcn (x)(:,columns(x) * (D-1) + (1:columns(x)));
y = movfun (@fcn, ...);

```

See also: [movslice], page 768, [prepad], page 524, [postpad], page 525, [permute], page 518, [reshape], page 518.
slcidx = movslice (N, wlen)
[slcidx, C, Cpre, Cpost, win] = movslice (...)
Generate indices to slice a vector of length \(N\) into windows of length wlen.
The input \(N\) must be a positive integer.
The moving window length input wlen can either be a scalar not equal to 1 or a 2 -element array of integers. For scalar values, if odd the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. If wlen is even the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. When wlenis a 2 -element array, [nb, na], the window includes \(n b\) elements to the left of the current element and na elements to the right of the current element.
The output slcidx is an array of indices of the slices that fit fully within the vector, where each column is an individual slice as the window moves from left to right. The slices have wlen elements for scalar wlen, or \(n b+n a+1\) elements for array valued wlen.

Optional output \(C\) is an row vector of window center positions where the window stays fully within the vector.
Optional outputs Cpre and Cpost contain the vector elements at the start and end of the vector, respectively, that result in the window extending beyond the ends of the vector.
Optional output win is a column vector with the same number of rows as slcidx that contains the moving window defined as a center relative position stencil.

See also: [movfun], page 766.
```

y = movmad (x, wlen)
y = movmad (x, [nb, na])
y = movmad (..., dim)
y = movmad (..., "nancond")
y = movmad (..., property, value)

```

Calculate the moving mean absolute deviation over a sliding window of length wlen on data \(x\).

If wlen is a scalar, the function mad is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movmad uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movmad uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to mad. The default is "includenan". Caution: the "omitnan" option is not yet implemented. The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=\operatorname{mad}(x(1: 2))\), and \(y(e n d)=\operatorname{mad}(x(\) end \(-1:\) end \())\).

Any y values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN . For example, with a window of length \(3, \mathrm{y}(1)=\) \(\operatorname{mad}([\mathrm{NaN}, x(1: 2)])\), and \(y(\) end \()=\operatorname{mad}([x(e n d-1: e n d)\), \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how mad handles NaN, and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\)
\(=\operatorname{mad}([\) user_value, \(x(1: 2)])\) and \(y(e n d)=\operatorname{mad}\) ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\operatorname{mad}([x(1), x(1: 2)])\), and \(y\) (end)
\(=\operatorname{mad}([x(\) end \(-1:\) end \(), x(\) end \()])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\operatorname{mad}([x(e n d), x(1: 2)])\), and \(y(\) end \()=\operatorname{mad}([x(\) end \(-1:\) end \(), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmax], page 770, [movmean], page 772 , [movmedian], page 773 , [movmin], page 775 , [movprod], page 776 , [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movmax (x, wlen)
y = movmax (x, [nb, na])
y = movmax (..., dim)
y = movmax (..., "nancond")
y = movmax (..., property, value)

```

Calculate the moving maximum over a sliding window of length wlen on data \(x\).
If wlen is a scalar, the function max is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movmax uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movmax uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [ \(n b, n a\) ], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).

If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to max. The default is "includenan". Caution: the "omitnan" option is not yet implemented.

The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=\max (x(1: 2))\), and \(y(\) end \()=\max (x(\) end -1 :end \()\) ).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN . For example, with a window of length \(3, y(1)=\) \(\max ([\mathrm{NaN}, x(1: 2)])\), and \(y(\) end \()=\max ([x(\) end \(-1:\) end), \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how max handles NaN , and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length 3, \(y(1)\) \(=\max \left(\left[u s e r_{-}\right.\right.\)value, \(\left.\left.x(1: 2)\right]\right)\), and \(y(e n d)=\max\) ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\max ([x(1), x(1: 2)])\), and \(y(e n d)\) \(=\max ([x(\) end \(-1:\) end \(), x(e n d)])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\max ([x(e n d), x(1: 2)])\), and \(y(\) end \()=\max ([x(\) end -1 :end \(), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmean], page 772, [movmedian], page 773, [movmin], page 775, [movprod], page 776, [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movmean (x, wlen)
y = movmean (x, [nb,na])
y = movmean (..., dim)
y = movmean (..., "nancond")
y = movmean (..., property, value)

```

Calculate the moving average over a sliding window of length wlen on data x .
If wlen is a scalar, the function mean is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen-1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movmean uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movmean uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to mean. The default is "includenan". Caution: the "omitnan" option is not yet implemented.
The calculation can be controlled by specifying property/value pairs. Valid properties are

\section*{"Endpoints"}

This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=\) mean ( \(x(1: 2)\) ), and \(y(\) end \()=\) mean \((x(\) end -1 :end \()\) ).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) mean ([NaN, \(x(1: 2)])\), and \(y(e n d)=\) mean ([x(end-1:end), \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how mean handles NaN, and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\) \(=\) mean \(\left(\left[u s e r_{-}\right.\right.\)value, \(\left.\left.x(1: 2)\right]\right)\), and \(y(e n d)=\) mean ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\) mean \(([x(1), x(1: 2)])\), and \(y(\) end \()=\) mean ([x(end-1:end), \(x(e n d)]\) ).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\) mean ( \([x(\) end),\(x(1: 2)]\) ), and \(y(\) end \()=\) mean \(([x(\) end -1 :end \(), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.
See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmedian], page 773, [movmin], page 775, [movprod], page 776, [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movmedian (x, wlen)
y = movmedian (x, [nb, na])
y = movmedian (..., dim)
y = movmedian (..., "nancond")
y = movmedian (..., property, value)

```

Calculate the moving median over a sliding window of length wlen on data x .
If wlen is a scalar, the function movmedian is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen-1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movmedian uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movmedian uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current
element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to movmedian. The default is "includenan". Caution: the "omitnan" option is not yet implemented.
The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=\) movmedian ( \(x(1: 2)\) ), and \(y(e n d)=\) movmedian ( \(x(\) end \(-1:\) end \()\) ).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN . For example, with a window of length \(3, \mathrm{y}(1)=\) movmedian ([NaN, x(1:2)]), and \(y\) (end) = movmedian ([x(end-1:end), NaN]). This option usually results in y having NaN values at the boundaries, although it is influenced by how movmedian handles NaN , and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\) \(=\) movmedian ([user_value, \(x(1: 2)]\) ), and \(y(e n d)=\) movmedian ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\) movmedian \(([x(1), x(1: 2)])\), and \(y\) (end) \(=\) movmedian ( \([x\) (end-1:end), \(x(e n d)]\).

The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\) movmedian ( \([x\) (end), \(x(1: 2)])\), and \(y(e n d)=\) movmedian ([x(end-1:end), \(x(1)]\) ).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmean], page 772, [movmin], page 775, [movprod], page 776, [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movmin (x, wlen)
y = movmin (x, [nb, na])
y = movmin (..., dim)
y = movmin (..., "nancond")
y = movmin (..., property, value)

```

Calculate the moving minimum over a sliding window of length wlen on data \(x\).
If wlen is a scalar, the function min is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen-1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movmin uses data elements [4, 5, 6]. If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movmin uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to min. The default is "includenan". Caution: the "omitnan" option is not yet implemented. The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For
example, with a window of length \(3, y(1)=\min (x(1: 2))\), and \(y(\) end \()=\min (x(\) end \(-1:\) end \())\).
"discard"
Any y values that use a window extending beyond the original data array are deleted. For example, with a 10-element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices [0, 1, 2] and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) \(\min ([N a N, x(1: 2)])\), and \(y(e n d)=\min ([x(e n d-1: e n d)\), NaN]). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how min handles NaN , and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\) \(=\min \left(\left[u s e r_{-}\right.\right.\)value,\(\left.\left.x(1: 2)\right]\right)\) and \(y(e n d)=\min\) ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\min ([x(1), x(1: 2)])\), and \(y(e n d)\) \(=\min ([x(\) end \(-1:\) end \(), x(\) end \()])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\min ([x(e n d), x(1: 2)])\), and \(y(\) end \()=\min ([x(\) end \(-1:\) end \(), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.
See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770 , [movmean], page 772 , [movmedian], page 773 , [movprod], page 776 , [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movprod (x, wlen)
y = movprod (x, [nb, na])
y = movprod (..., dim)
y = movprod (..., "nancond")

```
```

y = movprod (..., property, value)

```

Calculate the moving product over a sliding window of length wlen on data \(x\).
If wlen is a scalar, the function movprod is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movprod uses data elements [4, 5, 6]. If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movprod uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to movprod. The default is "includenan". Caution: the "omitnan" option is not yet implemented.
The calculation can be controlled by specifying property/value pairs. Valid properties are

\section*{"Endpoints"}

This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, \mathrm{y}(1)=\) movprod \((x(1: 2))\), and \(y(e n d)=\operatorname{movprod}(x(e n d-1: e n d))\).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)\) \(=\operatorname{movprod}([\operatorname{NaN}, x(1: 2)])\) and \(y(e n d)=\operatorname{movprod}\) ([x(end-1:end), NaN]). This option usually results in \(y\) having NaN values at the boundaries, although it is
influenced by how movprod handles NaN , and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)=\) movprod ([user_value, \(x(1: 2)]\) ), and \(y\) (end) = movprod ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\operatorname{movprod}([x(1), x(1: 2)])\), and \(y(e n d)\) \(=\operatorname{movprod}([x(e n d-1:\) end \(), x(e n d)])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\) movprod \(([x(\) end \(), x(1: 2)])\), and \(y(e n d)=\operatorname{movprod}([x(e n d-1: e n d), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmean], page 772, [movmedian], page 773 , [movmin], page 775 , [movstd], page 778, [movsum], page 780, [movvar], page 782.
```

y = movstd (x, wlen)
y = movstd (x, [nb, na])
y = movstd (..., opt)
y = movstd (..., opt, dim)
y = movstd (..., "nancond")
y = movstd (..., property, value)

```

Calculate the moving standard deviation over a sliding window of length wlen on data x.

If wlen is a scalar, the function movstd is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movstd uses data elements [4, 5, 6]. If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movstd uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is
always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
The optional argument opt determines the type of normalization to use. Valid values are

0: \(\quad\) normalize with \(N-1\), provides the square root of the best unbiased estimator of the variance [default]
1: \(\quad\) normalize with \(N\), this provides the square root of the second moment around the mean

If the optional argument dim is given, operate along this dimension. The normalization argument opt must be given before the dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to movstd. The default is "includenan". Caution: the "omitnan" option is not yet implemented.
The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, \mathrm{y}(1)=\) movstd \((x(1: 2))\), and \(y(e n d)=\) movstd \((x(e n d-1: e n d))\).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) movstd ([NaN, \(x(1: 2)]\) ), and \(y\) (end) \(=\) movstd ([x(end-1:end), NaN]). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how movstd handles NaN, and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)=\) movstd ([user_value, \(x(1: 2)]\) ), and \(y(e n d)=\) movstd
([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\operatorname{movstd}([x(1), x(1: 2)])\), and \(y\) (end) \(=\operatorname{movstd}([x(e n d-1:\) end \(), x(e n d)])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\) movstd \(([x(\) end \(), x(1: 2)])\), and \(y(e n d)=\operatorname{movstd}([x(e n d-1: e n d), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmean], page 772, [movmedian], page 773 , [movmin], page 775 , [movprod], page 776, [movsum], page 780, [movvar], page 782.
```

y = movsum (x, wlen)
y = movsum (x, [nb, na])
y = movsum (..., dim)
y = movsum (..., "nancond")
y = movsum (..., property, value)

```

Calculate the moving sum over a sliding window of length wlen on data \(x\).
If wlen is a scalar, the function movsum is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movsum uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movsum uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
If the optional argument dim is given, operate along this dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to movsum. The default is "includenan". Caution: the "omitnan" option is not yet implemented.

The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, \mathrm{y}(1)=\) movsum ( \(x(1: 2)\) ), and \(y(\) end \()=\) movsum ( \(x(\) end \(-1:\) end) \()\).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) movsum ([NaN, \(x(1: 2)]\) ), and \(y\) (end) \(=\) movsum ([x(end-1:end), \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how movsum handles NaN, and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)=\) movsum ([user_value, \(x(1: 2)]\) ), and \(y(e n d)=\) movsum ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\) movsum \(([x(1), x(1: 2)])\), and \(y(e n d)\) = movsum ([x(end-1:end), \(x(e n d)]\) ).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\) movsum ( \([x(e n d), x(1: 2)]\) ), and \(y(e n d)=\) movsum \(([x(\) end -1 :end \(), x(1)])\).
"SamplePoints"
Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmean], page 772, [movmedian], page 773, [movmin], page 775, [movprod], page 776, [movstd], page 778, [movvar], page 782.
```

y = movvar (x, wlen)
y = movvar (x, [nb, na])
y = movvar (...,opt)
y = movvar (..., opt, dim)
y = movvar (..., "nancond")
y = movvar (..., property, value)

```

Calculate the moving variance over a sliding window of length wlen on data \(x\).
If wlen is a scalar, the function var is applied to a moving window of length wlen. When wlen is an odd number the window is symmetric and includes (wlen - 1) / 2 elements on either side of the central element. For example, when calculating the output at index 5 with a window length of 3 , movvar uses data elements \([4,5,6]\). If wlen is an even number, the window is asymmetric and has wlen/2 elements to the left of the central element and wlen/2-1 elements to the right of the central element. For example, when calculating the output at index 5 with a window length of 4 , movvar uses data elements \([3,4,5,6]\).
If wlen is an array with two elements [nb, na], the function is applied to a moving window -nb:na. This window includes nb number of elements before the current element and na number of elements after the current element. The current element is always included. For example, given wlen \(=[3,0]\), the data used to calculate index 5 is \([2,3,4,5]\).
The optional argument opt determines the type of normalization to use. Valid values are

0: \(\quad\) normalize with \(N-1\), provides the best unbiased estimator of the variance [default]
1: \(\quad\) normalizes with \(N\), this provides the second moment around the mean
If the optional argument dim is given, operate along this dimension. The normalization argument opt must be given before the dimension.
The optional string argument "nancond" controls whether NaN and NA values should be included ("includenan"), or excluded ("omitnan"), from the data passed to var. The default is "includenan". Caution: the "omitnan" option is not yet implemented.
The calculation can be controlled by specifying property/value pairs. Valid properties are
"Endpoints"
This property controls how results are calculated at the boundaries (endpoints) of the window. Possible values are:
"shrink" (default)
The window is truncated at the beginning and end of the array to exclude elements for which there is no source data. For example, with a window of length \(3, y(1)=\operatorname{var}(x(1: 2))\), and \(y(\) end \()=\operatorname{var}(x(\) end \(-1:\) end \())\).
"discard"
Any \(y\) values that use a window extending beyond the original data array are deleted. For example, with a 10 -element data vector and a window of length 3 , the output will contain only 8 elements. The first element would require calculating the function over indices \([0,1,2]\) and is therefore discarded. The last element would require calculating the function over indices \([9,10,11]\) and is therefore discarded.
"fill" Any window elements outside the data array are replaced by NaN. For example, with a window of length \(3, y(1)=\) \(\operatorname{var}([\mathrm{NaN}, x(1: 2)])\), and \(y(\) end \()=\operatorname{var}([x(\) end \(-1:\) end) , \(\mathrm{NaN}]\) ). This option usually results in \(y\) having NaN values at the boundaries, although it is influenced by how var handles NaN , and also by the property "nancond".
user_value Any window elements outside the data array are replaced by the specified value user_value which must be a numeric scalar. For example, with a window of length \(3, y(1)\) \(=\operatorname{var}\left(\left[u s e r_{-}\right.\right.\)value, \(\left.\left.x(1: 2)\right]\right)\), and \(y(e n d)=\operatorname{var}\) ([x(end-1:end), user_value]). A common choice for user_value is 0 .
"same" Any window elements outside the data array are replaced by the value of \(x\) at the boundary. For example, with a window of length \(3, y(1)=\operatorname{var}([x(1), x(1: 2)])\), and \(y(e n d)\) \(=\operatorname{var}([x(e n d-1: e n d), x(e n d)])\).
"periodic"
The window is wrapped so that any missing data elements are taken from the other side of the data. For example, with a window of length \(3, y(1)=\operatorname{var}([x(\) end \(), x(1: 2)])\), and \(y(\) end \()=\operatorname{var}([x(\) end \(-1:\) end \(), x(1)])\).

\section*{"SamplePoints"}

Caution: This option is not yet implemented.
Programming Note: This function is a wrapper which calls movfun. For additional options and documentation, see [movfun], page 766.

See also: [movfun], page 766, [movslice], page 768, [movmad], page 768, [movmax], page 770, [movmean], page 772, [movmedian], page 773, [movmin], page 775, [movprod], page 776, [movstd], page 778, [movsum], page 780.

\subsection*{26.3 Basic Statistical Functions}

Octave supports various helpful statistical functions. Many are useful as initial steps to prepare a data set for further analysis. Others provide different measures from those of the basic descriptive statistics.
center ( \(x\) )
center ( \(x\), dim)
Center data by subtracting its mean.
If \(x\) is a vector, subtract its mean.
If \(x\) is a matrix, do the above for each column.
If the optional argument dim is given, operate along this dimension.
Programming Note: center has obvious application for normalizing statistical data. It is also useful for improving the precision of general numerical calculations. Whenever there is a large value that is common to a batch of data, the mean can be subtracted off, the calculation performed, and then the mean added back to obtain the final answer.

See also: [zscore], page 784.
\(z=\) zscore ( \(x\) )
\(z=\) zscore (x, opt)
\(z=\) zscore (x, opt, dim)
[z, mu, sigma] = zscore (...)
Compute the Z score of x .
If \(x\) is a vector, subtract its mean and divide by its standard deviation. If the standard deviation is zero, divide by 1 instead.
The optional parameter opt determines the normalization to use when computing the standard deviation and has the same definition as the corresponding parameter for std.
If \(x\) is a matrix, calculate along the first non-singleton dimension. If the third optional argument dim is given, operate along this dimension.
The optional outputs mu and sigma contain the mean and standard deviation.
See also: [mean], page 757, [std], page 760, [center], page 783.
\(n=\) histc ( \(x\), edges )
\(n=\) histc ( \(x\), edges, dim)
[n, idx] = histc (...)
Compute histogram counts.
When \(x\) is a vector, the function counts the number of elements of \(x\) that fall in the histogram bins defined by edges. This must be a vector of monotonically increasing values that define the edges of the histogram bins. \(n(k)\) contains the number of elements in x for which \(\operatorname{edges}(k)<=\mathrm{x}<\operatorname{edges}(k+1)\). The final element of \(n\) contains the number of elements of \(x\) exactly equal to the last element of edges.
When x is an \(N\)-dimensional array, the computation is carried out along dimension dim. If not specified dim defaults to the first non-singleton dimension.
When a second output argument is requested an index matrix is also returned. The idx matrix has the same size as \(x\). Each element of idx contains the index of the histogram bin in which the corresponding element of \(x\) was counted.
See also: [hist], page 323.
unique function documented at [unique], page 791, is often useful for statistics.
\(c=\) nchoosek (n, k)
\(c=\) nchoosek (set, k)
Compute the binomial coefficient of \(n\) or list all possible combinations of a set of items.
If \(n\) is a scalar then calculate the binomial coefficient of \(n\) and \(k\) which is defined as
\[
\binom{n}{k}=\frac{n(n-1)(n-2) \cdots(n-k+1)}{k!}=\frac{n!}{k!(n-k)!}
\]

This is the number of combinations of \(n\) items taken in groups of size \(k\).
If the first argument is a vector, set, then generate all combinations of the elements of set, taken \(k\) at a time, with one row per combination. The result \(c\) has \(k\) columns and nchoosek (length (set), k) rows.
For example:
How many ways can three items be grouped into pairs?
```

nchoosek (3, 2)
=>3

```

What are the possible pairs?
```

nchoosek (1:3, 2)
\# 1 2
1 3
2 3

```

Programming Note: When calculating the binomial coefficient nchoosek works only for non-negative, integer arguments. Use bincoeff for non-integer and negative scalar arguments, or for computing many binomial coefficients at once with vector inputs for \(n\) or \(k\).

See also: [bincoeff], page 570, [perms], page 785.
perms (v)
Generate all permutations of vector \(v\) with one row per permutation.
Results are returned in inverse lexicographic order. The result has size factorial ( \(n\) ) * \(n\), where \(n\) is the length of \(v\). Any repetitions are included in the output. To generate just the unique permutations use unique (perms (v), "rows") (end:-1:1,:).
Example
\begin{tabular}{lll} 
perms \\
\(\Rightarrow\) & \(([1,2,3])\) \\
3 & 2 & 1 \\
3 & 1 & 2 \\
2 & 3 & 1 \\
2 & 1 & 3 \\
1 & 3 & 2 \\
1 & 2 & 3
\end{tabular}

Programming Note: The maximum length of \(v\) should be less than or equal to 10 to limit memory consumption.
See also: [permute], page 518, [randperm], page 537, [nchoosek], page 785.
ranks (x)
ranks (x, dim)
ranks (x, dim, rtype)
Return the ranks (in the sense of order statistics) of \(x\) along the first non-singleton dimension adjusted for ties.

If the optional dim argument is given, operate along this dimension.
The optional parameter rtype determines how ties are handled. All examples below assume an input of [1, 2, 2, 4].
0 or "fractional" (default) for fractional ranking (1, 2.5, 2.5, 4);

1 or "competition" for competition ranking (1, 2, 2, 4);
2 or "modified" for modified competition ranking (1, 3, 3, 4);
3 or "ordinal" for ordinal ranking ( \(1,2,3,4\) );
4 or "dense" for dense ranking (1, 2, 2, 3).
See also: [spearman], page 788, [kendall], page 788.
run_count ( \(x, n\) )
run_count ( \(x, n\), dim)
Count the upward runs along the first non-singleton dimension of \(x\) of length 1,2 , \(\ldots, n-1\) and greater than or equal to \(n\).
If the optional argument dim is given then operate along this dimension.
See also: [runlength], page 786.
```

count = runlength (x)
[count, value] = runlength (x)

```

Find the lengths of all sequences of common values.
count is a vector with the lengths of each repeated value.
The optional output value contains the value that was repeated in the sequence.
```

runlength ([2, 2, 0, 4, 4, 4, 0, 1, 1, 1, 1])
m

```

See also: [run_count], page 786.

\subsection*{26.4 Correlation and Regression Analysis}
```

cov (x)
cov (x,opt)
cov (x, y)
cov (x, y, opt)

```

Compute the covariance matrix.
If each row of \(x\) and \(y\) is an observation, and each column is a variable, then the ( \(i, j\) )-th entry of cov ( \(x, y\) ) is the covariance between the \(i\)-th variable in \(x\) and the \(j\)-th variable in \(y\).
\[
\sigma_{i j}=\frac{1}{N-1} \sum_{i=1}^{N}\left(x_{i}-\bar{x}\right)\left(y_{i}-\bar{y}\right)
\]
where \(\bar{x}\) and \(\bar{y}\) are the mean values of \(x\) and \(y\).
If called with one argument, compute cov ( \(x, x\) ), the covariance between the columns of \(x\).
The argument opt determines the type of normalization to use. Valid values are
\(0: \quad\) normalize with \(N-1\), provides the best unbiased estimator of the covariance [default]
1: \(\quad\) normalize with \(N\), this provides the second moment around the mean
Compatibility Note:: Octave always treats rows of \(x\) and \(y\) as multivariate random variables. For two inputs, however, matlab treats \(x\) and \(y\) as two univariate distributions regardless of their shapes, and will calculate \(\operatorname{cov}([x(:), y(:)])\) whenever the number of elements in \(x\) and \(y\) are equal. This will result in a \(2 x 2\) matrix. Code relying on matlab's definition will need to be changed when running in Octave.
See also: [corr], page 787.
corr ( \(x\) )
\(\operatorname{corr}(x, y)\)
Compute matrix of correlation coefficients.
If each row of \(x\) and \(y\) is an observation and each column is a variable, then the ( \(i, j\) )-th entry of corr ( \(x, y\) ) is the correlation between the \(i\)-th variable in \(x\) and the \(j\)-th variable in \(y\).
\[
\operatorname{corr}(x, y)=\frac{\operatorname{cov}(x, y)}{\operatorname{std}(x) \operatorname{std}(y)}
\]

If called with one argument, compute corr \((x, x)\), the correlation between the columns of \(x\).

See also: [cov], page 786.
\(r=\operatorname{corrcoef}(x)\)
\(r=\operatorname{corrcoef}\) ( \(x, y\) )
\(r=\) corrcoef (..., param, value, ...)
[r, p] = corrcoef (...)
[r, p, lci, hci] = corrcoef (...)
Compute a matrix of correlation coefficients.
\(x\) is an array where each column contains a variable and each row is an observation.
If a second input \(y\) (of the same size as \(x\) ) is given then calculate the correlation coefficients between \(x\) and \(y\).
param, value are optional pairs of parameters and values which modify the calculation. Valid options are:
"alpha" Confidence level used for the bounds of the confidence interval, lci and hci. Default is 0.05 , i.e., \(95 \%\) confidence interval.
"rows" Determine processing of NaN values. Acceptable values are "all", "complete", and "pairwise". Default is "all". With "complete", only the rows without NaN values are considered. With "pairwise", the selection of NaN-free rows is made for each pair of variables.

Output \(r\) is a matrix of Pearson's product moment correlation coefficients for each pair of variables.
Output \(p\) is a matrix of pair-wise p -values testing for the null hypothesis of a correlation coefficient of zero.
Outputs lci and hci are matrices containing, respectively, the lower and higher bounds of the \(95 \%\) confidence interval of each correlation coefficient.
See also: [corr], page 787, [cov], page 786.
spearman ( \(x\) )
spearman ( \(x, y\) )
Compute Spearman's rank correlation coefficient \(\rho\).
For two data vectors \(x\) and \(y\), Spearman's \(\rho\) is the correlation coefficient of the ranks of \(x\) and \(y\).
If \(x\) and \(y\) are drawn from independent distributions, \(\rho\) has zero mean and variance \(1 /(N-1)\), where \(N\) is the length of the x and y vectors, and is asymptotically normally distributed.
spearman ( \(x\) ) is equivalent to spearman ( \(x, x\) ).
See also: [ranks], page 786, [kendall], page 788.
kendall ( x )
kendall (x, y)
Compute Kendall's \(\tau\).
For two data vectors \(x, y\) of common length \(N\), Kendall's \(\tau\) is the correlation of the signs of all rank differences of \(x\) and \(y\); i.e., if both \(x\) and \(y\) have distinct entries, then
\[
\tau=\frac{1}{N(N-1)} \sum_{i, j} \operatorname{sign}\left(q_{i}-q_{j}\right) \operatorname{sign}\left(r_{i}-r_{j}\right)
\]
in which the \(q_{i}\) and \(r_{i}\) are the ranks of \(x\) and \(y\), respectively.
If \(x\) and \(y\) are drawn from independent distributions, Kendall's \(\tau\) is asymptotically normal with mean 0 and variance \(\frac{2(2 N+5)}{9 N(N-1)}\).
kendall ( \(x\) ) is equivalent to kendall ( \(x, x\) ).
See also: [ranks], page 786, [spearman], page 788.

\subsection*{26.5 Distributions}

Octave has functions for computing the Probability Density Function (PDF), the Cumulative Distribution function (CDF), and the quantile (the inverse of the CDF) for arbitrary user-defined distributions (discrete) and for experimental data (empirical).

The following table summarizes the supported distributions (in alphabetical order).
\begin{tabular}{|l|l|l|l|}
\hline Distribution & PDF & CDF & Quantile \\
\hline Univariate Discrete & discrete_pdf & discrete_cdf & discrete_inv \\
Empirical & empirical_pdf & empirical_cdf & empirical_inv \\
\hline
\end{tabular}
discrete_pdf (x, v, p)
For each element of \(x\), compute the probability density function (PDF) at \(x\) of a univariate discrete distribution which assumes the values in \(v\) with probabilities \(p\).
discrete_cdf (x, v, p)
For each element of \(x\), compute the cumulative distribution function (CDF) at \(x\) of a univariate discrete distribution which assumes the values in \(v\) with probabilities \(p\).
discrete_inv ( \(x, v, p\) )
For each element of \(x\), compute the quantile (the inverse of the CDF) at \(x\) of the univariate distribution which assumes the values in \(v\) with probabilities \(p\).
empirical_pdf (x, data)
For each element of \(x\), compute the probability density function (PDF) at \(x\) of the empirical distribution obtained from the univariate sample data.
empirical_cdf (x, data)
For each element of \(x\), compute the cumulative distribution function (CDF) at \(x\) of the empirical distribution obtained from the univariate sample data.
empirical_inv (x, data)
For each element of \(x\), compute the quantile (the inverse of the CDF) at \(x\) of the empirical distribution obtained from the univariate sample data.

\subsection*{26.6 Random Number Generation}

Octave can generate random numbers from a large number of distributions. The random number generators are based on the random number generators described in Section 16.3 [Special Utility Matrices], page 525.

The following table summarizes the available random number generators (in alphabetical order).
\begin{tabular}{|l|l|}
\hline Distribution & Function \\
\hline Univariate Discrete Distribution & discrete_rnd \\
Empirical Distribution & empirical_rnd \\
Exponential Distribution & rande \\
Gamma Distribution & randg \\
Poisson Distribution & randp \\
Standard Normal Distribution & randn \\
Uniform Distribution & rand \\
Uniform Distribution (integers) & randi \\
\hline
\end{tabular}
```

discrete_rnd (v, p)
discrete_rnd (v, p,r)
discrete_rnd (v, p,r, c,...)
discrete_rnd (v, p, [sz])

```

Return a matrix of random samples from the univariate distribution which assumes the values in \(v\) with probabilities \(p\).

When called with a single size argument, return a square matrix with the dimension specified. When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions. The size may also be specified with a vector of dimensions sz . If no size arguments are given then the result matrix is the common size of \(v\) and \(p\).
```

empirical_rnd (data)
empirical_rnd (data, r)
empirical_rnd (data, r, c, ...)
empirical_rnd (data, [sz])

```

Return a matrix of random samples from the empirical distribution obtained from the univariate sample data.
When called with a single size argument, return a square matrix with the dimension specified. When called with more than one scalar argument the first two arguments are taken as the number of rows and columns and any further arguments specify additional matrix dimensions. The size may also be specified with a vector of dimensions sz.
If no size arguments are given then the result matrix is a random ordering of the sample data.

\section*{27 Sets}

Octave has a number of functions for managing sets of data. A set is defined as a collection of unique elements and is typically represented by a vector of numbers sorted in ascending order. Any vector or matrix can be converted to a set by removing duplicates through the use of the unique function. However, it isn't necessary to explicitly create a set as all of the functions which operate on sets will convert their input to a set before proceeding.
```

unique (x)
unique (x, "rows")
unique (..., "sorted")
unique (..., "stable")
[y, i, j] = unique (...)
[y, i, j] = unique (..., "first")
[y, i, j] = unique (..., "last")
[y, i, j] = unique (..., "legacy")

```

Return the unique elements of \(x\).
If the input \(x\) is a column vector then return a column vector; Otherwise, return a row vector. \(x\) may also be a cell array of strings.
If the optional argument "rows" is given then return the unique rows of \(x\). The input must be a 2-D numeric matrix to use this option.
The optional argument "sorted"/"stable" controls the order in which unique values appear in the output. The default is "sorted" and values in the output are placed in ascending order. The alternative "stable" preserves the order found in the input \(x\). If requested, return column index vectors \(i\) and \(j\) such that \(y=x(i)\) and \(x=y(j)\).
Additionally, if \(i\) is a requested output then one of the flags "first" or "last" may be given. If "last" is specified, return the highest possible indices in \(i\), otherwise, if "first" is specified, return the lowest. The default is "first".
Example 1: sort order
```

unique ([3, 1, 1, 2])

# [1, 2, 3]

unique ([3, 1, 1, 2], "stable")

# [3, 1, 2]

```

Example 2 : index selection
```

[~, i] = unique ([3, 1, 1, 2], "first")

# i = [2; 4; 1]

[~, i] = unique ([3, 1, 1, 2], "last")

# i = [3; 4; 1]

```

Programming Notes: The input flag "legacy" changes the algorithm to be compatible with matlab releases prior to R2012b. Specifically, The index ordering flag is changed to "last", and the shape of the outputs \(i, j\) will follow the shape of the input \(x\) rather than always being column vectors.
The third output, \(j\), has not been implemented yet when the sort order is "stable".
See also: [union], page 793, [intersect], page 793, [setdiff], page 794, [setxor], page 794, [ismember], page 795.
\(c=\) uniquetol (A)
\(c=\) uniquetol ( \(A\), tol)
c = uniquetol (..., property, value)
[c, ia, ic] = uniquetol (...)
Return the unique elements of \(A\) within tolerance tol.
Two values, \(x\) and \(y\), are within relative tolerance if abs \((x-y)<=t o l * \max\) (abs (A(:))).
The input \(A\) must be a floating point type (double or single).
If tol is unspecified, the default tolerance is \(1 \mathrm{e}-12\) for double precision input or \(1 \mathrm{e}-6\) for single precision input.
The function may also be called with the following optional property/value pairs. Property/value pairs must be passed after other input arguments:
"ByRows" (default: false)
When true, return the unique rows of \(A\). A must be a 2-D array to use this option. For rows, the criteria for uniqueness is changed to all (abs \((x-y)<=t o l * \max (\operatorname{abs}(A),[], 1)\) ) which compares each column component of a row against a column-specific tolerance.
"DataScale"
The tolerance test is changed to abs \((x-y)<=t o l * D S\) where \(D S\) is a scalar unless the property "ByRows" is true. In that case, \(D S\) can either be a scalar or a vector with a length equal to the number of columns in \(A\). Using a value of 1.0 for \(D S\) will change the tolerance from a relative one to an absolute tolerance. Using a value of Inf will disable testing.
"OutputAllIndices" (default: false)
When true, ia is a cell array (not a vector) that contains the indices for all elements in \(A\) that are within tolerance of a value in \(C\). That is, each cell in ia corresponds to a single unique value in \(C\), and the values in each cell correspond to locations in \(A\).

The output \(c\) is a row vector if the input \(A\) is a row vector. For all other cases, a column vector is returned.
The optional output ia is a column index vector such that \(c=A(i a)\). If the "ByRows" property is true, the condition is \(c=A(i a,:)\). If the "OutputAllIndices" property is true, then the values \(A(i a\{i\})\) are all within tolerance of the unique value \(c(i)\).
The optional output ic is a column index vector such that \(A=c(i c)\) when \(A\) is a vector. When \(A\) is a matrix, \(A(:)=c(i c)\). If the "ByRows" property is true then \(A\) \(=c(i c,:)\).
Example: small round-off errors require uniquetol, not unique
```

x = [1:5];

## Inverse_Function (Function (x)) should return exactly x

y = exp (log (x));
D = unique ([x, y])

# [llllllllll

C = uniquetol ([x, y])
=>[$$
\begin{array}{llllll}{1}&{2}&{3}&{4}&{5]}\end{array}
$$]

```

See also: [unique], page 791, [union], page 793, [intersect], page 793, [setdiff], page 794, [setxor], page 794, [ismember], page 795.

\subsection*{27.1 Set Operations}

Octave supports several basic set operations. Octave can compute the union, intersection, and difference of two sets. Octave also supports the Exclusive Or set operation.

The functions for set operations all work in the same way by accepting two input sets and returning a third set. As an example, assume that a and b contains two sets, then union ( \(\mathrm{a}, \mathrm{b}\) )
computes the union of the two sets.
Finally, determining whether elements belong to a set can be done with the ismember function. Because sets are ordered this operation is very efficient and is of order \(\mathrm{O}(\log 2(\mathrm{n}))\) which is preferable to the find function which is of order \(\mathrm{O}(\mathrm{n})\).
```

c = intersect (a, b)
c = intersect (a, b, "rows")
c = intersect (..., "sorted")
c = intersect (..., "stable")
c = intersect (..., "legacy")
[c, ia, ib] = intersect (...)

```
    Return the unique elements common to both \(a\) and \(b\).

If \(a\) and \(b\) are both row vectors then return a row vector; Otherwise, return a column vector. The inputs may also be cell arrays of strings.
If the optional input "rows" is given then return the common rows of \(a\) and \(b\). The inputs must be 2-D numeric matrices to use this option.
The optional argument "sorted"/"stable" controls the order in which unique values appear in the output. The default is "sorted" and values in the output are placed in ascending order. The alternative "stable" preserves the order found in the input. If requested, return column index vectors ia and ib such that \(c=a(i a)\) and \(c=\) b(ib).
Programming Note: The input flag "legacy" changes the algorithm to be compatible with matlab releases prior to R2012b.
See also: [unique], page 791, [union], page 793, [setdiff], page 794, [setxor], page 794, [ismember], page 795.
\(c=\) union \((a, b)\)
\(c=\) union ( \(a, b\), "rows")
\(c=\) union (..., "sorted")
c = union (..., "stable")
c = union (..., "legacy")
[c, ia, ib] = union (...)
Return the unique elements that are in either a or \(b\).
If \(a\) and \(b\) are both row vectors then return a row vector; Otherwise, return a column vector. The inputs may also be cell arrays of strings.

If the optional input "rows" is given then return rows that are in either a or \(b\). The inputs must be 2-D numeric matrices to use this option.
The optional argument "sorted"/"stable" controls the order in which unique values appear in the output. The default is "sorted" and values in the output are placed in ascending order. The alternative "stable" preserves the order found in the input. The optional outputs ia and ib are column index vectors such that \(a(i a)\) and \(b(i b)\) are disjoint sets whose union is \(c\).
Programming Note: The input flag "legacy" changes the algorithm to be compatible with matlab releases prior to R2012b.
See also: [unique], page 791, [intersect], page 793, [setdiff], page 794, [setxor], page 794, [ismember], page 795.
\(c=\operatorname{setdiff}(a, b)\)
\(c=\operatorname{setdiff}(a, b\), "rows")
\(c=\) setdiff (..., "sorted")
c = setdiff (..., "stable")
\(c=\) setdiff (..., "legacy")
[c, ia] = setdiff (...)
Return the unique elements in a that are not in \(b\).
If \(a\) is a row vector return a row vector; Otherwise, return a column vector. The inputs may also be cell arrays of strings.
If the optional input "rows" is given then return the rows in a that are not in \(b\). The inputs must be 2-D numeric matrices to use this option.
The optional argument "sorted"/"stable" controls the order in which unique values appear in the output. The default is "sorted" and values in the output are placed in ascending order. The alternative "stable" preserves the order found in the input.
If requested, return the index vector ia such that \(c=a(i a)\).
Programming Note: The input flag "legacy" changes the algorithm to be compatible with matlab releases prior to R2012b.

See also: [unique], page 791, [union], page 793, [intersect], page 793, [setxor], page 794, [ismember], page 795.
\(c=\operatorname{setxor}(a, b)\)
\(c=\operatorname{setxor}(a, b\), rows")
\(c=\) setxor (..., "sorted")
c \(=\) setxor (..., "stable")
c = setxor (..., "legacy")
[c, ia, ib] = setxor (...)
Return the unique elements exclusive to sets \(a\) or \(b\).
If \(a\) and \(b\) are both row vectors then return a row vector; Otherwise, return a column vector. The inputs may also be cell arrays of strings.
If the optional input "rows" is given then return the rows exclusive to sets \(a\) and \(b\). The inputs must be 2-D numeric matrices to use this option.

The optional argument "sorted"/"stable" controls the order in which unique values appear in the output. The default is "sorted" and values in the output are placed in ascending order. The alternative "stable" preserves the order found in the input. The optional outputs ia and ib are column index vectors such that \(a(i a)\) and \(b(i b)\) are disjoint sets whose union is \(c\).
Programming Note: The input flag "legacy" changes the algorithm to be compatible with matlab releases prior to R2012b.
See also: [unique], page 791, [union], page 793, [intersect], page 793, [setdiff], page 794, [ismember], page 795.
\(t f=\) ismember ( \(a, s\) )
\(t f=\) ismember (a, s, "rows")
[tf, s_idx] = ismember (...)
Return a logical matrix \(t f\) with the same shape as a which is true (1) if the element in \(a\) is found in \(s\) and false ( 0 ) if it is not.
If a second output argument is requested then the index into \(s\) of each matching element is also returned.
```

a = [3, 10, 1];
s = [0:9];
[tf, s_idx] = ismember (a, s)
t tf = [1, 0, 1]
\# s_idx = [4, 0, 2]

```

The inputs a and \(s\) may also be cell arrays.
```

a = {"abc"};
s = {"abc", "def"};
[tf, s_idx] = ismember (a, s)
tff=1
| s_idx = 1

```

If the optional third argument "rows" is given then compare rows in a with rows in s. The inputs must be 2-D matrices with the same number of columns to use this option.
```

a = [1:3; 5:7; 4:6];
s = [0:2; 1:3; 2:4; 3:5; 4:6];
[tf, s_idx] = ismember (a, s, "rows")
tf = logical ([1; 0; 1])
| s_idx = [2; 0; 5];

```

See also: [lookup], page 514, [unique], page 791, [union], page 793, [intersect], page 793, [setdiff], page 794, [setxor], page 794.
powerset (a)
powerset (a, "rows")
Compute the powerset (all subsets) of the set a.
The set a must be a numerical matrix or a cell array of strings. The output will always be a cell array of either vectors or strings.

With the optional argument "rows", each row of the set a is considered one element of the set. The input must be a 2-D numeric matrix to use this argument.

See also: [unique], page 791, [union], page 793, [intersect], page 793, [setdiff], page 794, [setxor], page 794, [ismember], page 795.

\section*{28 Polynomial Manipulations}

In Octave, a polynomial is represented by its coefficients (arranged in descending order). For example, a vector \(c\) of length \(N+1\) corresponds to the following polynomial of order \(N\)
\[
p(x)=c_{1} x^{N}+\ldots+c_{N} x+c_{N+1}
\]

\subsection*{28.1 Evaluating Polynomials}

The value of a polynomial represented by the vector \(c\) can be evaluated at the point \(x\) very easily, as the following example shows:
```

N = length (c) - 1;
val = dot (x.^(N:-1:0), c);

```

While the above example shows how easy it is to compute the value of a polynomial, it isn't the most stable algorithm. With larger polynomials you should use more elegant algorithms, such as Horner's Method, which is exactly what the Octave function polyval does.

In the case where \(x\) is a square matrix, the polynomial given by \(c\) is still well-defined. As when \(x\) is a scalar the obvious implementation is easily expressed in Octave, but also in this case more elegant algorithms perform better. The polyvalm function provides such an algorithm.
\(y=\operatorname{polyval}(p, x)\)
\(y=\operatorname{polyval}(p, x,[], m u)\)
\([y, d y]=\) polyval \((p, x, s)\)
\([y, d y]=\) polyval \((p, x, s, m u)\)
Evaluate the polynomial \(p\) at the specified values of \(x\).
If \(x\) is a vector or matrix, the polynomial is evaluated for each of the elements of \(x\).
When \(m u\) is present, evaluate the polynomial for \((x-m u(1)) / m u(2)\).
In addition to evaluating the polynomial, the second output represents the prediction interval, \(y+/-d y\), which contains at least \(50 \%\) of the future predictions. To calculate the prediction interval, the structured variable \(s\), originating from polyfit, must be supplied.

See also: [polyvalm], page 797, [polyaffine], page 803, [polyfit], page 803, [roots], page 798, [poly], page 813.
polyvalm ( \(c, x\) )
Evaluate a polynomial in the matrix sense.
polyvalm ( \(c, x\) ) will evaluate the polynomial in the matrix sense, i.e., matrix multiplication is used instead of element by element multiplication as used in polyval.

The argument \(x\) must be a square matrix.
See also: [polyval], page 797, [roots], page 798, [poly], page 813.

\subsection*{28.2 Finding Roots}

Octave can find the roots of a given polynomial. This is done by computing the companion matrix of the polynomial (see the compan function for a definition), and then finding its eigenvalues.
roots (c)
Compute the roots of the polynomial \(c\).
For a vector \(c\) with \(N\) components, return the roots of the polynomial
\[
c_{1} x^{N-1}+\cdots+c_{N-1} x+c_{N} .
\]

As an example, the following code finds the roots of the quadratic polynomial
\[
p(x)=x^{2}-5 .
\]
```

c = [1, 0, -5];
roots (c)
=> 2.2361
=> -2.2361

```

Note that the true result is \(\pm \sqrt{5}\) which is roughly \(\pm 2.2361\).
See also: [poly], page 813, [compan], page 798, [fzero], page 644.
\(z=\) polyeig (CO, C1, ..., Cl)
[v, z] = polyeig (CO, C1, ..., Cl)
Solve the polynomial eigenvalue problem of degree \(l\).
Given an nxn matrix polynomial
\(C(s)=C 0+C 1 s+\ldots+C l s^{\wedge} 1\)
polyeig solves the eigenvalue problem
\(\left(C 0+C 1 z+\ldots+C l z^{\wedge} 1\right) v=0\).
Note that the eigenvalues \(z\) are the zeros of the matrix polynomial. \(z\) is a row vector with \(n * l\) elements. \(v\) is a matrix ( \(n \times n^{*} l\) ) with columns that correspond to the eigenvectors.
See also: [eig], page 587, [eigs], page 686, [compan], page 798.
compan (c)
Compute the companion matrix corresponding to polynomial coefficient vector \(c\).
The companion matrix is
\[
A=\left[\begin{array}{ccccc}
-c_{2} / c_{1} & -c_{3} / c_{1} & \cdots & -c_{N} / c_{1} & -c_{N+1} / c_{1} \\
1 & 0 & \cdots & 0 & 0 \\
0 & 1 & \cdots & 0 & 0 \\
\vdots & \vdots & \ddots & \vdots & \vdots \\
0 & 0 & \cdots & 1 & 0
\end{array}\right]
\]

The eigenvalues of the companion matrix are equal to the roots of the polynomial.
See also: [roots], page 798, [poly], page 813, [eig], page 587.
```

[multp, idxp] = mpoles (p)

```
[multp, idxp] = mpoles ( \(p\), tol)
[multp, idxp] = mpoles ( \(p\), tol, reorder)

Identify unique poles in \(p\) and their associated multiplicity.
The output is ordered from pole with largest magnitude to smallest magnitude.
If the relative difference of two poles is less than tol then they are considered to be multiples. The default value for \(t o l\) is 0.001 .
If the optional parameter reorder is zero, poles are not sorted.
The output multp is a vector specifying the multiplicity of the poles. multp(n) refers to the multiplicity of the Nth pole \(p(i d x p(n))\).
For example:
```

p = [2 3 1 1 1 1 2];
[m, n] = mpoles (p)
m = [1; 1; 2; 1; 2]
\# n = [2; 5; 1; 4; 3]
mp(n) = [3, 2, 2, 1, 1]

```

See also: [residue], page 801, [poly], page 813, [roots], page 798, [conv], page 799, [deconv], page 800.

\subsection*{28.3 Products of Polynomials}
```

conv (a, b)
conv (a, b, shape)

```

Convolve two vectors \(a\) and \(b\).
When \(a\) and \(b\) are the coefficient vectors of two polynomials, the convolution represents the coefficient vector of the product polynomial.

The size of the result is determined by the optional shape argument which takes the following values
shape \(=\) "full"
Return the full convolution. (default) The result is a vector with length equal to length (a) + length (b) -1 .
shape \(=\) "same"
Return the central part of the convolution with the same size as a.
shape \(=\) "valid"
Return only the parts which do not include zero-padded edges. The size of the result is max (size (a) - size (b) \(+1,0\) ).

See also: [deconv], page 800, [conv2], page 800, [convn], page 799, [fftconv], page 848.
\(C=\operatorname{convn}(A, B)\)
\(C=\) convn (A, B, shape)
Return the n-D convolution of \(A\) and \(B\).

The size of the result is determined by the optional shape argument which takes the following values
shape \(=\) "full"
Return the full convolution. (default)
shape \(=\) "same"
Return central part of the convolution with the same size as \(A\). The central part of the convolution begins at the indices floor ([size(B)/2] \(+1)\).
shape \(=\) "valid"
Return only the parts which do not include zero-padded edges. The size of the result is max (size (A) - size (B) \(+1,0\) ).

See also: [conv2], page 800, [conv], page 799.
\(b=\operatorname{deconv}(y, a)\)
\([b, r]=\operatorname{deconv}(y, a)\)
Deconvolve two vectors (polynomial division).
\([b, r]=\operatorname{deconv}(y, a)\) solves for \(b\) and \(r\) such that \(y=\operatorname{conv}(a, b)+r\).
If \(y\) and a are polynomial coefficient vectors, \(b\) will contain the coefficients of the polynomial quotient and \(r\) will be a remainder polynomial of lowest order.
See also: [conv], page 799, [residue], page 801.
```

conv2 (A, B)
conv2 (v1, v2, m)
conv2 (..., shape)

```

Return the 2-D convolution of \(A\) and \(B\).
The size of the result is determined by the optional shape argument which takes the following values
shape \(=\) "full"
Return the full convolution. (default)
shape \(=\) "same"
Return the central part of the convolution with the same size as \(A\). The central part of the convolution begins at the indices floor ([size(B)/2] \(+1)\).
shape \(=\) "valid"
Return only the parts which do not include zero-padded edges. The size of the result is max (size (A) - size (B) \(+1,0\) ).

When the third argument is a matrix, return the convolution of the matrix \(m\) by the vector v1 in the column direction and by the vector v2 in the row direction.

See also: [conv], page 799, [convn], page 799.
\(q=\operatorname{polygcd}(b, a)\)
\(q=\operatorname{polygcd}(b, a, t o l)\)
Find the greatest common divisor of two polynomials.

This is equivalent to the polynomial found by multiplying together all the common roots. Together with deconv, you can reduce a ratio of two polynomials.
The tolerance tol defaults to sqrt (eps).
Caution: This is a numerically unstable algorithm and should not be used on large polynomials.
Example code:
```

polygcd (poly (1:8), poly (3:12)) - poly (3:8)

# [ 0, 0, 0, 0, 0, 0, 0 ]

deconv (poly (1:8), polygcd (poly (1:8), poly (3:12))) - poly (1:2)
=> [ 0, 0, 0 ]

```

See also: [poly], page 813, [roots], page 798, [conv], page 799, [deconv], page 800, [residue], page 801.
\([r, p, k, e]=\) residue \((b, a)\)
[b, a] = residue ( \(r, p, k\) )
\([b, a]=\operatorname{residue}(r, p, k, e)\)
The first calling form computes the partial fraction expansion for the quotient of the polynomials, \(b\) and \(a\).
The quotient is defined as
\[
\frac{B(s)}{A(s)}=\sum_{m=1}^{M} \frac{r_{m}}{\left(s-p_{m}\right)_{m}^{e}}+\sum_{i=1}^{N} k_{i} s^{N-i} .
\]
where \(M\) is the number of poles (the length of the \(r, p\), and \(e\) ), the \(k\) vector is a polynomial of order \(N-1\) representing the direct contribution, and the e vector specifies the multiplicity of the m-th residue's pole.
For example,
\[
\begin{aligned}
& \mathrm{b}=[1,1,1] ; \\
& \mathrm{a}=[1,-5,8,-4] ; \\
& {[\mathrm{r}, \mathrm{p}, \mathrm{k}, \mathrm{e}]=\text { residue }(\mathrm{b}, \mathrm{a})} \\
& \quad \Rightarrow \mathrm{r}=[-2 ; 7 ; 3] \\
& \\
& \quad \Rightarrow \mathrm{p}=[2 ; 2 ; 1] \\
& \\
& \\
& \\
&
\end{aligned} \Rightarrow \mathrm{k}=[](0 \mathrm{x} 0) \mathrm{e}=[1 ; 2 ; 1] .
\]
which represents the following partial fraction expansion
\[
\frac{s^{2}+s+1}{s^{3}-5 s^{2}+8 s-4}=\frac{-2}{s-2}+\frac{7}{(s-2)^{2}}+\frac{3}{s-1}
\]

The second calling form performs the inverse operation and computes the reconstituted quotient of polynomials, \(b(\mathrm{~s}) / a(\mathrm{~s})\), from the partial fraction expansion; represented by the residues, poles, and a direct polynomial specified by \(r, p\) and \(k\), and the pole multiplicity \(e\).
If the multiplicity, \(e\), is not explicitly specified the multiplicity is determined by the function mpoles.

For example:
\[
\begin{aligned}
& \mathrm{r}=[-2 ; 7 ; 3] ; \\
& \mathrm{p}=[2 ; 2 ; 1] ; \\
& \mathrm{k}=[1,0] ; \\
& {[\mathrm{b}, \mathrm{a}]=\text { residue }(\mathrm{r}, \mathrm{p}, \mathrm{k})} \\
& \quad \Rightarrow \mathrm{b}=[1,-5,9,-3,1] \\
& \\
& \quad \Rightarrow \mathrm{a}=[1,-5,8,-4]
\end{aligned}
\]
```

where mpoles is used to determine e = [1; 2; 1]

```

Alternatively the multiplicity may be defined explicitly, for example,
\[
\begin{aligned}
& \mathrm{r}=[7 ; 3 ;-2] ; \\
& \mathrm{p}=[2 ; 1 ; 2] ; \\
& \mathrm{k}=[1,0] ; \\
& \mathrm{e}=[2 ; 1 ; 1] ; \\
& {[\mathrm{b}, \mathrm{a}]=\text { residue }(\mathrm{r}, \mathrm{p}, \mathrm{k}, \mathrm{e})} \\
& \quad \Rightarrow \mathrm{b}=[1,-5,9,-3,1] \\
& \quad \Rightarrow \mathrm{a}=[1,-5,8,-4]
\end{aligned}
\]
which represents the following partial fraction expansion
\[
\frac{-2}{s-2}+\frac{7}{(s-2)^{2}}+\frac{3}{s-1}+s=\frac{s^{4}-5 s^{3}+9 s^{2}-3 s+1}{s^{3}-5 s^{2}+8 s-4}
\]

See also: [mpoles], page 799, [poly], page 813, [roots], page 798, [conv], page 799, [deconv], page 800.

\subsection*{28.4 Derivatives / Integrals / Transforms}

Octave comes with functions for computing the derivative and the integral of a polynomial. The functions polyder and polyint both return new polynomials describing the result. As an example we'll compute the definite integral of \(p(x)=x^{2}+1\) from 0 to 3 .
```

c = [1, 0, 1];
integral = polyint (c);
area = polyval (integral, 3) - polyval (integral, 0)

# 12

```
polyder ( \(p\) )
[k] = polyder ( \(a, b\) )
\([q, d]=\operatorname{polyder}(b, a)\)

Return the coefficients of the derivative of the polynomial whose coefficients are given by the vector \(p\).

If a pair of polynomials is given, return the derivative of the product \(a * b\).
If two inputs and two outputs are given, return the derivative of the polynomial quotient \(b / a\). The quotient numerator is in \(q\) and the denominator in \(d\).
See also: [polyint], page 803, [polyval], page 797, [polyreduce], page 813.
```

polyint (p)

```
polyint ( \(p, k\) )

Return the coefficients of the integral of the polynomial whose coefficients are represented by the vector \(p\).
The variable \(k\) is the constant of integration, which by default is set to zero.
See also: [polyder], page 802, [polyval], page 797.
polyaffine (f, mu)
Return the coefficients of the polynomial vector \(f\) after an affine transformation.
If \(f\) is the vector representing the polynomial \(\mathrm{f}(\mathrm{x})\), then \(g=\) polyaffine ( \(f, \mathrm{mu}\) ) is the vector representing:
\[
g(x)=f((x-m u(1)) / m u(2))
\]

See also: [polyval], page 797, [polyfit], page 803.

\subsection*{28.5 Polynomial Interpolation}

Octave comes with good support for various kinds of interpolation, most of which are described in Chapter 29 [Interpolation], page 815. One simple alternative to the functions described in the aforementioned chapter, is to fit a single polynomial, or a piecewise polynomial (spline) to some given data points. To avoid a highly fluctuating polynomial, one most often wants to fit a low-order polynomial to data. This usually means that it is necessary to fit the polynomial in a least-squares sense, which just is what the polyfit function does.
\(p=\) polyfit (x, y, n)
[p,s] = polyfit (x, y, n)
[p, s, mu] = polyfit (x, y, n)
Return the coefficients of a polynomial \(p(x)\) of degree \(n\) that minimizes the least-squares-error of the fit to the points \([x(:), y(:)]\).
\(n\) is typically an integer \(\geq 0\) specifying the degree of the approximating polynomial. If \(n\) is a logical vector, it is used as a mask to selectively force the corresponding polynomial coefficients to be used or ignored.
The polynomial coefficients are returned in the row vector \(p\). The output \(p\) may be directly used with polyval to estimate values using the fitted polynomial.
The optional output \(s\) is a structure containing the following fields:
' \(y f\) ' The values of the polynomial for each value of \(x\).
' X ' The Vandermonde matrix used to compute the polynomial coefficients.
' R ' Triangular factor R from the QR decomposition.
' C ' The unscaled covariance matrix, formally equal to the inverse of \(x\) '* \(x\), but computed in a way minimizing roundoff error propagation.
'df' The degrees of freedom.
'normr' The norm of the residuals.

The second output may be used by polyval to calculate the statistical error limits of the predicted values. In particular, the standard deviation of \(p\) coefficients is given by
sqrt (diag (s.C)/s.df) * s.normr.
When the third output, \(m u\), is present the original data is centered and scaled which can improve the numerical stability of the fit. The coefficients \(p\) are associated with a polynomial in
```

xhat = (x-mu(1)) / mu(2)

```
where \(m u(1)=\) mean \((x)\), and \(m u(2)=\operatorname{std}(x)\).
Example 1: logical \(n\) and integer \(n\)
```

f = @(x) x.^2 + 5; \# data-generating function
x = 0:5;
y = f (x);

## Fit data to polynomial A*x^3 + B*x^1

p = polyfit (x, y, logical ([1, 0, 1, 0]))
p = [ 0.0680, 0, 4.2444, 0 ]

## Fit data to polynomial using all terms up to x^3

p = polyfit (x, y, 3)
p = [ -4.9608e-17, 1.0000e+00, -3.3813e-15, 5.0000e+00 ]

```

See also: [polyval], page 797, [polyaffine], page 803, [roots], page 798, [vander], page 544, [zscore], page 784.

In situations where a single polynomial isn't good enough, a solution is to use several polynomials pieced together. The function splinefit fits a piecewise polynomial (spline) to a set of data.
```

pp = splinefit (x, y, breaks)
pp = splinefit (x, y, p)
pp = splinefit (..., "periodic", periodic)
pp = splinefit (..., "robust", robust)
pp = splinefit (..., "beta", beta)
pp = splinefit (..., "order", order)
pp = splinefit (..., "constraints", constraints)

```

Fit a piecewise cubic spline with breaks (knots) breaks to the noisy data, \(x\) and \(y\).
\(x\) is a vector, and \(y\) is a vector or N-D array. If \(y\) is an N-D array, then \(x(j)\) is matched to \(y(:, \ldots, ., \dot{j})\).
\(p\) is a positive integer defining the number of intervals along \(x\), and \(p+1\) is the number of breaks. The number of points in each interval differ by no more than 1 .
The optional property periodic is a logical value which specifies whether a periodic boundary condition is applied to the spline. The length of the period is max (breaks) - min (breaks). The default value is false.

The optional property robust is a logical value which specifies if robust fitting is to be applied to reduce the influence of outlying data points. Three iterations of weighted least squares are performed. Weights are computed from previous residuals. The sensitivity of outlier identification is controlled by the property beta. The value of
beta is restricted to the range, \(0<\) beta \(<1\). The default value is beta \(=1 / 2\). Values close to 0 give all data equal weighting. Increasing values of beta reduce the influence of outlying data. Values close to unity may cause instability or rank deficiency.

The fitted spline is returned as a piecewise polynomial, \(p p\), and may be evaluated using ppval.

The splines are constructed of polynomials with degree order. The default is a cubic, order \(=3\). A spline with P pieces has \(\mathrm{P}+\) order degrees of freedom. With periodic boundary conditions the degrees of freedom are reduced to P .

The optional property, constraints, is a structure specifying linear constraints on the fit. The structure has three fields, "xc", "yc", and "cc".
"xc" Vector of the x-locations of the constraints.
"ус" Constraining values at the locations xc. The default is an array of zeros.
"cc" Coefficients (matrix). The default is an array of ones. The number of rows is limited to the order of the piecewise polynomials, order.

Constraints are linear combinations of derivatives of order 0 to order- 1 according to
\[
c c(1, j) \cdot y(x c(j))+c c(2, j) \cdot y \prime(x c(j))+\ldots=y c(:, \ldots,:, j) .
\]

See also: [interp1], page 815, [unmkpp], page 811, [ppval], page 812, [spline], page 818, [pchip], page 854, [ppder], page 812, [ppint], page 812, [ppjumps], page 812.

The number of breaks (or knots) used to construct the piecewise polynomial is a significant factor in suppressing the noise present in the input data, \(x\) and \(y\). This is demonstrated by the example below.
```

x = 2 * pi * rand (1, 200);
y = sin (x) + sin (2 * x) + 0.2 * randn (size (x));

## Uniform breaks

breaks = linspace (0, 2 * pi, 41); % 41 breaks, 40 pieces
pp1 = splinefit (x, y, breaks);

## Breaks interpolated from data

pp2 = splinefit (x, y, 10); % 11 breaks, 10 pieces

## Plot

xx = linspace (0, 2 * pi, 400);
y1 = ppval (pp1, xx);
y2 = ppval (pp2, xx);
plot (x, y, ".", xx, [y1; y2])
axis tight
ylim auto
legend ({"data", "41 breaks, 40 pieces", "11 breaks, 10 pieces"})

```

The result of which can be seen in Figure 28.1.


Figure 28.1: Comparison of a fitting a piecewise polynomial with 41 breaks to one with 11 breaks. The fit with the large number of breaks exhibits a fast ripple that is not present in the underlying function.

The piecewise polynomial fit, provided by splinefit, has continuous derivatives up to the order-1. For example, a cubic fit has continuous first and second derivatives. This is demonstrated by the code
```


## Data (200 points)

x = 2 * pi * rand (1, 200);
y = sin (x) + sin (2 * x) + 0.1 * randn (size (x));

## Piecewise constant

pp1 = splinefit (x, y, 8, "order", 0);

## Piecewise linear

pp2 = splinefit (x, y, 8, "order", 1);

## Piecewise quadratic

pp3 = splinefit (x, y, 8, "order", 2);

## Piecewise cubic

pp4 = splinefit (x, y, 8, "order", 3);

## Piecewise quartic

pp5 = splinefit (x, y, 8, "order", 4);

## Plot

xx = linspace (0, 2 * pi, 400);
y1 = ppval (pp1, xx);
y2 = ppval (pp2, xx);
y3 = ppval (pp3, xx);
y4 = ppval (pp4, xx);
y5 = ppval (pp5, xx);
plot (x, y, ".", xx, [y1; y2; y3; y4; y5])
axis tight

```
```

ylim auto
legend ({"data", "order 0", "order 1", "order 2", "order 3", "order 4"})

```

The result of which can be seen in Figure 28.2.


Figure 28.2: Comparison of a piecewise constant, linear, quadratic, cubic, and quartic polynomials with 8 breaks to noisy data. The higher order solutions more accurately represent the underlying function, but come with the expense of computational complexity.

When the underlying function to provide a fit to is periodic, splinefit is able to apply the boundary conditions needed to manifest a periodic fit. This is demonstrated by the code below.
```


## Data (100 points)

x = 2 * pi * [0, (rand (1, 98)), 1];
y = sin (x) - cos (2 * x) + 0.2 * randn (size (x));

## No constraints

pp1 = splinefit (x, y, 10, "order", 5);

## Periodic boundaries

pp2 = splinefit (x, y, 10, "order", 5, "periodic", true);

## Plot

xx = linspace (0, 2 * pi, 400);
y1 = ppval (pp1, xx);
y2 = ppval (pp2, xx);
plot (x, y, ".", xx, [y1; y2])
axis tight
ylim auto
legend ({"data", "no constraints", "periodic"})

```

The result of which can be seen in Figure 28.3.


Figure 28.3: Comparison of piecewise polynomial fits to a noisy periodic function with, and without, periodic boundary conditions.

More complex constraints may be added as well. For example, the code below illustrates a periodic fit with values that have been clamped at the endpoints, and a second periodic fit which is hinged at the endpoints.
```


## Data (200 points)

x = 2 * pi * rand (1, 200);
y = sin (2 * x) + 0.1 * randn (size (x));

## Breaks

breaks = linspace (0, 2 * pi, 10);

## Clamped endpoints, y = y' = 0

xc = [0, 0, 2*pi, 2*pi];
cc = [(eye (2)), (eye (2))];
con = struct ("xc", xc, "cc", cc);
pp1 = splinefit (x, y, breaks, "constraints", con);

## Hinged periodic endpoints, y = 0

con = struct ("xc", 0);
pp2 = splinefit (x, y, breaks, "constraints", con, "periodic", true);

## Plot

xx = linspace (0, 2 * pi, 400);
y1 = ppval (pp1, xx);
y2 = ppval (pp2, xx);
plot (x, y, ".", xx, [y1; y2])
axis tight
ylim auto
legend ({"data", "clamped", "hinged periodic"})

```

The result of which can be seen in Figure 28.4.


Figure 28.4: Comparison of two periodic piecewise cubic fits to a noisy periodic signal. One fit has its endpoints clamped and the second has its endpoints hinged.

The splinefit function also provides the convenience of a robust fitting, where the effect of outlying data is reduced. In the example below, three different fits are provided. Two with differing levels of outlier suppression and a third illustrating the non-robust solution.
```


## Data

x = linspace (0, 2*pi, 200);
y = sin (x) + sin (2 * x) + 0.05 * randn (size (x));

## Add outliers

x = [x, linspace(0,2*pi,60)];
y = [y, -ones(1,60)];

## Fit splines with hinged conditions

con = struct ("xc", [0, 2*pi]);

## Robust fitting, beta = 0.25

pp1 = splinefit (x, y, 8, "constraints", con, "beta", 0.25);

## Robust fitting, beta = 0.75

pp2 = splinefit (x, y, 8, "constraints", con, "beta", 0.75);

## No robust fitting

pp3 = splinefit (x, y, 8, "constraints", con);

## Plot

xx = linspace (0, 2*pi, 400);
y1 = ppval (pp1, xx);
y2 = ppval (pp2, xx);
y3 = ppval (pp3, xx);
plot (x, y, ".", xx, [y1; y2; y3])
legend ({"data with outliers","robust, beta = 0.25", ...
"robust, beta = 0.75", "no robust fitting"})
axis tight

```
ylim auto
The result of which can be seen in Figure 28.5.


Figure 28.5: Comparison of two different levels of robust fitting (beta \(=0.25\) and 0.75 ) to noisy data combined with outlying data. A conventional fit, without robust fitting (beta \(=0\) ) is also included.

A very specific form of polynomial interpretation is the Padé approximant. For control systems, a continuous-time delay can be modeled very simply with the approximant.
[num, den] = padecoef \((T)\)
[num, den] = padecoef ( \(T, N\) )
Compute the Nth-order Padé approximant of the continuous-time delay \(T\) in transfer function form.
The Padé approximant of \(e^{-s T}\) is defined by the following equation
\[
e^{-s T} \approx \frac{P_{n}(s)}{Q_{n}(s)}
\]
where both \(P_{n}(s)\) and \(Q_{n}(s)\) are \(N^{t h}\)-order rational functions defined by the following expressions
\[
\begin{gathered}
P_{n}(s)=\sum_{k=0}^{N} \frac{(2 N-k)!N!}{(2 N)!k!(N-k)!}(-s T)^{k} \\
Q_{n}(s)=P_{n}(-s)
\end{gathered}
\]

The inputs \(T\) and \(N\) must be non-negative numeric scalars. If \(N\) is unspecified it defaults to 1 .
The output row vectors num and den contain the numerator and denominator coefficients in descending powers of s. Both are Nth-order polynomials.

For example:
```

t = 0.1;
n = 4;
[num, den] = padecoef (t, n)

# num =

    1.0000e-04 -2.0000e-02 1.8000e+00 -8.4000e+01 1.6800e+03
    | den =
1.0000e-04 2.0000e-02 1.8000e+00 8.4000e+01 1.6800e+03

```

The function, ppval, evaluates the piecewise polynomials, created by mkpp or other means, and unmkpp returns detailed information about the piecewise polynomial.

The following example shows how to combine two linear functions and a quadratic into one function. Each of these functions is expressed on adjoined intervals.
```

    x = [-2, -1, 1, 2];
    p = [ 0, 1, 0;
        1, -2, 1;
    0, -1, 1];
    pp = mkpp (x, p);
    xi = linspace (-2, 2, 50);
    yi = ppval (pp, xi);
    plot (xi, yi);
    pp = mkpp (breaks, coefs)
pp = mkpp (breaks, coefs,d)

```

Construct a piecewise polynomial (pp) structure from sample points breaks and coefficients coefs.
breaks must be a vector of strictly increasing values. The number of intervals is given by \(n i=\) length (breaks) -1 .
When \(m\) is the polynomial order coefs must be of size: ni-by- \((m+1)\).
The i-th row of coefs, coefs(i,:), contains the coefficients for the polynomial over the \(i\)-th interval, ordered from highest \((m)\) to lowest ( 0 ) degree.
coefs may also be a multi-dimensional array, specifying a vector-valued or array-valued polynomial. In that case the polynomial order \(m\) is defined by the length of the last dimension of coefs. The size of first dimension(s) are given by the scalar or vector \(d\). If \(d\) is not given it is set to 1 . In this case \(p(r, i,:)\) contains the coefficients for the \(r\)-th polynomial defined on interval \(i\). In any case coefs is reshaped to a 2-D matrix of size \([n i * \operatorname{prod}(d) \mathrm{m}]\).
Programming Note: ppval evaluates polynomials at xi - breaks(i), i.e., it subtracts the lower endpoint of the current interval from xi. This must be taken into account when creating piecewise polynomials objects with mkpp.

See also: [unmkpp], page 811, [ppval], page 812, [spline], page 818, [pchip], page 854, [ppder], page 812, [ppint], page 812, [ppjumps], page 812.
\([x, p, n, k, d]=\) unmkpp \((p p)\)
Extract the components of a piecewise polynomial structure pp.

This function is the inverse of mkpp: it extracts the inputs to mkpp needed to create the piecewise polynomial structure pp. The code below makes this relation explicit:
```

[breaks, coefs, numinter, order, dim] = unmkpp (pp);
pp2 = mkpp (breaks, coefs, dim);

```

The piecewise polynomial structure pp2 obtained in this way, is identical to the original pp . The same can be obtained by directly accessing the fields of the structure pp.
The components are:
\(x \quad\) Sample points or breaks.
\(p \quad\) Polynomial coefficients for points in sample interval. \(p(i,:)\) contains the coefficients for the polynomial over interval \(i\) ordered from highest to lowest degree. If \(d>1\), then \(p\) is a matrix of size \([n * \operatorname{prod}(d) m\), where the \(i+(1: d)\) rows are the coefficients of all the \(d\) polynomials in the interval \(i\).
\(n \quad\) Number of polynomial pieces or intervals, \(n=\) length ( \(x\) ) - 1 .
\(k \quad\) Order of the polynomial plus 1.
d Number of polynomials defined for each interval.
See also: [mkpp], page 811, [ppval], page 812, [spline], page 818, [pchip], page 854.
yi = ppval (pp, xi)
Evaluate the piecewise polynomial structure \(p p\) at the points xi.
If \(p p\) describes a scalar polynomial function, the result is an array of the same shape as xi. Otherwise, the size of the result is [pp.dim, length (xi)] if xi is a vector, or [pp.dim, \(\operatorname{size}(x i)]\) if it is a multi-dimensional array.
See also: [mkpp], page 811, [unmkpp], page 811, [spline], page 818, [pchip], page 854.
ppd \(=\) ppder ( \(p p\) )
ppd \(=\) ppder ( \(p p, m\) )
Compute the piecewise m-th derivative of a piecewise polynomial struct \(p p\).
If \(m\) is omitted the first derivative is calculated.
See also: [mkpp], page 811, [ppval], page 812, [ppint], page 812.
ppi = ppint (pp)
ppi \(=\) ppint ( \(p\) p, c)
Compute the integral of the piecewise polynomial struct \(p p\).
\(c\), if given, is the constant of integration.
See also: [mkpp], page 811, [ppval], page 812, [ppder], page 812.
jumps = ppjumps (pp)
Evaluate the boundary jumps of a piecewise polynomial.
If there are \(n\) intervals, and the dimensionality of \(p p\) is \(d\), the resulting array has dimensions [d, \(\mathrm{n}-1\) ].
See also: [mkpp], page 811.

\subsection*{28.6 Miscellaneous Functions}
```

poly (A)

```
poly (x)

If \(A\) is a square \(N\)-by- \(N\) matrix, poly (A) is the row vector of the coefficients of det
( \(\mathrm{z} *\) eye ( N ) - A), the characteristic polynomial of \(A\).
For example, the following code finds the eigenvalues of \(A\) which are the roots of poly
(A).
```

roots (poly (eye (3)))
=>1.00001 + 0.00001i
1.00001 - 0.00001i
0.99999 + 0.00000i

```

In fact, all three eigenvalues are exactly 1 which emphasizes that for numerical performance the eig function should be used to compute eigenvalues.
If \(x\) is a vector, poly ( \(x\) ) is a vector of the coefficients of the polynomial whose roots are the elements of \(x\). That is, if \(c\) is a polynomial, then the elements of \(d=\) roots (poly (c)) are contained in \(c\). The vectors \(c\) and \(d\) are not identical, however, due to sorting and numerical errors.
See also: [roots], page 798, [eig], page 587.
polyout (c)
polyout ( \(c, x\) )
str = polyout (...)
Display a formatted version of the polynomial \(c\).
The formatted polynomial
\[
c(x)=c_{1} x^{n}+\ldots+c_{n} x+c_{n+1}
\]
is returned as a string or written to the screen if nargout is zero.
The second argument \(x\) specifies the variable name to use for each term and defaults to the string "s".
See also: [polyreduce], page 813.
polyreduce (c)
Reduce a polynomial coefficient vector to a minimum number of terms by stripping off any leading zeros.

See also: [polyout], page 813.

\section*{29 Interpolation}

\subsection*{29.1 One-dimensional Interpolation}

Octave supports several methods for one-dimensional interpolation, most of which are described in this section. Section 28.5 [Polynomial Interpolation], page 803, and Section 30.4 [Interpolation on Scattered Data], page 838, describe additional methods.
```

yi = interp1 (x, y, xi)
yi = interp1 (y, xi)
yi = interp1 (..., method)
yi = interp1 (..., extrap)
yi = interp1 (..., "left")
yi = interp1 (..., "right")
pp = interp1 (..., "pp")

```

One-dimensional interpolation.
Interpolate input data to determine the value of \(y i\) at the points xi. If not specified, \(x\) is taken to be the indices of \(y\) (1:length ( \(y\) ) ). If \(y\) is a matrix or an N-dimensional array, the interpolation is performed on each column of \(y\).
The interpolation method is one of:
```

"nearest"
Return the nearest neighbor.
"previous"
Return the previous neighbor.
"next" Return the next neighbor.
"linear" (default)

```

Linear interpolation from nearest neighbors.
"pchip" Piecewise cubic Hermite interpolating polynomial-shape-preserving interpolation with smooth first derivative.
"cubic" Cubic interpolation (same as "pchip").
"spline" Cubic spline interpolation - smooth first and second derivatives throughout the curve.

Adding ' \({ }^{*}\) ' to the start of any method above forces interp1 to assume that x is uniformly spaced, and only \(x(1)\) and \(x(2)\) are referenced. This is usually faster, and is never slower. The default method is "linear".
If extrap is the string "extrap", then extrapolate values beyond the endpoints using the current method. If extrap is a number, then replace values beyond the endpoints with that number. When unspecified, extrap defaults to NA.
If the string argument "pp" is specified, then xi should not be supplied and interp1 returns a piecewise polynomial object. This object can later be used with ppval to evaluate the interpolation. There is an equivalence, such that ppval (interp1 ( \(x\), y, method, "pp"), xi) == interp1 (x, y, xi, method, "extrap").

Duplicate points in \(x\) specify a discontinuous interpolant. There may be at most 2 consecutive points with the same value. If \(x\) is increasing, the default discontinuous interpolant is right-continuous. If \(x\) is decreasing, the default discontinuous interpolant is left-continuous. The continuity condition of the interpolant may be specified by using the options "left" or "right" to select a left-continuous or right-continuous interpolant, respectively. Discontinuous interpolation is only allowed for "nearest" and "linear" methods; in all other cases, the \(x\)-values must be unique.
An example of the use of interp1 is
```

xf = [0:0.05:10];
yf = sin (2*pi*xf/5);
xp = [0:10];
yp = sin (2*pi*xp/5);
lin = interp1 (xp, yp, xf);
near = interp1 (xp, yp, xf, "nearest");
pch = interp1 (xp, yp, xf, "pchip");
spl = interp1 (xp, yp, xf, "spline");
plot (xf,yf,"r", xf,near,"g", xf,lin,"b", xf,pch,"c", xf,spl,"m",
xp,yp,"r*");
legend ("original", "nearest", "linear", "pchip", "spline");

```

See also: [pchip], page 854, [spline], page 818, [interpft], page 817, [interp2], page 819, [interp3], page 820, [interpn], page 821.

There are some important differences between the various interpolation methods. The "spline" method enforces that both the first and second derivatives of the interpolated values have a continuous derivative, whereas the other methods do not. This means that the results of the "spline" method are generally smoother. If the function to be interpolated is in fact smooth, then "spline" will give excellent results. However, if the function to be evaluated is in some manner discontinuous, then "pchip" interpolation might give better results.

This can be demonstrated by the code
```

t = -2:2;
dt = 1;
ti =-2:0.025:2;
dti = 0.025;
y = sign (t);
ys = interp1 (t,y,ti,"spline");
yp = interp1 (t,y,ti,"pchip");
ddys = diff (diff (ys)./dti) ./ dti;
ddyp = diff (diff (yp)./dti) ./ dti;
figure (1);
plot (ti,ys,"r-", ti,yp,"g-");
legend ("spline", "pchip", 4);
figure (2);
plot (ti,ddys,"r+", ti,ddyp,"g*");
legend ("spline", "pchip");

```

The result of which can be seen in Figure 29.1 and Figure 29.2.


Figure 29.1: Comparison of "pchip" and "spline" interpolation methods for a step function


Figure 29.2: Comparison of the second derivative of the "pchip" and "spline" interpolation methods for a step function

Fourier interpolation, is a resampling technique where a signal is converted to the frequency domain, padded with zeros and then reconverted to the time domain.
interpft ( \(x, n\) )
interpft ( \(x, n\), dim)
Fourier interpolation.

If \(x\) is a vector then \(x\) is resampled with \(n\) points. The data in \(x\) is assumed to be equispaced. If \(x\) is a matrix or an N -dimensional array, the interpolation is performed on each column of \(x\).

If \(\operatorname{dim}\) is specified, then interpolate along the dimension dim.
interpft assumes that the interpolated function is periodic, and so assumptions are made about the endpoints of the interpolation.

See also: [interp1], page 815.

There are two significant limitations on Fourier interpolation. First, the function signal is assumed to be periodic, and so non-periodic signals will be poorly represented at the edges. Second, both the signal and its interpolation are required to be sampled at equispaced points. An example of the use of interpft is
```

t = 0 : 0.3 : pi; dt = t(2)-t(1);
n = length (t); k = 100;
ti = t(1) + [0 : k-1]*dt*n/k;
y = sin (4*t + 0.3).* cos (3*t - 0.1);
yp = sin (4*ti + 0.3).* cos (3*ti - 0.1);
plot (ti, yp, "g", ti, interp1 (t, y, ti, "spline"), "b", ...
ti, interpft (y, k), "c", t, y, "r+");
legend ("sin(4t+0.3)cos(3t-0.1)", "spline", "interpft", "data");

```
which demonstrates the poor behavior of Fourier interpolation for non-periodic functions, as can be seen in Figure 29.3.


Figure 29.3: Comparison of interp1 and interpft for non-periodic data
In addition, the support functions spline and lookup that underlie the interp 1 function can be called directly.
```

pp = spline (x, y)
yi = spline (x, y, xi)

```

Return the cubic spline interpolant of points \(x\) and \(y\).
When called with two arguments, return the piecewise polynomial \(p p\) that may be used with ppval to evaluate the polynomial at specific points.
When called with a third input argument, spline evaluates the spline at the points xi. The third calling form spline ( \(x, y, x i\) ) is equivalent to ppval (spline ( \(x\), \(y), x i)\).
The variable x must be a vector of length \(n\).
\(y\) can be either a vector or array. If \(y\) is a vector it must have a length of either \(n\) or \(n+2\). If the length of \(y\) is \(n\), then the "not-a-knot" end condition is used. If the length of \(y\) is \(n+2\), then the first and last values of the vector \(y\) are the values of the first derivative of the cubic spline at the endpoints.
If \(y\) is an array, then the size of \(y\) must have the form
\[
\left[s_{1}, s_{2}, \cdots, s_{k}, n\right]
\]
or
\[
\left[s_{1}, s_{2}, \cdots, s_{k}, n+2\right] .
\]

The array is reshaped internally to a matrix where the leading dimension is given by
\[
s_{1} s_{2} \cdots s_{k}
\]
and each row of this matrix is then treated separately. Note that this is exactly the opposite of interp1 but is done for matlab compatibility.

See also: [pchip], page 854, [ppval], page 812, [mkpp], page 811, [unmkpp], page 811.

\subsection*{29.2 Multi-dimensional Interpolation}

There are three multi-dimensional interpolation functions in Octave, with similar capabilities. Methods using Delaunay tessellation are described in Section 30.4 [Interpolation on Scattered Data], page 838.
```

zi = interp2 (x, y, z, xi, yi)
zi = interp2 (z, xi, yi)
zi = interp2 (z, n)
zi = interp2 (z)
zi = interp2 (..., method)
zi = interp2 (..., method, extrap)
Two-dimensional interpolation.

```

Interpolate reference data \(x, y, z\) to determine \(z i\) at the coordinates xi, yi. The reference data \(x, y\) can be matrices, as returned by meshgrid, in which case the sizes of \(x, y\), and \(z\) must be equal. If \(x, y\) are vectors describing a grid then length ( \(x\) ) \(==\) columns ( \(z\) ) and length ( \(y\) ) == rows ( \(z\) ). In either case the input data must be strictly monotonic.

If called without \(x, y\), and just a single reference data matrix \(z\), the 2-D region \(x=\) 1: columns ( \(z\) ) , \(y=1\) :rows ( \(z\) ) is assumed. This saves memory if the grid is regular and the distance between points is not important.
If called with a single reference data matrix \(z\) and a refinement value \(n\), then perform interpolation over a grid where each original interval has been recursively subdivided \(n\) times. This results in \(2^{\wedge} n-1\) additional points for every interval in the original grid. If \(n\) is omitted a value of 1 is used. As an example, the interval \([0,1]\) with \(n==2\) results in a refined interval with points at \([0,1 / 4,1 / 2,3 / 4,1]\).
The interpolation method is one of:
```

"nearest"
Return the nearest neighbor.
"linear" (default)

```

Linear interpolation from nearest neighbors.
"pchip" Piecewise cubic Hermite interpolating polynomial-shape-preserving interpolation with smooth first derivative.
"cubic" Cubic interpolation (same as "pchip").
"spline" Cubic spline interpolation-smooth first and second derivatives throughout the curve.
extrap is a scalar number. It replaces values beyond the endpoints with extrap. Note that if extrap is used, method must be specified as well. If extrap is omitted and the method is "spline", then the extrapolated values of the "spline" are used. Otherwise the default extrap value for any other method is "NA".
See also: [interp1], page 815, [interp3], page 820, [interpn], page 821, [meshgrid], page 373.
```

vi = interp3 (x, y, z, v, xi, yi, zi)
vi = interp3 (v, xi, yi, zi)
vi = interp3 (v, n)
vi = interp3 (v)
vi = interp3 (..., method)
vi = interp3 (..., method, extrapval)

```

Three-dimensional interpolation.
Interpolate reference data \(x, y, z, v\) to determine vi at the coordinates xi, yi, zi. The reference data \(x, y, z\) can be matrices, as returned by meshgrid, in which case the sizes of \(x, y, z\), and \(v\) must be equal. If \(x, y, z\) are vectors describing a cubic grid then length ( \(x\) ) == columns (v), length ( \(y\) ) == rows (v), and length ( \(z\) ) == size ( \(v\), 3 ). In either case the input data must be strictly monotonic.
If called without \(x, y, z\), and just a single reference data matrix \(v\), the 3 -D region \(x=\) 1: columns (v), y = 1:rows (v), \(z=1\) : size ( \(v, 3\) ) is assumed. This saves memory if the grid is regular and the distance between points is not important.
If called with a single reference data matrix \(v\) and a refinement value \(n\), then perform interpolation over a 3-D grid where each original interval has been recursively subdivided \(n\) times. This results in \(2^{\wedge} n-1\) additional points for every interval in the
original grid. If \(n\) is omitted a value of 1 is used. As an example, the interval \([0,1]\) with \(n==2\) results in a refined interval with points at \([0,1 / 4,1 / 2,3 / 4,1]\).
The interpolation method is one of:
```

"nearest"
Return the nearest neighbor.
"linear" (default)

```

Linear interpolation from nearest neighbors.
"cubic" Piecewise cubic Hermite interpolating polynomial-shape-preserving interpolation with smooth first derivative (not implemented yet).
"spline" Cubic spline interpolation-smooth first and second derivatives throughout the curve.
extrapval is a scalar number. It replaces values beyond the endpoints with extrapval. Note that if extrapval is used, method must be specified as well. If extrapval is omitted and the method is "spline", then the extrapolated values of the "spline" are used. Otherwise the default extrapval value for any other method is "NA".

See also: [interp1], page 815, [interp2], page 819, [interpn], page 821, [meshgrid], page 373 .
```

vi = interpn (x1, x2,..., v, y1, y2,...)
vi = interpn (v, y1, y2,...)
vi = interpn (v, m)
vi = interpn (v)
vi = interpn (..., method)
vi = interpn (..., method, extrapval)

```

Perform \(n\)-dimensional interpolation, where \(n\) is at least two.
Each element of the n-dimensional array \(v\) represents a value at a location given by the parameters \(x 1, \mathrm{x} 2, \ldots\), xn . The parameters \(\mathrm{x} 1, \mathrm{x} 2, \ldots\), xn are either \(n\)-dimensional arrays of the same size as the array \(v\) in the "ndgrid" format or vectors.
The parameters \(y 1, y 2, \ldots, y n\) represent the points at which the array vi is interpolated. They can be vectors of the same length and orientation in which case they are interpreted as coordinates of scattered points. If they are vectors of differing orientation or length, they are used to form a grid in "ndgrid" format. They can also be \(n\)-dimensional arrays of equal size.
If \(x 1, \ldots\), \(x n\) are omitted, they are assumed to be \(x 1=1\) : size ( \(v, 1\) ), etc. If \(m\) is specified, then the interpolation adds a point half way between each of the interpolation points. This process is performed \(m\) times. If only \(v\) is specified, then \(m\) is assumed to be 1 .
The interpolation method is one of:
"nearest"
Return the nearest neighbor.
"linear" (default)
Linear interpolation from nearest neighbors.
"pchip" Piecewise cubic Hermite interpolating polynomial-shape-preserving interpolation with smooth first derivative (not implemented yet).
"cubic" Cubic interpolation (same as "pchip" [not implemented yet]).
"spline" Cubic spline interpolation-smooth first and second derivatives throughout the curve.

The default method is "linear".
extrapval is a scalar number. It replaces values beyond the endpoints with extrapval. Note that if extrapval is used, method must be specified as well. If extrapval is omitted and the method is "spline", then the extrapolated values of the "spline" are used. Otherwise the default extrapval value for any other method is NA.

See also: [interp1], page 815, [interp2], page 819, [interp3], page 820, [spline], page 818, [ndgrid], page 374.

A significant difference between interpn and the other two multi-dimensional interpolation functions is the fashion in which the dimensions are treated. For interp2 and interp3, the y -axis is considered to be the columns of the matrix, whereas the x -axis corresponds to the rows of the array. As Octave indexes arrays in column major order, the first dimension of any array is the columns, and so interpn effectively reverses the ' \(x\) ' and ' \(y\) ' dimensions. Consider the example,
```

x = y = z = -1:1;
f = @(x,y,z) x.`2 - y - z.^2;
[xx, yy, zz] = meshgrid (x, y, z);
v = f (xx,yy,zz);
xi = yi = zi = -1:0.1:1;
[xxi, yyi, zzi] = meshgrid (xi, yi, zi);
vi = interp3 (x, y, z, v, xxi, yyi, zzi, "spline");
[xxi, yyi, zzi] = ndgrid (xi, yi, zi);
vi2 = interpn (x, y, z, v, xxi, yyi, zzi, "spline");
mesh (zi, yi, squeeze (vi2(1,:,:)));

```
where vi and vi2 are identical. The reversal of the dimensions is treated in the meshgrid and ndgrid functions respectively. The result of this code can be seen in Figure 29.4.


Figure 29.4: Demonstration of the use of interpn

\section*{30 Geometry}

Much of the geometry code in Octave is based on the Qhull library \({ }^{1}\). Some of the documentation for Qhull, particularly for the options that can be passed to delaunay, voronoi and convhull, etc., is relevant to Octave users.

\subsection*{30.1 Delaunay Triangulation}

The Delaunay triangulation is constructed from a set of circum-circles. These circum-circles are chosen so that there are at least three of the points in the set to triangulation on the circumference of the circum-circle. None of the points in the set of points falls within any of the circum-circles.

In general there are only three points on the circumference of any circum-circle. However, in some cases, and in particular for the case of a regular grid, 4 or more points can be on a single circum-circle. In this case the Delaunay triangulation is not unique.
```

tri = delaunay (x, y)
tetr = delaunay (x, y, z)
tri = delaunay (x)
tri = delaunay (..., options)
Compute the Delaunay triangulation for a 2-D or 3-D set of points.

```

For 2-D sets, the return value tri is a set of triangles which satisfies the Delaunay circum-circle criterion, i.e., no data point from \([x, y]\) is within the circum-circle of the defining triangle. The set of triangles tri is a matrix of size [n, 3]. Each row defines a triangle and the three columns are the three vertices of the triangle. The value of \(\operatorname{tri}(i, j)\) is an index into \(x\) and \(y\) for the location of the \(j\)-th vertex of the i -th triangle.
For 3-D sets, the return value tetr is a set of tetrahedrons which satisfies the Delaunay circum-circle criterion, i.e., no data point from \([x, y, z]\) is within the circum-circle of the defining tetrahedron. The set of tetrahedrons is a matrix of size \([\mathrm{n}, 4]\). Each row defines a tetrahedron and the four columns are the four vertices of the tetrahedron. The value of \(\operatorname{tetr}(i, j)\) is an index into \(x, y, z\) for the location of the \(j\)-th vertex of the i-th tetrahedron.

The input \(x\) may also be a matrix with two or three columns where the first column contains x-data, the second y-data, and the optional third column contains z-data.
An optional final argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options. The default options are \{"Qt", "Qbb", "Qc"\}. If Qhull fails for 2-D input the triangulation is attempted again with the options \{"Qt", "Qbb", "Qc", "Qz"\} which may result in reduced accuracy.
If options is not present or [] then the default arguments are used. Otherwise, options replaces the default argument list. To append user options to the defaults it

\footnotetext{
\({ }^{1}\) Barber, C.B., Dobkin, D.P., and Huhdanpaa, H.T., The Quickhull Algorithm for Convex Hulls, ACM Trans. on Mathematical Software, 22(4):469-483, Dec 1996, http://www.qhull.org
}
is necessary to repeat the default arguments in options. Use a null string to pass no arguments.
```

x = rand (1, 10);
y = rand (1, 10);
tri = delaunay (x, y);
triplot (tri, x, y);
hold on;
plot (x, y, "r*");
axis ([0,1,0,1]);

```

See also: [delaunayn], page 826, [convhull], page 836, [voronoi], page 832, [triplot], page 827, [trimesh], page 828, [tetramesh], page 829, [trisurf], page 828.

For 3-D inputs delaunay returns a set of tetrahedra that satisfy the Delaunay circumcircle criteria. Similarly, delaunayn returns the N-dimensional simplex satisfying the Delaunay circum-circle criteria. The N -dimensional extension of a triangulation is called a tessellation.
```

T = delaunayn (pts)
T = delaunayn (pts, options)

```

Compute the Delaunay triangulation for an N-dimensional set of points.
The Delaunay triangulation is a tessellation of the convex hull of a set of points such that no N -sphere defined by the N -triangles contains any other points from the set.
The input matrix pts of size \([\mathrm{n}, \operatorname{dim}]\) contains n points in a space of dimension dim. The return matrix \(T\) has size \([\mathrm{m}, \operatorname{dim}+1]\). Each row of \(T\) contains a set of indices back into the original set of points pts which describes a simplex of dimension dim. For example, a 2-D simplex is a triangle and 3-D simplex is a tetrahedron.

An optional second argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options. The default options depend on the dimension of the input:
- 2-D and 3-D: options = \{"Qt", "Qbb", "Qc"\}
- 4-D and higher: options = \{"Qt", "Qbb", "Qc", "Qx"\}

If Qhull fails for 2-D input the triangulation is attempted again with the options \{"Qt", "Qbb", "Qc", "Qz"\} which may result in reduced accuracy.
If options is not present or [] then the default arguments are used. Otherwise, options replaces the default argument list. To append user options to the defaults it is necessary to repeat the default arguments in options. Use a null string to pass no arguments.

See also: [delaunay], page 825, [convhulln], page 837, [voronoin], page 833, [trimesh], page 828 , [tetramesh], page 829.

An example of a Delaunay triangulation of a set of points is
```

rand ("state", 1);
x = rand (1, 10);
y = rand (1, 10);
T = delaunay (x, y);
X = [ x(T(:,1)); x(T(:,2)); x(T(:,3)); x(T(:,1)) ];
Y = [ y(T(:,1)); y(T(:,2)); y(T(:,3)); y(T(:,1)) ];
axis ([0, 1, 0, 1]);
plot (X, Y, "b", x, y, "r*");

```

The result of which can be seen in Figure 30.1.


Figure 30.1: Delaunay triangulation of a random set of points

\subsection*{30.1.1 Plotting the Triangulation}

Octave has the functions triplot, trimesh, and trisurf to plot the Delaunay triangulation of a 2-dimensional set of points. tetramesh will plot the triangulation of a 3-dimensional set of points.
triplot (tri, \(x, y\) )
triplot (tri, x, y, linespec)
h = triplot (...)
Plot a 2-D triangular mesh.
tri is typically the output of a Delaunay triangulation over the grid of \(x, y\). Every row of tri represents one triangle and contains three indices into \([x, y]\) which are the vertices of the triangles in the \(x-y\) plane.

The linestyle to use for the plot can be defined with the argument linespec of the same format as the plot command.
The optional return value \(h\) is a graphics handle to the created patch object.
See also: [plot], page 316, [trimesh], page 828, [trisurf], page 828, [delaunay], page 825.
```

trimesh (tri, x, y, z, c)
trimesh (tri, x, y, z)
trimesh (tri, x, y)
trimesh (..., prop, val, ...)
h = trimesh (...)

```

Plot a 3-D triangular wireframe mesh.
In contrast to mesh, which plots a mesh using rectangles, trimesh plots the mesh using triangles.
tri is typically the output of a Delaunay triangulation over the grid of \(x, y\). Every row of tri represents one triangle and contains three indices into \([x, y]\) which are the vertices of the triangles in the \(x-y\) plane. \(z\) determines the height above the plane of each vertex. If no \(z\) input is given then the triangles are plotted as a 2-D figure.
The color of the trimesh is computed by linearly scaling the \(z\) values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally, the color of the mesh can be specified independently of \(z\) by supplying \(c\), which is a vector for colormap data, or a matrix with three columns for RGB data. The number of colors specified in \(c\) must either equal the number of vertices in \(z\) or the number of triangles in tri.
Any property/value pairs are passed directly to the underlying patch object. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.
The optional return value \(h\) is a graphics handle to the created patch object.
See also: [mesh], page 358, [tetramesh], page 829, [triplot], page 827, [trisurf], page 828 , [delaunay], page 825 , [patch], page 425 , [hidden], page 360 .
```

trisurf (tri, x, y, z, c)

```
trisurf (tri, x, y, z)
trisurf (..., prop, val, ...)
\(h=\) trisurf (...)

Plot a 3-D triangular surface.
In contrast to surf, which plots a surface mesh using rectangles, trisurf plots the mesh using triangles.
tri is typically the output of a Delaunay triangulation over the grid of \(x, y\). Every row of tri represents one triangle and contains three indices into \([x, y]\) which are the vertices of the triangles in the \(x-y\) plane. \(z\) determines the height above the plane of each vertex.
The color of the trisurf is computed by linearly scaling the \(z\) values to fit the range of the current colormap. Use caxis and/or change the colormap to control the appearance.
Optionally, the color of the mesh can be specified independently of \(z\) by supplying \(c\), which is a vector for colormap data, or a matrix with three columns for RGB data. The number of colors specified in \(c\) must either equal the number of vertices in \(z\) or the number of triangles in tri. When specifying the color at each vertex the triangle will be colored according to the color of the first vertex only (see patch documentation and the "FaceColor" property when set to "flat").

Any property/value pairs are passed directly to the underlying patch object. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.

The optional return value \(h\) is a graphics handle to the created patch object.
See also: [surf], page 360, [triplot], page 827, [trimesh], page 828, [delaunay], page 825, [patch], page 425, [shading], page 382.
tetramesh ( \(T, X\) )
tetramesh ( \(T, X, C\) )
tetramesh (..., property, val, ...)
\(h=\) tetramesh (...)
Display the tetrahedrons defined in the m-by-4 matrix \(T\) as 3-D patches.
\(T\) is typically the output of a Delaunay triangulation of a 3-D set of points. Every row of \(T\) contains four indices into the n-by- 3 matrix \(X\) of the vertices of a tetrahedron. Every row in \(X\) represents one point in 3-D space.

The vector \(C\) specifies the color of each tetrahedron as an index into the current colormap. The default value is \(1: \mathrm{m}\) where m is the number of tetrahedrons; the indices are scaled to map to the full range of the colormap. If there are more tetrahedrons than colors in the colormap then the values in \(C\) are cyclically repeated.

Calling tetramesh (..., "property", "value", ...) passes all property/value pairs directly to the patch function as additional arguments. The full list of properties is documented at Section 15.3.3.8 [Patch Properties], page 457.

The optional return value \(h\) is a vector of patch handles where each handle represents one tetrahedron in the order given by \(T\). A typical use case for \(h\) is to turn the respective patch "visible" property "on" or "off".

Type demo tetramesh to see examples on using tetramesh.
See also: [trimesh], page 828, [delaunay], page 825, [delaunayn], page 826, [patch], page 425 .

The difference between triplot, and trimesh or trisurf, is that the former only plots the 2-dimensional triangulation itself, whereas the second two plot the value of a function \(f(x, y)\). An example of the use of the triplot function is
```

rand ("state", 2)
x = rand (20, 1);
y = rand (20, 1);
tri = delaunay (x, y);
triplot (tri, x, y);

```
which plots the Delaunay triangulation of a set of random points in 2-dimensions. The output of the above can be seen in Figure 30.2.


Figure 30.2: Delaunay triangulation of a random set of points

\subsection*{30.1.2 Identifying Points in Triangulation}

It is often necessary to identify whether a particular point in the N-dimensional space is within the Delaunay tessellation of a set of points in this N-dimensional space, and if so which N -simplex contains the point and which point in the tessellation is closest to the desired point. The functions tsearch and dsearch perform this function in a triangulation, and tsearchn and dsearchn in an N-dimensional tessellation.

To identify whether a particular point represented by a vector \(p\) falls within one of the simplices of an N-simplex, we can write the Cartesian coordinates of the point in a parametric form with respect to the N -simplex. This parametric form is called the Barycentric Coordinates of the point. If the points defining the N -simplex are given by \(N+1\) vectors \(t(i,:)\), then the Barycentric coordinates defining the point \(p\) are given by
```

p = beta * t

```
where beta contains \(N+1\) values that together as a vector represent the Barycentric coordinates of the point \(p\). To ensure a unique solution for the values of beta an additional criteria of
```

sum (beta) == 1

```
is imposed, and we can therefore write the above as
```

p - t(end, :) = beta(1:end-1) * (t(1:end-1, :)
- ones (N, 1) * t(end, :)

```

Solving for beta we can then write
```

beta(1:end-1) = (p - t(end, :)) /
(t(1:end-1, :) - ones (N, 1) * t(end, :))
beta(end) = sum (beta(1:end-1))

```
which gives the formula for the conversion of the Cartesian coordinates of the point \(p\) to the Barycentric coordinates beta. An important property of the Barycentric coordinates is that for all points in the N -simplex
```

0<= beta(i) <= 1

```

Therefore, the test in tsearch and tsearchn essentially only needs to express each point in terms of the Barycentric coordinates of each of the simplices of the N -simplex and test the values of beta. This is exactly the implementation used in tsearchn. tsearch is optimized for 2-dimensions and the Barycentric coordinates are not explicitly formed.
```

idx = tsearch (x, y, t, xi, yi)

```

Search for the enclosing Delaunay convex hull.
For \(t=\) delaunay ( \(x, y\) ), finds the index in \(t\) containing the points (xi, yi). For points outside the convex hull, idx is NaN.
See also: [delaunay], page 825, [delaunayn], page 826 .
\(i d x=t s e a r c h n(x, t, x i)\)
[idx, p] = tsearchn (x, t, xi)
Search for the enclosing Delaunay convex hull.
For \(t=\) delaunayn ( \(x\) ), finds the index in \(t\) containing the points xi. For points outside the convex hull, idx is NaN.
If requested tsearchn also returns the Barycentric coordinates \(p\) of the enclosing triangles.
See also: [delaunay], page 825, [delaunayn], page 826.
An example of the use of tsearch can be seen with the simple triangulation
```

x = [-1; -1; 1; 1];
y = [-1; 1; -1; 1];
tri = [1, 2, 3; 2, 3, 4];

```
consisting of two triangles defined by tri. We can then identify which triangle a point falls in like
```

tsearch (x, y, tri, -0.5, -0.5)

# 1

tsearch (x, y, tri, 0.5, 0.5)
=> 2

```
and we can confirm that a point doesn't lie within one of the triangles like
```

tsearch (x, y, tri, 2, 2)

# NaN

```

The dsearch and dsearchn find the closest point in a tessellation to the desired point. The desired point does not necessarily have to be in the tessellation, and even if it the returned point of the tessellation does not have to be one of the vertices of the N -simplex within which the desired point is found.
```

idx = dsearch (x, y, tri, xi, yi)
idx = dsearch (x, y, tri, xi, yi, s)

```

Return the index idx of the closest point in \(x, y\) to the elements [xi(:), yi(:)].

The variable \(s\) is accepted for compatibility but is ignored.
See also: [dsearchn], page 832, [tsearch], page 831.
\(i d x=d s e a r c h n(x, t r i, x i)\)
idx = dsearchn (x, tri, xi, outval)
\(i d x=\) dsearchn ( \(x, x i\) )
[idx, d] = dsearchn (...)
Return the index idx of the closest point in \(x\) to the elements xi.
If outval is supplied, then the values of xi that are not contained within one of the simplices tri are set to outval. Generally, tri is returned from delaunayn (x).
The optional output \(d\) contains a column vector of distances between the query points xi and the nearest simplex points \(x\).
See also: [dsearch], page 831, [tsearch], page 831.
An example of the use of dsearch, using the above values of \(x, y\) and tri is
```

dsearch (x, y, tri, -2, -2)

# 1

```

If you wish the points that are outside the tessellation to be flagged, then dsearchn can be used as
```

dsearchn ([x, y], tri, [-2, -2], NaN)

# NaN

dsearchn ([x, y], tri, [-0.5, -0.5], NaN)

# 1

```
where the point outside the tessellation are then flagged with NaN.

\subsection*{30.2 Voronoi Diagrams}

A Voronoi diagram or Voronoi tessellation of a set of points \(s\) in an N-dimensional space, is the tessellation of the \(N\)-dimensional space such that all points in \(v(p)\), a partitions of the tessellation where \(p\) is a member of \(s\), are closer to \(p\) than any other point in \(s\). The Voronoi diagram is related to the Delaunay triangulation of a set of points, in that the vertices of the Voronoi tessellation are the centers of the circum-circles of the simplices of the Delaunay tessellation.
```

voronoi (x, y)
voronoi (x, y, options)
voronoi (..., "linespec")
voronoi (hax,...)
h = voronoi (...)
[vx, vy] = voronoi (...)
Plot the Voronoi diagram of points ( $x, y$ ).

```

The Voronoi facets with points at infinity are not drawn.
The options argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options.

If "linespec" is given it is used to set the color and line style of the plot.
If an axes graphics handle hax is supplied then the Voronoi diagram is drawn on the specified axes rather than in a new figure.
If a single output argument is requested then the Voronoi diagram will be plotted and a graphics handle \(h\) to the plot is returned.
\([v x, v y]=\) voronoi \((\ldots)\) returns the Voronoi vertices instead of plotting the diagram.
```

x = rand (10, 1);
y = rand (size (x));
h = convhull (x, y);
[vx, vy] = voronoi (x, y);
plot (vx, vy, "-b", x, y, "o", x(h), y(h), "-g");
legend ("", "points", "hull");

```

See also: [voronoin], page 833, [delaunay], page 825, [convhull], page 836.
\([C, F]=\) voronoin (pts)
\([C, F]=\) voronoin (pts, options)
Compute N-dimensional Voronoi facets.
The input matrix pts of size [n, dim] contains n points in a space of dimension dim.
\(C\) contains the points of the Voronoi facets. The list \(F\) contains, for each facet, the indices of the Voronoi points.
An optional second argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options. The default options depend on the dimension of the input:
- 2-D and 3-D: options \(=\{\) "Qbb" \(\}\)
- 4-D and higher: options = \{"Qbb", "Qx"\}

If options is not present or [] then the default arguments are used. Otherwise, options replaces the default argument list. To append user options to the defaults it is necessary to repeat the default arguments in options. Use a null string to pass no arguments.
See also: [voronoi], page 832, [convhulln], page 837, [delaunayn], page 826.
An example of the use of voronoi is
```

rand ("state",9);
x = rand (10,1);
y = rand (10,1);
tri = delaunay (x, y);
[vx, vy] = voronoi (x, y, tri);
triplot (tri, x, y, "b");
hold on;
plot (vx, vy, "r");

```

The result of which can be seen in Figure 30.3. Note that the circum-circle of one of the triangles has been added to this figure, to make the relationship between the Delaunay tessellation and the Voronoi diagram clearer.


Figure 30.3: Delaunay triangulation (blue lines) and Voronoi diagram (red lines) of a random set of points

Additional information about the size of the facets of a Voronoi diagram, and which points of a set of points is in a polygon can be had with the polyarea and inpolygon functions respectively.
polyarea ( \(x, y\) )
polyarea ( \(x, y, d i m\) )
Determine area of a polygon by triangle method.
The variables \(x\) and \(y\) define the vertex pairs, and must therefore have the same shape. They can be either vectors or arrays. If they are arrays then the columns of \(x\) and \(y\) are treated separately and an area returned for each.
If the optional dim argument is given, then polyarea works along this dimension of the arrays \(x\) and \(y\).

An example of the use of polyarea might be
```

rand ("state", 2);
x = rand (10, 1);
y = rand (10, 1);
[c, f] = voronoin ([x, y]);
af = zeros (size (f));
for i = 1 : length (f)
af(i) = polyarea (c (f {i, :}, 1), c (f {i, :}, 2));
endfor

```

Facets of the Voronoi diagram with a vertex at infinity have infinity area. A simplified version of polyarea for rectangles is available with rectint
area \(=\operatorname{rectint}(a, b)\)
Compute area or volume of intersection of rectangles or N-D boxes.

Compute the area of intersection of rectangles in a and rectangles in \(b\). N-dimensional boxes are supported in which case the volume, or hypervolume is computed according to the number of dimensions.

2-dimensional rectangles are defined as [xpos ypos width height] where xpos and ypos are the position of the bottom left corner. Higher dimensions are supported where the coordinates for the minimum value of each dimension follow the length of the box in that dimension, e.g., [xpos ypos zpos kpos ... width height depth k_length ...].

Each row of \(a\) and \(b\) define a rectangle, and if both define multiple rectangles, then the output, area, is a matrix where the i-th row corresponds to the i-th row of a and the j -th column corresponds to the j -th row of b .

See also: [polyarea], page 834 .
in = inpolygon ( \(x, y, x v, y v\) )
[in, on] = inpolygon ( \(x, y, x v, y v\) )
For a polygon defined by vertex points ( \(x v, y v\) ), return true if the points ( \(x, y\) ) are inside (or on the boundary) of the polygon; Otherwise, return false.

The input variables \(x\) and \(y\), must have the same dimension.

The optional output on returns true if the points are exactly on the polygon edge, and false otherwise.

See also: [delaunay], page 825.

An example of the use of inpolygon might be
```

randn ("state", 2);
x = randn (100, 1);
y = randn (100, 1);
vx = cos (pi * [-1 : 0.1: 1]);
vy = sin (pi * [-1 : 0.1 : 1]);
in = inpolygon (x, y, vx, vy);
plot (vx, vy, x(in), y(in), "r+", x(!in), y(!in), "bo");
axis ([-2, 2, -2, 2]);

```

The result of which can be seen in Figure 30.4.


Figure 30.4: Demonstration of the inpolygon function to determine the points inside a polygon

\subsection*{30.3 Convex Hull}

The convex hull of a set of points is the minimum convex envelope containing all of the points. Octave has the functions convhull and convhulln to calculate the convex hull of 2 -dimensional and N -dimensional sets of points.
```

H = convhull (x, y)
H = convhull (x, y, z)
H = convhull (x)
H = convhull (..., options)
[H,V] = convhull (...)

```

Compute the convex hull of a 2-D or 3-D set of points.
The hull \(H\) is a linear index vector into the original set of points that specifies which points form the enclosing hull. For 2-D inputs only, the output is ordered in a counterclockwise manner around the hull.
The input \(x\) may also be a matrix with two or three columns where the first column contains x-data, the second y-data, and the optional third column contains z-data.
An optional final argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options. The default option is \{"Qt"\}.
If options is not present or [] then the default arguments are used. Otherwise, options replaces the default argument list. To append user options to the defaults it is necessary to repeat the default arguments in options. Use a null string to pass no arguments.

If the second output \(V\) is requested the volume of the enclosing convex hull is calculated.

See also: [convhulln], page 837, [delaunay], page 825, [voronoi], page 832.
```

h = convhulln (pts)
h = convhulln (pts, options)
[h, v] = convhulln (...)

```

Compute the convex hull of the set of points pts.
\(p t s\) is a matrix of size \([\mathrm{n}, \operatorname{dim}]\) containing n points in a space of dimension dim.
The hull \(h\) is an index vector into the set of points and specifies which points form the enclosing hull.

An optional second argument, which must be a string or cell array of strings, contains options passed to the underlying qhull command. See the documentation for the Qhull library for details http://www.qhull.org/html/qh-quick.htm\#options. The default options depend on the dimension of the input:
- 2D, 3D, 4D: options = \{"Qt"\}
- 5D and higher: options \(=\{\) "Qt", "Qx" \(\}\)

If options is not present or [] then the default arguments are used. Otherwise, options replaces the default argument list. To append user options to the defaults it is necessary to repeat the default arguments in options. Use a null string to pass no arguments.

If the second output \(v\) is requested the volume of the enclosing convex hull is calculated.

See also: [convhull], page 836, [delaunayn], page 826, [voronoin], page 833.

An example of the use of convhull is
```

x = -3:0.05:3;
y = abs (sin (x));
k = convhull (x, y);
plot (x(k), y(k), "r-", x, y, "b+");
axis ([-3.05, 3.05, -0.05, 1.05]);

```

The output of the above can be seen in Figure 30.5.


Figure 30.5: The convex hull of a simple set of points

\subsection*{30.4 Interpolation on Scattered Data}

An important use of the Delaunay tessellation is that it can be used to interpolate from scattered data to an arbitrary set of points. To do this the N -simplex of the known set of points is calculated with delaunay or delaunayn. Then the simplices in to which the desired points are found are identified. Finally the vertices of the simplices are used to interpolate to the desired points. The functions that perform this interpolation are griddata, griddata3 and griddatan.
```

zi = griddata (x, y, z, xi, yi)
zi = griddata (x, y, z, xi, yi, method)
[xi, yi, zi] = griddata (...)
vi = griddata (x, y, z, v, xi, yi, zi)
vi = griddata (x, y, z, v, xi, yi, zi, method)
vi = griddata (x, y, z, v, xi, yi, zi, method, options)

```

Interpolate irregular 2-D and 3-D source data at specified points.
For 2-D interpolation, the inputs \(x\) and \(y\) define the points where the function \(z=f\) ( \(x, y\) ) is evaluated. The inputs \(x, y, z\) are either vectors of the same length, or the unequal vectors \(x, y\) are expanded to a 2-D grid with meshgrid and \(z\) is a 2-D matrix matching the resulting size of the X-Y grid.
The interpolation points are (xi, yi). If, and only if, xi is a row vector and yi is a column vector, then meshgrid will be used to create a mesh of interpolation points.
For 3-D interpolation, the inputs \(x, y\), and \(z\) define the points where the function \(v=\) \(f(x, y, z)\) is evaluated. The inputs \(x, y, z\) are either vectors of the same length, or if they are of unequal length, then they are expanded to a 3-D grid with meshgrid. The size of the input \(v\) must match the size of the original data, either as a vector or a matrix.

The optional input interpolation method can be "nearest", "linear", or for 2-D data "v4". When the method is "nearest", the output vi will be the closest point in the original data ( \(x, y, z\) ) to the query point (xi, yi, zi). When the method is "linear", the output vi will be a linear interpolation between the two closest points in the original source data in each dimension. For 2-D cases only, the "v4" method is also available which implements a biharmonic spline interpolation. If method is omitted or empty, it defaults to "linear".
For 3-D interpolation, the optional argument options is passed directly to Qhull when computing the Delaunay triangulation used for interpolation. For more information on the defaults and how to pass different values, see [delaunayn], page 826.
Programming Notes: If the input is complex the real and imaginary parts are interpolated separately. Interpolation is normally based on a Delaunay triangulation. Any query values outside the convex hull of the input points will return NaN. However, the "v4" method does not use the triangulation and will return values outside the original data (extrapolation).
See also: [griddata3], page 839, [griddatan], page 839, [delaunay], page 825.
vi = griddata3 (x, y, z, v, xi, yi, zi)
vi = griddata3 (x, y, z, v, xi, yi, zi, method)
vi = griddata3 (x, y, z, v, xi, yi, zi, method, options)
Interpolate irregular 3-D source data at specified points.
The inputs \(x, y\), and \(z\) define the points where the function \(v=f(x, y, z)\) is evaluated. The inputs \(x, y, z\) are either vectors of the same length, or if they are of unequal length, then they are expanded to a 3-D grid with meshgrid. The size of the input \(v\) must match the size of the original data, either as a vector or a matrix.
The interpolation points are specified by xi, yi, zi.
The optional input interpolation method can be "nearest" or "linear". When the method is "nearest", the output vi will be the closest point in the original data ( \(x\), \(y, z\) ) to the query point (xi, yi, zi). When the method is "linear", the output vi will be a linear interpolation between the two closest points in the original source data in each dimension. If method is omitted or empty, it defaults to "linear".
The optional argument options is passed directly to Qhull when computing the Delaunay triangulation used for interpolation. See delaunayn for information on the defaults and how to pass different values.
Programming Notes: If the input is complex the real and imaginary parts are interpolated separately. Interpolation is based on a Delaunay triangulation and any query values outside the convex hull of the input points will return NaN.

See also: [griddata], page 838, [griddatan], page 839, [delaunayn], page 826.
yi \(=\) griddatan ( \(x, y, x i\) )
yi = griddatan (x, y, xi, method)
yi = griddatan (x, y, xi, method, options)
Interpolate irregular source data \(x, y\) at points specified by xi.
The input \(x\) is an MxN matrix representing M points in an N-dimensional space. The input \(y\) is a single-valued column vector (Mx1) representing a function evaluated at
the points \(x\), i.e., \(y=f c n(x)\). The input \(x i\) is a list of points for which the function output yi should be approximated through interpolation. xi must have the same number of columns \((N)\) as \(x\) so that the dimensionality matches.

The optional input interpolation method can be "nearest" or "linear". When the method is "nearest", the output yi will be the closest point in the original data \(x\) to the query point xi. When the method is "linear", the output yi will be a linear interpolation between the two closest points in the original source data. If method is omitted or empty, it defaults to "linear".

The optional argument options is passed directly to Qhull when computing the Delaunay triangulation used for interpolation. See delaunayn for information on the defaults and how to pass different values.

Example
```


## Evaluate sombrero() function at irregular data points

x = 16*gallery ("uniformdata", [200,1], 1) - 8;
y = 16*gallery ("uniformdata", [200,1], 11) - 8;
z = sin (sqrt (x.^2 + y.^2)) ./ sqrt (x.^2 + y. ^2);

## Create a regular grid and interpolate data

[xi, yi] = ndgrid (linspace (-8, 8, 50));
zi = griddatan ([x, y], z, [xi(:), yi(:)]);
zi = reshape (zi, size (xi));

## Plot results

clf ();
plot3 (x, y, z, "or");
hold on
surf (xi, yi, zi);
legend ("Original Data", "Interpolated Data");

```

Programming Notes: If the input is complex the real and imaginary parts are interpolated separately. Interpolation is based on a Delaunay triangulation and any query values outside the convex hull of the input points will return NaN. For 2-D and 3-D data additional interpolation methods are available by using the griddata function.

See also: [griddata], page 838, [griddata3], page 839, [delaunayn], page 826.
An example of the use of the griddata function is
```

rand ("state", 1);
x = 2*rand (1000,1) - 1;
y = 2*rand (size (x)) - 1;
z = sin (2*(x.^2+y.^2));
[xx,yy] = meshgrid (linspace (-1,1,32));
zz = griddata (x, y, z, xx, yy);
mesh (xx, yy, zz);

```
that interpolates from a random scattering of points, to a uniform grid. The output of the above can be seen in Figure 30.6.


Figure 30.6: Interpolation from a scattered data to a regular grid

\subsection*{30.5 Vector Rotation Matrices}

Also included in Octave's geometry functions are primitive functions to enable vector rotations in 3-dimensional space. Separate functions are provided for rotation about each of the principle axes, \(x, y\), and \(z\). According to Euler's rotation theorem, any arbitrary rotation, \(R\), of any vector, \(p\), can be expressed as a product of the three principle rotations:
\(p^{\prime}=R \cdot p=R_{z} \cdot R_{y} \cdot R_{x} \cdot p\)
\(T=\operatorname{rotx}(\) angle \()\)
rotx returns the \(3 x 3\) transformation matrix corresponding to an active rotation of a vector about the x -axis by the specified angle, given in degrees, where a positive angle corresponds to a counterclockwise rotation when viewing the \(y-z\) plane from the positive x side.

The form of the transformation matrix is:
\[
T=\left[\begin{array}{ccc}
1 & 0 & 0 \\
0 & \cos (\text { angle }) & -\sin (\text { angle }) \\
0 & \sin (\text { angle }) & \cos (\text { angle })
\end{array}\right] .
\]

This rotation matrix is intended to be used as a left-multiplying matrix when acting on a column vector, using the notation \(v=T * u\). For example, a vector, \(u\), pointing along the positive \(y\)-axis, rotated 90 -degrees about the x -axis, will result in a vector pointing along the positive z -axis:
```

>> u = [0 1 0]'
u =
0
1
0
>> T = rotx (90)
T =
1.00000 0.00000 0.00000
0.00000 0.00000 -1.00000
0.00000 1.00000 0.00000
>> v = T*u
v =
0.00000
0.00000
1.00000

```

See also: [roty], page 842, [rotz], page 843.
\(T=\operatorname{roty}(\) angle \()\)
roty returns the \(3 x 3\) transformation matrix corresponding to an active rotation of a vector about the y -axis by the specified angle, given in degrees, where a positive angle corresponds to a counterclockwise rotation when viewing the z -x plane from the positive y side.

The form of the transformation matrix is:
\[
T=\left[\begin{array}{ccc}
\cos (\text { angle }) & 0 & \sin (\text { angle }) \\
0 & 1 & 0 \\
-\sin (\text { angle }) & 0 & \cos (\text { angle })
\end{array}\right] .
\]

This rotation matrix is intended to be used as a left-multiplying matrix when acting on a column vector, using the notation \(v=T * u\). For example, a vector, \(u\), pointing along the positive z -axis, rotated 90 -degrees about the y -axis, will result in a vector pointing along the positive x -axis:
```

>> u = [l0 0 1]'
u =
0
0
1
>> T = roty (90)
T =
0.00000 0.00000 1.00000
0.00000 1.00000 0.00000
-1.00000 0.00000 0.00000
>> v = T*u
v =
1.00000
0.00000
0.00000

```

See also: [rotx], page 841, [rotz], page 843.

\section*{\(T=\operatorname{rotz}(\) angle \()\)}
rotz returns the \(3 x 3\) transformation matrix corresponding to an active rotation of a vector about the \(z\)-axis by the specified angle, given in degrees, where a positive angle corresponds to a counterclockwise rotation when viewing the \(x-y\) plane from the positive z side.

The form of the transformation matrix is:
\[
T=\left[\begin{array}{ccc}
\cos (\text { angle }) & -\sin (\text { angle }) & 0 \\
\sin (\text { angle }) & \cos (\text { angle }) & 0 \\
0 & 0 & 1
\end{array}\right] .
\]

This rotation matrix is intended to be used as a left-multiplying matrix when acting on a column vector, using the notation \(v=T * u\). For example, a vector, \(u\), pointing along the positive x -axis, rotated 90 -degrees about the z -axis, will result in a vector pointing along the positive y -axis:
```

>> u = [11 0 0}\mp@subsup{]}{}{\prime
u =
1
0
0
>> T = rotz (90)
T =
0.00000 -1.00000 0.00000
1.00000 0.00000 0.00000
0.00000 0.00000 1.00000
>> v = T*u
v =
0.00000
1.00000
0.00000

```

See also: [rotx], page 841, [roty], page 842.

\section*{31 Signal Processing}

This chapter describes the signal processing and fast Fourier transform functions available in Octave. Fast Fourier transforms are computed with the FFTW or FFTPACK libraries depending on how Octave is built.
fft (x)
fft ( \(x, n\) )
fft (x, n, dim)
Compute the discrete Fourier transform of x using a Fast Fourier Transform (FFT) algorithm.
The FFT is calculated along the first non-singleton dimension of the array. Thus if \(x\) is a matrix, \(f f t(x)\) computes the FFT for each column of \(x\).
If called with two arguments, \(n\) is expected to be an integer specifying the number of elements of \(x\) to use, or an empty matrix to specify that its value should be ignored. If \(n\) is larger than the dimension along which the FFT is calculated, then \(x\) is resized and padded with zeros. Otherwise, if \(n\) is smaller than the dimension along which the FFT is calculated, then \(x\) is truncated.
If called with three arguments, dim is an integer specifying the dimension of the matrix along which the FFT is performed.
See also: [ifft], page 845, [fft2], page 845, [fftn], page 846, [fftw], page 846.
ifft (x)
ifft (x, n)
ifft ( \(x, n\), dim)
Compute the inverse discrete Fourier transform of \(x\) using a Fast Fourier Transform (FFT) algorithm.
The inverse FFT is calculated along the first non-singleton dimension of the array. Thus if \(x\) is a matrix, \(f f t(x)\) computes the inverse FFT for each column of \(x\).
If called with two arguments, \(n\) is expected to be an integer specifying the number of elements of \(x\) to use, or an empty matrix to specify that its value should be ignored. If \(n\) is larger than the dimension along which the inverse FFT is calculated, then \(x\) is resized and padded with zeros. Otherwise, if \(n\) is smaller than the dimension along which the inverse FFT is calculated, then \(x\) is truncated.
If called with three arguments, dim is an integer specifying the dimension of the matrix along which the inverse FFT is performed.

See also: [fft], page 845, [ifft2], page 846, [ifftn], page 846, [fftw], page 846.
fft2 (A)
fft2 (A, m, n)
Compute the two-dimensional discrete Fourier transform of \(A\) using a Fast Fourier Transform (FFT) algorithm.
The optional arguments \(m\) and \(n\) may be used specify the number of rows and columns of \(A\) to use. If either of these is larger than the size of \(A, A\) is resized and padded with zeros.

If \(A\) is a multi-dimensional matrix, each two-dimensional sub-matrix of \(A\) is treated separately.

See also: [ifft2], page 846, [fft], page 845, [fftn], page 846, [fftw], page 846.
ifft2 (A)
ifft2 (A, m, n)
Compute the inverse two-dimensional discrete Fourier transform of \(A\) using a Fast Fourier Transform (FFT) algorithm.
The optional arguments \(m\) and \(n\) may be used specify the number of rows and columns of \(A\) to use. If either of these is larger than the size of \(A, A\) is resized and padded with zeros.
If \(A\) is a multi-dimensional matrix, each two-dimensional sub-matrix of \(A\) is treated separately.
See also: [fft2], page 845, [ifft], page 845, [ifftn], page 846, [fftw], page 846.
fftn (A)
fftn (A, size)
Compute the N -dimensional discrete Fourier transform of \(A\) using a Fast Fourier Transform (FFT) algorithm.
The optional vector argument size may be used specify the dimensions of the array to be used. If an element of size is smaller than the corresponding dimension of \(A\), then the dimension of \(A\) is truncated prior to performing the FFT. Otherwise, if an element of size is larger than the corresponding dimension then \(A\) is resized and padded with zeros.
See also: [ifftn], page 846, [fft], page 845, [fft2], page 845, [fftw], page 846.
ifftn (A)
ifftn (A, size)
Compute the inverse N-dimensional discrete Fourier transform of \(A\) using a Fast Fourier Transform (FFT) algorithm.
The optional vector argument size may be used specify the dimensions of the array to be used. If an element of size is smaller than the corresponding dimension of \(A\), then the dimension of \(A\) is truncated prior to performing the inverse FFT. Otherwise, if an element of size is larger than the corresponding dimension then \(A\) is resized and padded with zeros.
See also: [fftn], page 846, [ifft], page 845, [ifft2], page 846, [fftw], page 846.
Octave uses the FFTW libraries to perform FFT computations. When Octave starts up and initializes the FFTW libraries, they read a system wide file (on a Unix system, it is typically /etc/fftw/wisdom) that contains information useful to speed up FFT computations. This information is called the wisdom. The system-wide file allows wisdom to be shared between all applications using the FFTW libraries.

Use the fftw function to generate and save wisdom. Using the utilities provided together with the FFTW libraries (fftw-wisdom on Unix systems), you can even add wisdom generated by Octave to the system-wide wisdom file.
```

method = fftw ("planner")
fftw ("planner", method)
wisdom = fftw ("dwisdom")
fftw ("dwisdom", wisdom)
fftw ("threads", nthreads)
nthreads = fftw ("threads")

```
    Manage FFTW wisdom data.

Wisdom data can be used to significantly accelerate the calculation of the FFTs, but implies an initial cost in its calculation. When the FFTW libraries are initialized, they read a system wide wisdom file (typically in /etc/fftw/wisdom), allowing wisdom to be shared between applications other than Octave. Alternatively, the fftw function can be used to import wisdom. For example,
```

wisdom = fftw ("dwisdom")

```
will save the existing wisdom used by Octave to the string wisdom. This string can then be saved to a file and restored using the save and load commands respectively. This existing wisdom can be re-imported as follows
```

fftw ("dwisdom", wisdom)

```

If wisdom is an empty string, then the wisdom used is cleared.
During the calculation of Fourier transforms further wisdom is generated. The fashion in which this wisdom is generated is also controlled by the fftw function. There are five different manners in which the wisdom can be treated:
```

"estimate"

```

Specifies that no run-time measurement of the optimal means of calculating a particular is performed, and a simple heuristic is used to pick a (probably sub-optimal) plan. The advantage of this method is that there is little or no overhead in the generation of the plan, which is appropriate for a Fourier transform that will be calculated once.
"measure"
In this case a range of algorithms to perform the transform is considered and the best is selected based on their execution time.
"patient"
Similar to "measure", but a wider range of algorithms is considered.
"exhaustive"
Like "measure", but all possible algorithms that may be used to treat the transform are considered.
"hybrid" As run-time measurement of the algorithm can be expensive, this is a compromise where "measure" is used for transforms up to the size of 8192 and beyond that the "estimate" method is used.

The default method is "estimate". The current method can be queried with
```

method = fftw ("planner")

```
or set by using
```

fftw ("planner", method)

```

Note that calculated wisdom will be lost when restarting Octave. However, the wisdom data can be reloaded if it is saved to a file as described above. Saved wisdom files should not be used on different platforms since they will not be efficient and the point of calculating the wisdom is lost.
The number of threads used for computing the plans and executing the transforms can be set with
```

fftw ("threads", NTHREADS)

```

Note that octave must be compiled with multi-threaded FFTW support for this feature. The number of processors available to the current process is used per default.

See also: [fft], page 845, [ifft], page 845, [fft2], page 845, [ifft2], page 846, [fftn], page 846, [ifftn], page 846.
fftconv (x, y)
fftconv ( \(x, y, n\) )
Convolve two vectors using the FFT for computation.
\(c=f f t c o n v(x, y)\) returns a vector of length equal to length ( \(x\) ) + length ( \(y\) ) -
1. If \(x\) and \(y\) are the coefficient vectors of two polynomials, the returned value is the coefficient vector of the product polynomial.
The computation uses the FFT by calling the function fftfilt. If the optional argument \(n\) is specified, an N-point FFT is used.
See also: [deconv], page 800, [conv], page 799, [conv2], page 800.
fftfilt (b, x)
fftfilt (b, x, n)
Filter \(x\) with the FIR filter \(b\) using the FFT.
If \(x\) is a matrix, filter each column of the matrix.
Given the optional third argument, \(n\), fftfilt uses the overlap-add method to filter \(x\) with \(b\) using an N-point FFT. The FFT size must be an even power of 2 and must be greater than or equal to the length of \(b\). If the specified \(n\) does not meet these criteria, it is automatically adjusted to the nearest value that does.
See also: [filter], page 848, [filter2], page 849.
```

y = filter (b, a, x)

```
[y, sf] = filter (b, a, x, si)
[y, sf] = filter (b, a, x, [], dim)
[y, sf] = filter (b, a, x, si, dim)
Apply a 1-D digital filter to the data \(x\).
filter returns the solution to the following linear, time-invariant difference equation:
\[
\sum_{k=0}^{N} a_{k+1} y_{n-k}=\sum_{k=0}^{M} b_{k+1} x_{n-k}, \quad 1 \leq n \leq P
\]
where \(a \in \Re^{N-1}, b \in \Re^{M-1}\), and \(x \in \Re^{P}\). The result is calculated over the first non-singleton dimension of \(x\) or over dim if supplied.

An equivalent form of the equation is:
\[
y_{n}=-\sum_{k=1}^{N} c_{k+1} y_{n-k}+\sum_{k=0}^{M} d_{k+1} x_{n-k}, \quad 1 \leq n \leq P
\]
where \(c=a / a_{1}\) and \(d=b / a_{1}\).
If the fourth argument si is provided, it is taken as the initial state of the system and the final state is returned as sf. The state vector is a column vector whose length is equal to the length of the longest coefficient vector minus one. If si is not supplied, the initial state vector is set to all zeros.
In terms of the Z Transform, \(y\) is the result of passing the discrete-time signal \(x\) through a system characterized by the following rational system function:
\[
H(z)=\frac{\sum_{k=0}^{M} d_{k+1} z^{-k}}{1+\sum_{k+1}^{N} c_{k+1} z^{-k}}
\]

See also: [filter2], page 849, [fftfilt], page 848, [freqz], page 849.
```

y = filter2 (b, x)
y = filter2 (b, x, shape)

```

Apply the 2-D FIR filter \(b\) to \(x\).
If the argument shape is specified, return an array of the desired shape. Possible values are:
"full" pad \(x\) with zeros on all sides before filtering.
"same" unpadded \(x\) (default)
"valid" trim \(x\) after filtering so edge effects are no included.
Note this is just a variation on convolution, with the parameters reversed and \(b\) rotated 180 degrees.

See also: [conv2], page 800 .
```

[h, w] = freqz (b, a, n, "whole")
[h, w] = freqz (b)
[h, w] = freqz (b,a)
[h, w] = freqz (b,a, n)
h = freqz (b, a, w)
[h, w] = freqz (..., Fs)
freqz (...)

```

Return the complex frequency response \(h\) of the rational IIR filter whose numerator and denominator coefficients are \(b\) and \(a\), respectively.
The response is evaluated at \(n\) angular frequencies between 0 and \(2 \pi\).
The output value \(w\) is a vector of the frequencies.

If a is omitted, the denominator is assumed to be 1 (this corresponds to a simple FIR filter).
If \(n\) is omitted, a value of 512 is assumed. For fastest computation, \(n\) should factor into a small number of small primes.
If the fourth argument, "whole", is omitted the response is evaluated at frequencies between 0 and \(\pi\).
freqz (b, a, w)
Evaluate the response at the specific frequencies in the vector \(w\). The values for \(w\) are measured in radians.
[...] = freqz (..., Fs)
Return frequencies in Hz instead of radians assuming a sampling rate Fs. If you are evaluating the response at specific frequencies \(w\), those frequencies should be requested in Hz rather than radians.
freqz (...)
Plot the magnitude and phase response of \(h\) rather than returning them.
See also: [freqz_plot], page 850 .
freqz_plot (w, h)
freqz_plot (w, h, freq_norm)
Plot the magnitude and phase response of \(h\).
If the optional freq_norm argument is true, the frequency vector \(w\) is in units of normalized radians. If freq_norm is false, or not given, then \(w\) is measured in Hertz.

See also: [freqz], page 849.
sinc ( \(x\) )
Compute the sinc function.
Return \(\sin (\pi x) /(\pi x)\).
\(b=\) unwrap ( \(x\) )
\(b=\) unwrap ( \(x\), tol)
\(b=\) unwrap (x, tol, dim)
Unwrap radian phases by adding or subtracting multiples of \(2^{*}\) pi as appropriate to remove jumps greater than tol.
tol defaults to pi.
Unwrap will work along the dimension dim. If dim is unspecified it defaults to the first non-singleton dimension.
[a, b] = arch_fit (y, x, p, iter, gamma, a0, b0)
Fit an ARCH regression model to the time series \(y\) using the scoring algorithm in Engle's original ARCH paper.
The model is
```

y(t) = b(1) * x(t,1) + ... + b(k) * x(t,k) +e(t),
h(t) = a(1) +a(2) *e(t-1)^2 + ... +a(p+1) *e(t-p)^2

```
in which \(e(t)\) is \(N(0, h(t))\), given a time-series vector y up to time \(t-1\) and a matrix of (ordinary) regressors \(x\) up to \(t\). The order of the regression of the residual variance is specified by \(p\).
If invoked as arch_fit ( \(y, k, p\) ) with a positive integer \(k\), fit an \(\operatorname{ARCH}(k, p)\) process, i.e., do the above with the \(t\)-th row of \(x\) given by
\[
[1, y(t-1), \ldots, y(t-k)]
\]

Optionally, one can specify the number of iterations iter, the updating factor gamma, and initial values \(a 0\) and \(b 0\) for the scoring algorithm.
arch_rnd \((a, b, t)\)
Simulate an ARCH sequence of length \(t\) with AR coefficients \(b\) and CH coefficients a.

The result \(y(t)\) follows the model
\(\mathrm{y}(\mathrm{t})=\mathrm{b}(1)+\mathrm{b}(2) * \mathrm{y}(\mathrm{t}-1)+\ldots+\mathrm{b}(\mathrm{lb}) * \mathrm{y}(\mathrm{t}-\mathrm{lb}+1)+\mathrm{e}(\mathrm{t})\),
where \(e(t)\), given \(y\) up to time \(t-1\), is \(N(0, h(t))\), with
\(\mathrm{h}(\mathrm{t})=\mathrm{a}(1)+\mathrm{a}(2) * \mathrm{e}(\mathrm{t}-1)^{\wedge} 2+\ldots+\mathrm{a}(\mathrm{la}) * \mathrm{e}(\mathrm{t}-\mathrm{la}+1)^{\wedge} 2\)
[pval, lm] = arch_test (y, x, p)
For a linear regression model
\(\mathrm{y}=\mathrm{x} * \mathrm{~b}+\mathrm{e}\)
perform a Lagrange Multiplier (LM) test of the null hypothesis of no conditional heteroscedascity against the alternative of \(\mathrm{CH}(p)\).
I.e., the model is
\[
y(t)=b(1) * x(t, 1)+\ldots+b(k) * x(t, k)+e(t)
\]
given \(y\) up to \(t-1\) and \(x\) up to \(t, e(\mathrm{t})\) is \(N(0, h(t))\) with
\[
h(t)=v+a(1) * e(t-1)^{\wedge} 2+\ldots+a(p) * e(t-p)^{\wedge} 2,
\]
and the null is \(a(1)==\ldots==a(p)==0\).
If the second argument is a scalar integer, \(k\), perform the same test in a linear autoregression model of order \(k\), i.e., with
\[
[1, y(t-1), \ldots, y(t-k)]
\]
as the \(t\)-th row of \(x\).
Under the null, LM approximately has a chisquare distribution with \(p\) degrees of freedom and pval is the \(p\)-value ( 1 minus the CDF of this distribution at LM) of the test.
If no output argument is given, the \(p\)-value is displayed.
arma_rnd ( \(a, b, v, t, n)\)
Return a simulation of the ARMA model.
The ARMA model is defined by
```

x(n) = a(1) * x (n-1) + ... + a(k) * x (n-k)
+e(n) + b(1) * e(n-1) + ... + b(l) * e(n-l)

```
in which \(k\) is the length of vector a, \(l\) is the length of vector \(b\) and \(e\) is Gaussian white noise with variance \(v\). The function returns a vector of length \(t\).

The optional parameter \(n\) gives the number of dummy \(x(i)\) used for initialization, i.e., a sequence of length \(t+n\) is generated and \(x(n+1: t+n)\) is returned. If \(n\) is omitted, \(n\) \(=100\) is used.
autoreg_matrix (y, k)
Given a time series (vector) \(y\), return a matrix with ones in the first column and the first \(k\) lagged values of \(y\) in the other columns.
In other words, for \(t>k,[1, y(t-1), \ldots, y(t-k)]\) is the \(t\)-th row of the result. The resulting matrix may be used as a regressor matrix in autoregressions.
bartlett (m)
Return the filter coefficients of a Bartlett (triangular) window of length m.
For a definition of the Bartlett window see, e.g., A.V. Oppenheim \& R. W. Schafer, Discrete-Time Signal Processing.
blackman (m)
blackman (m, "periodic")
blackman (m, "symmetric")
Return the filter coefficients of a Blackman window of length \(m\).
If the optional argument "periodic" is given, the periodic form of the window is returned. This is equivalent to the window of length \(m+1\) with the last coefficient removed. The optional argument "symmetric" is equivalent to not specifying a second argument.
For a definition of the Blackman window, see, e.g., A.V. Oppenheim \& R. W. Schafer, Discrete-Time Signal Processing.
detrend ( \(x, p\) )
If \(x\) is a vector, detrend \((x, p)\) removes the best fit of a polynomial of order \(p\) from the data \(x\).
If \(x\) is a matrix, detrend ( \(x, p\) ) does the same for each column in \(x\).
The second argument \(p\) is optional. If it is not specified, a value of 1 is assumed. This corresponds to removing a linear trend.
The order of the polynomial can also be given as a string, in which case \(p\) must be either "constant" (corresponds to \(p=0\) ) or "linear" (corresponds to \(p=1\) ).
See also: [polyfit], page 803.
[d, dd] = diffpara (x, a, b)
Return the estimator \(d\) for the differencing parameter of an integrated time series.
The frequencies from \([2 * p i * a / t, 2 * p i * b / T]\) are used for the estimation. If \(b\) is omitted, the interval \([2 * p i / T, 2 * p i * a / T]\) is used. If both \(b\) and a are omitted then \(a=0.5 * \operatorname{sqrt}(T)\) and \(b=1.5 * \operatorname{sqrt}(T)\) is used, where \(T\) is the sample size. If x is a matrix, the differencing parameter of each column is estimated.
The estimators for all frequencies in the intervals described above is returned in \(d d\). The value of \(d\) is simply the mean of \(d d\).
Reference: P.J. Brockwell \& R.A. Davis. Time Series: Theory and Methods. Springer 1987.
durbinlevinson (c, oldphi, oldv)
Perform one step of the Durbin-Levinson algorithm.
The vector \(c\) specifies the autocovariances [gamma_0, ..., gamma_t] from lag 0 to \(t\), oldphi specifies the coefficients based on \(c(t-1)\) and oldv specifies the corresponding error.
If oldphi and oldv are omitted, all steps from 1 to \(t\) of the algorithm are performed.
fftshift (x)
fftshift (x, dim)
Perform a shift of the vector \(x\), for use with the fft and ifft functions, in order to move the frequency 0 to the center of the vector or matrix.
If x is a vector of \(N\) elements corresponding to \(N\) time samples spaced by \(d t\), then fftshift (fft (x)) corresponds to frequencies
\(\mathrm{f}=[-(\) ceil \(((\mathrm{N}-1) / 2):-1: 1), 0,(1: f l o o r((N-1) / 2))] * d f\)
where \(d f=1 /(N * d t)\).
If \(x\) is a matrix, the same holds for rows and columns. If \(x\) is an array, then the same holds along each dimension.
The optional dim argument can be used to limit the dimension along which the permutation occurs.

See also: [ifftshift], page 853.
ifftshift (x)
ifftshift (x, dim)
Undo the action of the fftshift function.
For even length x , fftshift is its own inverse, but odd lengths differ slightly.
See also: [fftshift], page 853.
fractdiff ( \(x, d\) )
Compute the fractional differences \((1-L)^{d} x\) where \(L\) denotes the lag-operator and \(d\) is greater than -1 .
hamming ( \(m\) )
hamming ( \(m\), "periodic")
hamming (m, "symmetric")
Return the filter coefficients of a Hamming window of length \(m\).
If the optional argument "periodic" is given, the periodic form of the window is returned. This is equivalent to the window of length \(m+1\) with the last coefficient removed. The optional argument "symmetric" is equivalent to not specifying a second argument.
For a definition of the Hamming window see, e.g., A.V. Oppenheim \& R. W. Schafer, Discrete-Time Signal Processing.
hanning ( \(m\) )
hanning ( \(m\), "periodic")
hanning ( \(m\), "symmetric")
Return the filter coefficients of a Hanning window of length \(m\).

If the optional argument "periodic" is given, the periodic form of the window is returned. This is equivalent to the window of length \(m+1\) with the last coefficient removed. The optional argument "symmetric" is equivalent to not specifying a second argument.
For a definition of the Hanning window see, e.g., A.V. Oppenheim \& R. W. Schafer, Discrete-Time Signal Processing.
hurst (x)
Estimate the Hurst parameter of sample \(x\) via the rescaled range statistic.
If \(x\) is a matrix, the parameter is estimated for every column.
\(p p=\operatorname{pchip}(x, y)\)
yi = pchip (x, y, xi)
Return the Piecewise Cubic Hermite Interpolating Polynomial (pchip) of points \(x\) and \(y\).
If called with two arguments, return the piecewise polynomial \(p p\) that may be used with ppval to evaluate the polynomial at specific points.
When called with a third input argument, pchip evaluates the pchip polynomial at the points xi. The third calling form is equivalent to ppval (pchip ( \(x, y\) ), xi).
The variable \(x\) must be a strictly monotonic vector (either increasing or decreasing) of length \(n\).
\(y\) can be either a vector or array. If \(y\) is a vector then it must be the same length \(n\) as \(x\). If \(y\) is an array then the size of \(y\) must have the form
\[
\left[s_{1}, s_{2}, \cdots, s_{k}, n\right]
\]

The array is reshaped internally to a matrix where the leading dimension is given by
\[
s_{1} s_{2} \cdots s_{k}
\]
and each row of this matrix is then treated separately. Note that this is exactly opposite to interp1 but is done for MATLAB compatibility.
See also: [spline], page 818, [ppval], page 812, [mkpp], page 811, [unmkpp], page 811.
[Pxx, w] = periodogram (x)
[Pxx, w] = periodogram (x, win)
[Pxx, w] = periodogram (x, win, nfft)
[Pxx, f] = periodogram (x, win, nfft, Fs)
[Pxx, f] = periodogram (..., "range")
periodogram (...)
Return the periodogram (Power Spectral Density) of \(x\).
The possible inputs are:
\(x\)
data vector. If \(x\) is real-valued a one-sided spectrum is estimated. If \(x\) is complex-valued, or "range" specifies "twosided", the full spectrum is estimated.
win window weight data. If window is empty or unspecified a default rectangular window is used. Otherwise, the window is applied to the signal ( \(x\) .* win) before computing the periodogram. The window data must be a vector of the same length as \(x\).
nfft number of frequency bins. The default is 256 or the next higher power of 2 greater than the length of \(x(\max (256,2 . \wedge\) nextpow2 (length (x))) ). If \(n f f t\) is greater than the length of the input then \(x\) will be zero-padded to the length of \(n f f t\).
Fs sampling rate. The default is 1.
range range of spectrum. "onesided" computes spectrum from [0:nfft/2+1]. "twosided" computes spectrum from [0:nfft-1].
The optional second output \(w\) are the normalized angular frequencies. For a one-sided calculation \(w\) is in the range \([0, \mathrm{pi}]\) if \(n f f t\) is even and \([0, \mathrm{pi})\) if \(n f f t\) is odd. Similarly, for a two-sided calculation \(w\) is in the range \(\left[0,2^{*} \mathrm{pi}\right]\) or \(\left[0,2^{*} \mathrm{pi}\right)\) depending on \(n f f t\).
If a sampling frequency is specified, \(F s\), then the output frequencies \(f\) will be in the range \(\left[0, F_{s} / 2\right]\) or \(\left[0, F_{s} / 2\right)\) for one-sided calculations. For two-sided calculations the range will be \([0, F s)\).
When called with no outputs the periodogram is immediately plotted in the current figure window.
See also: [fft], page 845.
sinetone (freq, rate, sec, ampl)
Return a sinetone of frequency freq with a length of sec seconds at sampling rate rate and with amplitude ampl.
The arguments freq and ampl may be vectors of common size.
The defaults are rate \(=8000, \mathrm{sec}=1\), and ampl \(=64\).
See also: [sinewave], page 855.
sinewave ( \(m, n, d\) )
Return an m-element vector with \(i\)-th element given by \(\sin (2 * \mathrm{pi} *(i+d-1) / n)\). The default value for \(d\) is 0 and the default value for \(n\) is \(m\).

See also: [sinetone], page 855.
spectral_adf (c)
spectral_adf (c, win)
spectral_adf (c, win, b)
Return the spectral density estimator given a vector of autocovariances \(c\), window name win, and bandwidth, \(b\).
The window name, e.g., "triangle" or "rectangle" is used to search for a function called win_lw.
If win is omitted, the triangle window is used.
If \(b\) is omitted, \(1 /\) sqrt (length (x)) is used.
See also: [spectral_xdf], page 856 .
spectral_xdf (x)
spectral_xdf (x, win)
spectral_xdf (x, win, b)
Return the spectral density estimator given a data vector \(x\), window name win, and bandwidth, \(b\).
The window name, e.g., "triangle" or "rectangle" is used to search for a function called win_sw.
If win is omitted, the triangle window is used.
If \(b\) is omitted, \(1 /\) sqrt (length ( \(x\) ) ) is used.
See also: [spectral_adf], page 855 .
spencer ( \(x\) )
Return Spencer's 15 -point moving average of each column of \(x\).
\(y=\operatorname{stft}(x)\)
\(y=s t f t(x\), win_size)
\(y=\operatorname{stft}(x\), win_size, inc)
\(y=\) stft (x, win_size, inc, num_coef)
\(y=\) stft ( \(x\), win_size, inc, num_coef, win_type)
[y, c] = stft (...)
Compute the short-time Fourier transform of the vector \(x\) with num_coef coefficients by applying a window of win_size data points and an increment of inc points.
Before computing the Fourier transform, one of the following windows is applied:
```

"hanning"
win_type $=1$
"hamming"
win_type $=2$
"rectangle"
win_type $=3$

```

The window names can be passed as strings or by the win_type number.
The following defaults are used for unspecified arguments: win_size \(=80\), inc \(=24\), num_coef \(=64\), and win_type \(=1\).
\(y=\operatorname{stft}(x, \ldots)\) returns the absolute values of the Fourier coefficients according to the num_coef positive frequencies.
\([y, c]=\operatorname{stft}(x, \ldots)\) returns the entire STFT-matrix \(y\) and a 3-element vector \(c\) containing the window size, increment, and window type, which is needed by the synthesis function.
See also: [synthesis], page 856.
\(x=\) synthesis ( \(y, c\) )
Compute a signal from its short-time Fourier transform \(y\) and a 3-element vector \(c\) specifying window size, increment, and window type.
The values \(y\) and \(c\) can be derived by
\([y, c]=\operatorname{stft}(x, \ldots)\)
See also: [stft], page 856 .
[a, v] = yulewalker (c)
Fit an AR (p)-model with Yule-Walker estimates given a vector \(c\) of autocovariances [gamma_0, ..., gamma_p].
Returns the AR coefficients, a, and the variance of white noise, \(v\).

\section*{32 Image Processing}

Since an image is basically a matrix, Octave is a very powerful environment for processing and analyzing images. To illustrate how easy it is to do image processing in Octave, the following example will load an image, smooth it by a 5 -by- 5 averaging filter, and compute the gradient of the smoothed image.
```

I = imread ("myimage.jpg");
S = conv2 (I, ones (5, 5) / 25, "same");
[Dx, Dy] = gradient (S);

```

In this example \(S\) contains the smoothed image, and Dx and Dy contains the partial spatial derivatives of the image.

\subsection*{32.1 Loading and Saving Images}

The first step in most image processing tasks is to load an image into Octave which is done with the imread function. The imwrite function is the corresponding function for writing images to the disk.

In summary, most image processing code will follow the structure of this code
```

    I = imread ("my_input_image.img");
    J = process_my_image (I);
    imwrite (J, "my_output_image.img");
    [img, map, alpha] = imread (filename)
[...] = imread (url)
[...] = imread (..., ext)
[...] = imread (..., idx)
[...] = imread (..., param1, value1,...)
Read images from various file formats.

```

Read an image as a matrix from the file filename or from the online resource url. If neither is given, but ext was specified, look for a file with the extension ext.
The size and class of the output depends on the format of the image. A color image is returned as an MxNx3 matrix. Grayscale and black-and-white images are of size MxN. Multipage images will have an additional 4th dimension.
The bit depth of the image determines the class of the output: "uint8", "uint16", or "single" for grayscale and color, and "logical" for black-and-white. Note that indexed images always return the indexes for a colormap, independent of whether map is a requested output. To obtain the actual RGB image, use ind2rgb. When more than one indexed image is being read, map is obtained from the first. In some rare cases this may be incorrect and imfinfo can be used to obtain the colormap of each image.
See the Octave manual for more information in representing images. (see Section 32.3 [Representing Images], page 867)
Some file formats, such as TIFF and GIF, are able to store multiple images in a single file. idx can be a scalar or vector specifying the index of the images to read. By default, Octave will read only the first page.

Depending on the file format, it is possible to configure the reading of images with parameter, value pairs. The following options are supported:
"Frames" or "Index"
This is an alternative method to specify idx. When specifying it in this way, its value can also be the string "all".
"Info" This option exists for matlab compatibility, but has no effect. For maximum performance when reading multiple images from a single file, use the "Index" option.
"PixelRegion"
Controls the image region that is read. The value must be a cell array with two arrays of 3 elements \{[rows], [cols]\}. The elements in the array are the start, increment, and end pixel to be read. If the increment value is omitted it defaults to 1 . For example, the following are all equivalent:
```

imread (filename, "PixelRegion", {[200 600], [300 700]});
imread (filename, "PixelRegion", {[[200 1 600], [300 1 700]});
imread (filename)(200:600, 300:700);

```

See also: [imwrite], page 860, [imfinfo], page 862, [imformats], page 864.
```

imwrite (img, filename)
imwrite (img, filename, ext)
imwrite (img, map, filename)
imwrite (..., param1, val1,...)

```

Write images in various file formats.
The image img can be a binary, grayscale, RGB, or multi-dimensional image. The size and class of img should be the same as what should be expected when reading it with imread: the 3rd and 4th dimensions reserved for color space, and multiple pages respectively. If it's an indexed image, the colormap map must also be specified.

If ext is not supplied, the file extension of filename is used to determine the format. The actual supported formats are dependent on options made during the build of Octave. Use imformats to check the support of the different image formats.
Depending on the file format, it is possible to configure the writing of images with param, val pairs. The following options are supported:
'Alpha' Alpha (transparency) channel for the image. This must be a matrix with same class, and number of rows and columns of img. In case of a multipage image, the size of the 4th dimension must also match and the third dimension must be a singleton. By default, image will be completely opaque.
'Compression'
Compression to use one the image. Can be one of the following: "none" (default), "bzip", "fax3", "fax4", "jpeg", "lzw", "rle", or "deflate". Note that not all compression types are available for all image formats in which it defaults to your Magick library.
'DelayTime'
For formats that accept animations (such as GIF), controls for how long a frame is displayed until it moves to the next one. The value must be scalar (which will applied to all frames in img), or a vector of length equal to the number of frames in im. The value is in seconds, must be between 0 and 655.35 , and defaults to 0.5 .
'DisposalMethod'
For formats that accept animations (such as GIF), controls what happens to a frame before drawing the next one. Its value can be one of the following strings: "doNotSpecify" (default); "leaveInPlace"; "restoreBG"; and "restorePrevious", or a cell array of those string with length equal to the number of frames in img.
'LoopCount'
For formats that accept animations (such as GIF), controls how many times the sequence is repeated. A value of Inf means an infinite loop (default), a value of 0 or 1 that the sequence is played only once (loops zero times), while a value of 2 or above loops that number of times (looping twice means it plays the complete sequence 3 times). This option is ignored when there is only a single image at the end of writing the file.
'Quality' Set the quality of the compression. The value should be an integer between 0 and 100, with larger values indicating higher visual quality and lower compression. Defaults to 75 .
'WriteMode'
Some file formats, such as TIFF and GIF, are able to store multiple images in a single file. This option specifies if img should be appended to the file (if it exists) or if a new file should be created for it (possibly overwriting an existing file). The value should be the string "Overwrite" (default), or "Append".
Despite this option, the most efficient method of writing a multipage image is to pass a 4 dimensional img to imwrite, the same matrix that could be expected when using imread with the option "Index" set to "all".

See also: [imread], page 859, [imfinfo], page 862, [imformats], page 864.
val = IMAGE_PATH ()
old_val = IMAGE_PATH (new_val)
IMAGE_PATH (new_val, "local")
Query or set the internal variable that specifies a colon separated list of directories in which to search for image files.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [EXEC_PATH], page 976, [OCTAVE_HOME], page 987, [OCTAVE_EXEC_HOME], page 987.

It is possible to get information about an image file on disk, without actually reading it into Octave. This is done using the imfinfo function which provides read access to many of the parameters stored in the header of the image file.
```

info = imfinfo (filename)
info = imfinfo (url)
info = imfinfo (..., ext)

```

Read image information from a file.
imfinfo returns a structure containing information about the image stored in the file filename. If there is no file filename, and ext was specified, it will look for a file named filename and extension ext, i.e., a file named filename.ext.
The output structure info contains the following fields:
'Filename'
The full name of the image file.
'FileModDate'
Date of last modification to the file.
'FileSize'
Number of bytes of the image on disk
'Format' Image format (e.g., "jpeg").
'Height' Image height in pixels.
'Width' Image Width in pixels.
'BitDepth'
Number of bits per channel per pixel.
'ColorType'
Image type. Value is "grayscale", "indexed", "truecolor", "CMYK", or "undefined".
'XResolution'
X resolution of the image.
'YResolution'
Y resolution of the image.
'ResolutionUnit'
Units of image resolution. Value is "Inch", "Centimeter", or
"undefined".
'DelayTime'
Time in \(1 / 100\) ths of a second ( 0 to 65535 ) which must expire before displaying the next image in an animated sequence.
'LoopCount'
Number of iterations to loop an animation.
'ByteOrder'
Endian option for formats that support it. Value is "little-endian",
"big-endian", or "undefined".
'Gamma' Gamma level of the image. The same color image displayed on two different workstations may look different due to differences in the display monitor.
'Quality' JPEG/MIFF/PNG compression level. Value is an integer in the range [0 100].
'DisposalMethod'
Only valid for GIF images, control how successive frames are rendered (how the preceding frame is disposed of) when creating a GIF animation. Values can be "doNotSpecify", "leaveInPlace", "restoreBG", or "restorePrevious". For non-GIF files, value is an empty string.
'Chromaticities'
Value is a 1 x 8 Matrix with the \(\mathrm{x}, \mathrm{y}\) chromaticity values for white, red, green, and blue points, in that order.
'Comment' Image comment.
'Compression'
Compression type. Value can be "none", "bzip", "fax3", "fax4", "jpeg", "lzw", "rle", "deflate", "lzma", "jpeg2000", "jbig2", "jbig2", or "undefined".
'Colormap'
Colormap for each image.
'Orientation'
The orientation of the image with respect to the rows and columns. Value is an integer between 1 and 8 as defined in the TIFF 6 specifications, and for matlab compatibility.
'Software'
Name and version of the software or firmware of the camera or image input device used to generate the image.
'Make' The manufacturer of the recording equipment. This is the manufacture of the DSC, scanner, video digitizer or other equipment that generated the image.
'Model' The model name or model number of the recording equipment as mentioned on the field "Make".
'DateTime'
The date and time of image creation as defined by the Exif standard, i.e., it is the date and time the file was changed.
'ImageDescription'
The title of the image as defined by the Exif standard.
'Artist' Name of the camera owner, photographer or image creator.
'Copyright'
Copyright notice of the person or organization claiming rights to the image.
'DigitalCamera'
A struct with information retrieved from the Exif tag.
'GPSInfo' A struct with geotagging information retrieved from the Exif tag.
See also: [imread], page 859, [imwrite], page 860, [imshow], page 865, [imformats], page 864 .

By default, Octave's image IO functions (imread, imwrite, and imfinfo) use the GraphicsMagick library for their operations. This means a vast number of image formats is supported but considering the large amount of image formats in science and its commonly closed nature, it is impossible to have a library capable of reading them all. Because of this, the function imformats keeps a configurable list of available formats, their extensions, and what functions should the image IO functions use. This allows one to expand Octave's image IO capabilities by creating functions aimed at acting on specific file formats.

While it would be possible to call the extra functions directly, properly configuring Octave with imformats allows one to keep a consistent code that is abstracted from file formats.

It is important to note that a file format is not actually defined by its file extension and that GraphicsMagick is capable to read and write more file formats than the ones listed by imformats. What this means is that even with an incorrect or missing extension the image may still be read correctly, and that even unlisted formats are not necessarily unsupported.
```

imformats ()
formats = imformats (ext)
formats = imformats (format)
formats = imformats ("add", format)
formats = imformats ("remove", ext)
formats = imformats ("update", ext, format)
formats = imformats ("factory")
Manage supported image formats.

```
formats is a structure with information about each supported file format, or from a specific format ext, the value displayed on the field ext. It contains the following fields:
ext The name of the file format. This may match the file extension but Octave will automatically detect the file format.
description
A long description of the file format.
isa A function handle to confirm if a file is of the specified format.
write A function handle to write if a file is of the specified format.
read A function handle to open files the specified format.
info A function handle to obtain image information of the specified format.
alpha Logical value if format supports alpha channel (transparency or matte).
multipage Logical value if format supports multipage (multiple images per file).

It is possible to change the way Octave manages file formats with the options "add", "remove", and "update", and supplying a structure format with the required fields. The option "factory" resets the configuration to the default.
This can be used by Octave packages to extend the image reading capabilities Octave, through use of the PKG_ADD and PKG_DEL commands.
See also: [imfinfo], page 862, [imread], page 859, [imwrite], page 860.

\subsection*{32.2 Displaying Images}

A natural part of image processing is visualization of an image. The most basic function for this is the imshow function that shows the image given in the first input argument.
```

imshow (im)
imshow (im, limits)
imshow (im, map)
imshow (rgb,...)
imshow (filename)
imshow (..., string_param1, value1,...)
h = imshow (...)
Display the image im, where im can be a 2-dimensional (grayscale image) or a 3-
dimensional (RGB image) matrix.

```

If limits is a 2 -element vector [low, high], the image is shown using a display range between low and high. If an empty matrix is passed for limits, the display range is computed as the range between the minimal and the maximal value in the image.
If map is a valid color map, the image will be shown as an indexed image using the supplied color map.
If a filename is given instead of an image, the file will be read and shown.
If given, the parameter string_param1 has value value1. string_param1 can be any of the following:
"displayrange"
value1 is the display range as described above.
"colormap"
value1 is the colormap to use when displaying an indexed image.
"xdata" If value1 is a 2-element vector, it must contain horizontal image limits in the form [xfirst, xlast], where xfirst and xlast are the abscissa of the centers of the corner pixels. Otherwise value1 must be a vector and only the first and last elements will be used for xfirst and xlast respectively.
"ydata" If value1 is a 2-element vector, it must contain vertical image limits in the form [yfirst, ylast], where yfirst and ylast are the ordinates of the center of the corner pixels. Otherwise value1 must be a vector and only the first and last elements will be used for yfirst and ylast respectively.

The optional return value \(h\) is a graphics handle to the image.
See also: [image], page 866, [imagesc], page 866, [colormap], page 870, [gray2ind], page 868, [rgb2ind], page 868.
```

image (img)
image (x, y, img)
image (..., "prop", val, ...)
image ("prop1", val1,...)
h = image (...)
Display a matrix as an indexed color image.

```

The elements of img are indices into the current colormap.
\(x\) and \(y\) are optional 2-element vectors, [min, max], which specify the coordinates of the centers of the corner pixels. If a range is specified as [max, min] then the image will be reversed along that axis. For convenience, \(x\) and \(y\) may be specified as N -element vectors matching the length of the data in img. However, only the first and last elements will be used to determine the axis limits.

Multiple property/value pairs may be specified for the image object, but they must appear in pairs.

The optional return value \(h\) is a graphics handle to the image.
Implementation Note: The origin \((0,0)\) for images is located in the upper left. For ordinary plots, the origin is located in the lower left. Octave handles this inversion by plotting the data normally, and then reversing the direction of the \(y\)-axis by setting the ydir property to "reverse". This has implications whenever an image and an ordinary plot need to be overlaid. The recommended solution is to display the image and then plot the reversed ydata using, for example, flipud (ydata).
Calling Forms: The image function can be called in two forms: High-Level and LowLevel. When invoked with normal options, the High-Level form is used which first calls newplot to prepare the graphic figure and axes. When the only inputs to image are property/value pairs the Low-Level form is used which creates a new instance of an image object and inserts it in the current axes.

Graphic Properties: The full list of properties is documented at Section 15.3.3.7 [Image Properties], page 455.

See also: [imshow], page 865, [imagesc], page 866, [colormap], page 870.
```

imagesc (img)
imagesc (x, y, img)
imagesc (..., climits)
imagesc (..., "prop",val, ...)
imagesc ("prop1", val1,...)
imagesc (hax,...)
h = imagesc (...)

```

Display a scaled version of the matrix img as a color image.
The colormap is scaled so that the entries of the matrix occupy the entire colormap. If climits \(=[10, h i]\) is given, then that range is set to the "clim" of the current axes.
\(x\) and \(y\) are optional 2-element vectors, [min, max], which specify the coordinates of the centers of the corner pixels. If a range is specified as [max, min] then the image will be reversed along that axis. For convenience, \(x\) and \(y\) may be specified
as N -element vectors matching the length of the data in img. However, only the first and last elements will be used to determine the image limits.
The optional return value \(h\) is a graphics handle to the image.
Calling Forms: The imagesc function can be called in two forms: High-Level and Low-Level. When invoked with normal options, the High-Level form is used which first calls newplot to prepare the graphic figure and axes. When the only inputs to image are property/value pairs the Low-Level form is used which creates a new instance of an image object and inserts it in the current axes. The full list of properties is documented at Section 15.3.3.7 [Image Properties], page 455.
See also: [image], page 866, [imshow], page 865, [caxis], page 350.

\subsection*{32.3 Representing Images}

In general Octave supports four different kinds of images, grayscale images, RGB images, binary images, and indexed images. A grayscale image is represented with an M-by-N matrix in which each element corresponds to the intensity of a pixel. An RGB image is represented with an M-by-N-by-3 array where each 3 -vector corresponds to the red, green, and blue intensities of each pixel.

The actual meaning of the value of a pixel in a grayscale or RGB image depends on the class of the matrix. If the matrix is of class double pixel intensities are between 0 and 1 , if it is of class uint8 intensities are between 0 and 255 , and if it is of class uint16 intensities are between 0 and 65535 .

A binary image is an M-by-N matrix of class logical. A pixel in a binary image is black if it is false and white if it is true.

An indexed image consists of an M-by-N matrix of integers and a C-by-3 color map. Each integer corresponds to an index in the color map, and each row in the color map corresponds to an RGB color. The color map must be of class double with values between 0 and 1.

The following convenience functions are available for conversion between image formats.
```

im2double (img)
im2double (img, "indexed")

```

Convert image to double precision.
The conversion of img to double precision, is dependent on the type of input image. The following input classes are supported:
'uint8, uint16, and int16'
The range of values from the class is scaled to the interval [01].
'logical' True and false values are assigned a value of 0 and 1 respectively.
'single' Values are cast to double.
'double' Returns the same image.
If img is an indexed image, then the second argument should be the string "indexed". If so, then img must either be of floating point class, or unsigned integer class and it will simply be cast to double. If it is an integer class, \(a+1\) offset is applied.
See also: [double], page 51.
img = gray2ind (I)
img = gray2ind (I, n)
img = gray2ind (BW)
img = gray2ind (BW, n)
[img, map] = gray2ind (...)
Convert a grayscale or binary intensity image to an indexed image.
The indexed image will consist of \(n\) different intensity values. If not given \(n\) defaults to 64 for grayscale images or 2 for binary black and white images.
The output img is of class uint 8 if \(n\) is less than or equal to 256 ; Otherwise the return class is uint16.
See also: [ind2gray], page 868, [rgb2ind], page 868.
\(I\) = ind2gray (x, map)
Convert a color indexed image to a grayscale intensity image.
The image \(x\) must be an indexed image which will be converted using the colormap map. If map does not contain enough colors for the image, pixels in \(x\) outside the range are mapped to the last color in the map before conversion to grayscale.
The output \(I\) is of the same class as the input \(x\) and may be one of uint8, uint16, single, or double.
Implementation Note: There are several ways of converting colors to grayscale intensities. This functions uses the luminance value obtained from rgb2gray which is \(\mathrm{I}=\) \(0.299 * R+0.587 * G+0.114 * B\). Other possibilities include the value component from rgb2hsv or using a single color channel from ind2rgb.
See also: [gray2ind], page 868, [ind2rgb], page 868.
[x, map] = rgb2ind (rgb)
[x, map] = rgb2ind ( \(R, G, B\) )
Convert an image in red-green-blue (RGB) color space to an indexed image.
The input image rgb can be specified as a single matrix of size MxNx 3 , or as three separate variables, \(R, G\), and \(B\), its three color channels, red, green, and blue.
It outputs an indexed image x and a colormap map to interpret an image exactly the same as the input. No dithering or other form of color quantization is performed. The output class of the indexed image \(x\) can be uint8, uint16 or double, whichever is required to specify the number of unique colors in the image (which will be equal to the number of rows in map) in order.
Multi-dimensional indexed images (of size MxNx3xK) are also supported, both via a single input (rgb) or its three color channels as separate variables.
See also: [ind2rgb], page 868, [rgb2hsv], page 879, [rgb2gray], page 879.
rgb \(=\) ind2rgb ( \(x\), map)
\([R, G, B]=\operatorname{ind} 2 r g b\) (x, map)
Convert an indexed image to red, green, and blue color components.
The image \(x\) must be an indexed image which will be converted using the colormap map. If map does not contain enough colors for the image, pixels in \(x\) outside the range are mapped to the last color in the map.

The output may be a single RGB image ( MxNx 3 matrix where M and N are the original image \(x\) dimensions, one for each of the red, green and blue channels). Alternatively, the individual red, green, and blue color matrices of size MxN may be returned.
Multi-dimensional indexed images (of size MxNx1xK) are also supported.
See also: [rgb2ind], page 868, [ind2gray], page 868, [hsv2rgb], page 879.
Octave also provides tools to produce and work with movie frame structures. Those structures encapsulate the image data ("cdata" field) together with the corresponding colormap ("colormap" field).
frame = getframe ()
frame \(=\) getframe (hax)
frame = getframe (hfig)
frame \(=\) getframe (..., rect)
Capture a figure or axes as a movie frame structure.
Without an argument, capture the current axes excluding ticklabels, title, and \(\mathrm{x} / \mathrm{y} / \mathrm{zlabels}\). The returned structure frame has a field cdata, which contains the actual image data in the form of an NxMx3 (RGB) uint8 matrix, and a field colormap which is provided for matlab compatibility but is always empty.
If the first argument hax is an axes handle, then capture this axes, rather than the current axes returned by gca.
If the first argument hfig is a figure handle then the entire corresponding figure canvas is captured.
Finally, if a second argument rect is provided it must be a four-element vector ([left bottom width height]) defining the region inside the figure to be captured. Regardless of the figure "units" property, rect must be defined in pixels.
See also: [im2frame], page 870, [frame2im], page 870, [movie], page 869.
movie (mov)
movie (mov, \(n\) )
movie (mov, n, fps)
movie (h, ...)
Play a movie defined by an array of frame structures.
The movie mov must be a struct array of frames with fields "cdata" and "colormap", as returned by the getframe function. By default all images are displayed once, at 12 fps , in the current axes.
The optional argument \(n\) is a scalar or vector of integers that controls the number of times the movie is displayed and which particular frames are shown:
First element:
\(n(1)>0 \quad\) Play the movie \(n(1)\) times.
\(n(1)<0 \quad\) Play the movie abs ( \(n(1)\) times alternatively in forward and backward order.
Other elements (if any):
Indices of the frames in mov that will be displayed.

If the first argument is a handle to a figure or axes \(h\), the movie is played in that figure or axes instead of the current axes.
See also: [getframe], page 869, [im2frame], page 870, [frame2im], page 870.
[x, map] = frame2im (frame)
Convert movie frame to indexed image.
A movie frame is simply a struct with the fields "cdata" and "colormap".
Support for N-dimensional images or movies is given when frame is a struct array. In such cases, \(x\) will be a \(M x N x 1 x K\) or \(M x N x 3 x K\) for indexed and RGB movies respectively, with each frame concatenated along the 4th dimension.

See also: [im2frame], page 870, [getframe], page 869.
im2frame (rgb)
im2frame (x, map)
Convert image to movie frame.
A movie frame is simply a struct with the fields "cdata" and "colormap".
Support for N -dimensional images is given when each image projection, matrix sizes of MxN and MxNx 3 for RGB images, is concatenated along the fourth dimension. In such cases, the returned value is a struct array.
See also: [frame2im], page 870.
The colormap function is used to change the colormap of the current axes or figure.
```

cmap = colormap ()
cmap = colormap (map)
cmap = colormap ("default")
cmap = colormap (map_name)
cmap = colormap (hax,...)
colormap map_name

```

Query or set the current colormap.
With no input arguments, colormap returns the current color map.
colormap (map) sets the current colormap to map. The colormap should be an \(n\) row by 3 column matrix. The columns contain red, green, and blue intensities respectively. All entries must be between 0 and 1 inclusive. The new colormap is returned.
colormap ("default") restores the default colormap (the viridis map with 64 entries). The default colormap is returned.

The map may also be specified by a string, map_name, which is the name of a function that returns a colormap.
If the first argument hax is an axes handle, then the colormap for those axes is queried or set.
For convenience, it is also possible to use this function with the command form, colormap map_name.
The list of built-in colormaps is:

colorcube equally spaced colors in RGB color space.
flag cyclic 4-color map of red, white, blue, black.
lines
prism
cyclic colormap with colors from axes "ColorOrder" property.
cyclic 6 -color map of red, orange, yellow, green, blue, violet.
white all white colormap (no colors).
See also: [viridis], page 876, [turbo], page 875, [jet], page 874, [cubehelix], page 873, [hsv], page 874, [rainbow], page 875, [hot], page 874, [cool], page 873, [spring], page 875, [summer], page 875, [autumn], page 872, [winter], page 876, [gray], page 873, [bone], page 872, [copper], page 873, [pink], page 875, [ocean], page 874, [colorcube], page 872, [flag], page 873, [lines], page 874, [prism], page 875, [white], page 876.

\section*{iscolormap (cmap)}

Return true if cmap is a colormap.
A colormap is a real matrix, of class single or double, with 3 columns. Each row represents a single color. The 3 columns contain red, green, and blue intensities respectively.
All values in a colormap should be in the [01] range but this is not enforced. Each function must decide what to do for values outside this range.

See also: [colormap], page 870, [rgbplot], page 872.
The following functions return predefined colormaps, the same that can be requested by name using the colormap function.
rgbplot (cmap)
rgbplot (cmap, style)
\(h=r g b p l o t(. .\).
Plot the components of a colormap.
Two different styles are available for displaying the cmap:
profile (default)
Plot the RGB line profile of the colormap for each of the channels (red, green and blue) with the plot lines colored appropriately. Each line represents the intensity of an RGB component across the colormap.
composite Draw the colormap across the X-axis so that the actual index colors are visible rather than the individual color components.

The optional return value \(h\) is a graphics handle to the created plot.
Run demo rgbplot to see an example of rgbplot and each style option.
See also: [colormap], page 870 .
map \(=\) autumn ()
map \(=\) autumn (n)
Create color colormap. This colormap ranges from red through orange to yellow.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
map \(=\) bone ()
map \(=\) bone ( \(n\) )
Create color colormap. This colormap varies from black to white with gray-blue shades.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
map = colorcube ()
map \(=\) colorcube ( \(n\) )
Create color colormap. This colormap is composed of as many equally spaced colors (not grays) in the RGB color space as possible.
If there are not a perfect number \(n\) of regularly spaced colors then the remaining entries in the colormap are gradients of pure red, green, blue, and gray.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
```

map = cool ()

```
\(\operatorname{map}=\operatorname{cool}(n)\)

Create color colormap. The colormap varies from cyan to magenta.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
map \(=\) copper ()
map \(=\operatorname{copper}(n)\)
Create color colormap. This colormap varies from black to a light copper tone.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map = cubehelix ()
map \(=\) cubehelix ( \(n\) )
map \(=\) cubehelix (n, start, rots, hue, gamma)
Create cubehelix colormap.
This colormap varies from black to white going though blue, green, and red tones while maintaining a monotonically increasing perception of intensity. This is achieved by traversing a color cube from black to white through a helix, hence the name cubehelix, while taking into account the perceived brightness of each channel according to the NTSC specifications from 1953.
```

rgbplot (cubehelix (256))

```

The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
Reference: Green, D. A., 2011, A colour scheme for the display of astronomical intensity images, Bulletin of the Astronomical Society of India, 39, 289.
See also: [colormap], page 870 .
map \(=\) flag ()
map \(=\) flag ( \(n\) )
Create color colormap. This colormap cycles through red, white, blue, and black with each index change.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map = gray ()
map \(=\) gray ( \(n\) )
Create gray colormap. This colormap varies from black to white with shades of gray.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
map \(=\) hot ()
\(\operatorname{map}=\operatorname{hot}(n)\)
Create color colormap. This colormap ranges from black through dark red, red, orange, yellow, to white.

The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
hsv (n)
Create color colormap. This colormap begins with red, changes through yellow, green, cyan, blue, and magenta, before returning to red.
It is useful for displaying periodic functions. The map is obtained by linearly varying the hue through all possible values while keeping constant maximum saturation and value. The equivalent code is hsv2rgb ([(0:N-1)'/N, ones \((N, 2)])\).
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
map \(=\) jet ()
\(\operatorname{map}=\) jet \((n)\)
Create color colormap. This colormap ranges from dark blue through blue, cyan, green, yellow, red, to dark red.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

Programming Note: The jet colormap is not perceptually uniform. Try the viridis colormap if that is important. For a drop-in replacement for jet with better perceptual characteristics try the turbo colormap.

See also: [colormap], page 870, [turbo], page 875.
map \(=\) lines ()
map \(=\) lines ( \(n\) )
Create color colormap. This colormap is composed of the list of colors in the current axes "ColorOrder" property. The default is blue, orange, yellow, purple, green, light blue, and dark red.

The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
map \(=\) ocean ()
\(\operatorname{map}=\operatorname{ocean}(n)\)
Create color colormap. This colormap varies from black to white with shades of blue. The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
```

map = pink ()

```
map \(=\) pink ( \(n\) )

Create color colormap. This colormap varies from black to white with shades of gray-pink.
This colormap gives a sepia tone when used on grayscale images.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map \(=\) prism ()
map \(=\) prism (n)
Create color colormap. This colormap cycles through red, orange, yellow, green, blue and violet with each index change.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map = rainbow ()
map \(=\) rainbow ( \(n\) )
Create color colormap. This colormap ranges from red through orange, yellow, green, blue, to violet.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map \(=\) spring ()
map \(=\operatorname{spring}(n)\)
Create color colormap. This colormap varies from magenta to yellow.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870 .
map \(=\) summer ()
map \(=\) summer ( \(n\) )
Create color colormap. This colormap varies from green to yellow.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
map = turbo ()
map \(=\) turbo ( \(n\) )
Create color colormap. This colormap ranges from dark blue through green to dark red; similar to the outdated jet colormap but perceptually uniform.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.
See also: [colormap], page 870 .
map \(=\) viridis ()
map \(=\) viridis (n)
Create color colormap. This colormap ranges from dark purplish-blue through blue, green, to yellow.

The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
map \(=\) white ()
map \(=\) white ( \(n\) )
Create color colormap. This colormap is completely white.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
map \(=\) winter ()
map \(=\) winter ( \(n\) )
Create color colormap. This colormap varies from blue to green.
The argument \(n\) must be a scalar. If unspecified, the length of the current colormap, or 64 , is used.

See also: [colormap], page 870.
cmap \(=\) contrast \((x)\)
cmap \(=\) contrast ( \(x, n\) )
Return a gray colormap that maximizes the contrast in an image.
The returned colormap will have \(n\) rows. If \(n\) is not defined then the size of the current colormap is used.

See also: [colormap], page 870, [brighten], page 876.
The following three functions modify the existing colormap rather than replace it.
```

map_out = brighten (beta)
map_out = brighten (map, beta)
map_out = brighten (h, beta)
brighten (...)
Brighten or darken a colormap.

```

The argument beta must be a scalar between -1 and 1 , where a negative value darkens and a positive value brightens the colormap.

If the map argument is omitted, the function is applied to the current colormap.
The first argument can also be a valid graphics handle \(h\), in which case brighten is applied to the colormap associated with this handle.
If no output is specified then the result is written to the current colormap.
See also: [colormap], page 870, [contrast], page 876.
```

spinmap ()

```
spinmap ( \(t\) )
spinmap ( \(t\), inc)
spinmap ("inf")

Cycle the colormap for \(t\) seconds with a color increment of inc.
Both parameters are optional. The default cycle time is 5 seconds and the default increment is 2 . If the option "inf" is given then cycle continuously until Control-C is pressed.
When rotating, the original color 1 becomes color 2, color 2 becomes color 3, etc. A positive or negative increment is allowed and a higher value of inc will cause faster cycling through the colormap.

See also: [colormap], page 870 .
```

whitebg ()
whitebg (color)
whitebg ("none")
whitebg (hfig)
whitebg (hfig, color)
whitebg (hfig, "none")

```

Invert the colors in the current color scheme.
The root properties are also inverted such that all subsequent plots will use the new color scheme.
If the optional argument color is present then the background color is set to color rather than inverted. color may be a string representing one of the eight known colors or an RGB triplet. The special string argument "none" restores the plot to the factory default colors.
If the first argument hfig is a figure handle or list of figure handles, then operate on these figures rather than the current figure returned by gcf. The root properties will not be changed unless 0 is in the list of figures.
Programming Note: whitebg operates by changing the color properties of the children of the specified figures. Only objects with a single color are affected. For example, a patch with a single "FaceColor" will be changed, but a patch with shading ("interp") will not be modified. For inversion, the new color is simply the inversion in RGB space: cnew \(=[1-R 1-G 1-B]\). When a color is specified, the axes and figure are set to the new color, and the color of child objects are then adjusted to have some contrast (visibility) against the new background.
See also: [reset], page 490, [get], page 430, [set], page 430.
The following functions can be used to manipulate colormaps.
```

[Y, newmap] = cmunique (X, map)
[Y, newmap] = cmunique (RGB)
[Y, newmap] = cmunique (I)
Convert an input image $X$ to an output indexed image $Y$ which uses the smallest colormap possible newmap.

```

When the input is an indexed image ( \(X\) with colormap map) the output is a colormap newmap from which any repeated rows have been eliminated. The output image, \(Y\), is the original input image with the indices adjusted to match the new, possibly smaller, colormap.

When the input is an RGB image (an MxNx 3 array), the output colormap will contain one entry for every unique color in the original image. In the worst case the new map could have as many rows as the number of pixels in the original image.

When the input is a grayscale image \(I\), the output colormap will contain one entry for every unique intensity value in the original image. In the worst case the new map could have as many rows as the number of pixels in the original image.
Implementation Details:
newmap is always an Mx3 matrix, even if the input image is an intensity grayscale image \(I\) (all three RGB planes are assigned the same value).

The output image is of class uint8 if the size of the new colormap is less than or equal to 256 . Otherwise, the output image is of class double.
See also: [rgb2ind], page 868, [gray2ind], page 868.
[ \(Y\), newmap] \(=\) cmpermute ( \(X\), map )
[ \(Y\), newmap] = cmpermute ( \(X\), map, index)
Reorder colors in a colormap.
When called with only two arguments, cmpermute randomly rearranges the colormap map and returns a new colormap newmap. It also returns the indexed image \(Y\) which is the equivalent of the original input image \(X\) when displayed using newmap.
When called with an optional third argument the order of colors in the new colormap is defined by index.
Caution: index should not have repeated elements or the function will fail.

\subsection*{32.4 Plotting on top of Images}

If gnuplot is being used to display images it is possible to plot on top of images. Since an image is a matrix it is indexed by row and column values. The plotting system is, however, based on the traditional \((x, y)\) system. To minimize the difference between the two systems Octave places the origin of the coordinate system in the point corresponding to the pixel at \((1,1)\). So, to plot points given by row and column values on top of an image, one should simply call plot with the column values as the first argument and the row values as the second. As an example the following code generates an image with random intensities between 0 and 1, and shows the image with red circles over pixels with an intensity above 0.99 .
```

I = rand (100, 100);
[row, col] = find (I > 0.99);
hold ("on");
imshow (I);
plot (col, row, "ro");
hold ("off");

```

\subsection*{32.5 Color Conversion}

Octave supports conversion from the RGB color system to the HSV color system and vice versa. It is also possible to convert from a color RGB image to a grayscale image.
hsv_map \(=\) rgb2hsv (rgb_map)
hsv_img = rgb2hsv (rgb_img)
Transform a colormap or image from RGB to HSV color space.
A color in the RGB space consists of red, green, and blue intensities.
A color in HSV space is represented by hue, saturation and value (brightness) levels in a cylindrical coordinate system. Hue is the azimuth and describes the dominant color. Saturation is the radial distance and gives the amount of hue mixed into the color. Value is the height and is the amount of light in the color.
Output class and size will be the same as input.
See also: [hsv2rgb], page 879, [rgb2ind], page 868, [rgb2gray], page 879.
rgb_map \(=\) hsv2rgb (hsv_map)
rgb_img = hsv2rgb (hsv_img)
Transform a colormap or image from HSV to RGB color space.
A color in HSV space is represented by hue, saturation and value (brightness) levels in a cylindrical coordinate system. Hue is the azimuth and describes the dominant color. Saturation is the radial distance and gives the amount of hue mixed into the color. Value is the height and is the amount of light in the color.
The input can be both a colormap or RGB image. In the case of floating point input, values are expected to be on the \(\left[\begin{array}{ll}0 & 1\end{array}\right]\) range. In the case of hue (azimuth), since the value corresponds to an angle, mod (h, 1) is used.
```

>> hsv2rgb ([0.5 1 1 1])
=>ans = 0 1 1
>> hsv2rgb ([2.5.5 1 1])
mans = 0 1 1
>> hsv2rgb ([[3.5 1 1])

# ans = 0 1 1

```

Output class and size will be the same as input.
See also: [rgb2hsv], page 879, [ind2rgb], page 868.
\(I=r g b 2 g r a y\left(r g b \_i m g\right)\)
gray_map \(=\) rgb2gray (rgb_map)
Transform an image or colormap from red-green-blue (RGB) color space to a grayscale intensity image.
The input may be of class uint8, int8, uint16, int16, single, or double. The output is of the same class as the input.
Implementation Note: The grayscale intensity is calculated as
\(I=0.298936 * R+0.587043 * G+0.114021 * B\)
which corresponds to the luminance channel when RGB is translated to YIQ as documented in https://en.wikipedia.org/wiki/YIQ.
See also: [rgb2hsv], page 879, [rgb2ind], page 868.

\section*{33 Audio Processing}

\subsection*{33.1 Audio File Utilities}

The following functions allow you to read, write and retrieve information about audio files. Various formats are supported including wav, flac and ogg vorbis.
info = audioinfo (filename)
Return information about an audio file specified by filename.
The output info is a structure containing the following fields:
'Filename'
Name of the audio file.
'CompressionMethod'
Audio compression method. Unused, only present for compatibility with matlab.
'NumChannels'
Number of audio channels.
'SampleRate'
Sample rate of the audio, in Hertz.
'TotalSamples'
Number of samples in the file.
'Duration'
Duration of the audio, in seconds.
'BitsPerSample'
Number of bits per sample.
'BitRate' Audio bit rate. Unused, only present for compatibility with matlab.
'Title' "Title" audio metadata value as a string, or empty if not present.
'Artist' "Artist" audio metadata value as a string, or empty if not present.
'Comment' "Comment" audio metadata value as a string, or empty if not present.
See also: [audioread], page 881, [audiowrite], page 882.
[y, fs] = audioread (filename)
[y, fs] = audioread (filename, samples)
[y, fs] = audioread (filename, datatype)
[y, fs] = audioread (filename, samples, datatype)
Read the audio file filename and return the audio data \(y\) and sampling rate fs.
The audio data is stored as matrix with rows corresponding to audio frames and columns corresponding to channels.
The optional two-element vector argument samples specifies starting and ending frames.
The optional argument datatype specifies the datatype to return. If it is "native", then the type of data depends on how the data is stored in the audio file.
See also: [audiowrite], page 882, [audioformats], page 882, [audioinfo], page 881.
audiowrite (filename, y, fs)
audiowrite (filename, y, fs, name, value, ...)
Write audio data from the matrix \(y\) to filename at sampling rate \(f_{S}\) with the file format determined by the file extension.
Additional name/value argument pairs may be used to specify the following options:
'BitsPerSample'
Number of bits per sample. Valid values are \(8,16,24\), and 32 . Default is 16.
'BitRate' Valid argument name, but ignored. Left for compatibility with matlab.
'Quality' Quality setting for the Ogg Vorbis compressor. Values can range between 0 and 100 with 100 being the highest quality setting. Default is 75 .
'Title' Title for the audio file.
'Artist' Artist name.
'Comment' Comment.
See also: [audioread], page 881, [audioformats], page 882, [audioinfo], page 881.
audioformats ()
audioformats (format)
Display information about all supported audio formats.
If the optional argument format is given, then display only formats with names that start with format.

See also: [audioread], page 881, [audiowrite], page 882.

\subsection*{33.2 Audio Device Information}
```

devinfo = audiodevinfo ()
devs = audiodevinfo (io)
name = audiodevinfo (io, id)
id = audiodevinfo (io, name)
driverversion = audiodevinfo (io, id, "DriverVersion")
id = audiodevinfo (io, rate, bits, chans)
supports = audiodevinfo (io, id, rate, bits, chans)

```

Return a structure describing the available audio input and output devices.
The devinfo structure has two fields "input" and "output". The value of each field is a structure array with fields "Name", "DriverVersion" and "ID" describing an audio device.
If the optional argument io is 1 , return information about input devices only. If it is 0 , return information about output devices only. If io is the only argument supplied, return the number of input or output devices available.
If the optional argument id is provided, return information about the corresponding device.
If the optional argument name is provided, return the ID of the named device.

If the optional argument "DriverVersion" is given, return the name of the driver for the specified device.
Given a sampling rate, bits per sample, and number of channels for an input or output device, return the ID of the first device that supports playback or recording using the specified parameters.
If also given a device ID, return true if the device supports playback or recording using those parameters.

\subsection*{33.3 Audio Player}

The following methods are used to create and use audioplayer objects. These objects can be used to play back audio data stored in Octave matrices and arrays. The audioplayer object supports playback from various devices available to the system, blocking and non-blocking playback, convenient pausing and resuming and much more.
```

player = audioplayer (y,fs)
player = audioplayer (y, fs, nbits)
player = audioplayer (y, fs, nbits, id)
player = audioplayer (recorder)
player = audioplayer (recorder, id)

```

Create an audioplayer object that will play back data \(y\) at sample rate \(f s\).
The optional arguments nbits, and id specify the bit depth and player device id, respectively. Device IDs may be found using the audiodevinfo function. Given an audioplayer object, use the data from the object to initialize the player.
The signal \(y\) can be a vector or a two-dimensional array.
The following example will create an audioplayer object that will play back one second of white noise at 44100 sample rate using 8 bits per sample.
```

y = 0.25 * randn (2, 44100);
player = audioplayer (y, 44100, 8);
play (player);

```

\subsection*{33.3.1 Playback}

The following methods are used to control player playback.
```

play (player)
play (player, start)
play (player, limits)

```

Play audio stored in the audioplayer object player without blocking.
Given optional argument start, begin playing at start samples in the recording. Given a two-element vector limits, begin and end playing at the number of samples specified by the elements of the vector.
```

playblocking (player)
playblocking (player, start)
playblocking (player, limits)

```

Play audio stored in the audioplayer object player with blocking.

Given optional argument start, begin playing at start samples in the recording. Given a two-element vector limits, begin and end playing at the number of samples specified by the elements of the vector.
```

pause (player)

```

Pause the audioplayer player.
resume (player)
Resume playback for the paused audioplayer object player.
stop (player)
Stop the playback for the audioplayer player and reset the relevant variables to their starting values.
isplaying (player)
Return true if the audioplayer object player is currently playing back audio and false otherwise.

\subsection*{33.3.2 Properties}

The remaining couple of methods are used to get and set various properties of the audioplayer object.
```

value = get (player, name)
values = get (player)

```

Return the value of the property identified by name.
If name is a cell array return the values of the properties identified by the elements of the cell array. Given only the player object, return a scalar structure with values of all properties of player. The field names of the structure correspond to property names.
```

set (player, name, value)

```
set (player, properties)
properties = set (player)

Set the value of property specified by name to a given value.
If name and value are cell arrays, set each property to the corresponding value. Given a structure of properties with fields corresponding to property names, set the value of those properties to the field values. Given only the audioplayer object, return a structure of settable properties.

\subsection*{33.4 Audio Recorder}

The following methods are used to create and use audiorecorder objects. These objects can be used to record audio data from various devices available to the system. You can use convenient methods to retrieve that data or audioplayer objects created from that data. Methods for blocking and non-blocking recording, pausing and resuming recording and much more is available.
```

recorder = audiorecorder ()
recorder = audiorecorder (fs, nbits, channels)

```
recorder \(=\) audiorecorder (fs, nbits, channels, id)
Create an audiorecorder object recording 8 bit mono audio at 8000 Hz sample rate.
The optional arguments fs, nbits, channels, and id specify the sample rate, bit depth, number of channels and recording device id, respectively. Device IDs may be found using the audiodevinfo function.

\subsection*{33.4.1 Recording}

The following methods control the recording process.
record (recorder)
record (recorder, length)
Record audio without blocking using the audiorecorder object recorder until stopped or paused by the stop or pause method.
Given the optional argument length, record for length seconds.
```

recordblocking (recorder, length)

```

Record audio with blocking (synchronous I/O).
The length of the recording in seconds (length) must be specified.

\section*{pause (recorder)}

Pause recording with audiorecorder object recorder.
resume (recorder)
Resume recording with the paused audiorecorder object recorder.
stop (recorder)
Stop the audiorecorder object recorder and clean up any audio streams.
```

isrecording (recorder)

```

Return true if the audiorecorder object recorder is currently recording audio and false otherwise.

\subsection*{33.4.2 Data Retrieval}

The following methods allow you to retrieve recorded audio data in various ways.
```

data = getaudiodata (recorder)
data = getaudiodata (recorder, datatype)

```

Return recorder audio data as a matrix with values between -1.0 and 1.0 and with as many columns as there are channels in the recorder.
Given the optional argument datatype, convert the recorded data to the specified type, which may be one of "double", "single", "int16", "int8" or "uint8".
player = getplayer (recorder)
Return an audioplayer object with data recorded by the audiorecorder object recorder.
```

player = play (recorder)
player = play (recorder, start)
player = play (recorder, [start, end])

```

Play the audio recorded in recorder and return a corresponding audioplayer object.

If the optional argument start is provided, begin playing start seconds in to the recording.
If the optional argument end is provided, stop playing at end seconds in the recording.

\subsection*{33.4.3 Properties}

The remaining two methods allow you to read or alter the properties of audiorecorder objects.
```

value = get (recorder, name)
values = get (recorder)

```

Return the value of the property identified by name.
If name is a cell array, return the values of the properties corresponding to the elements of the cell array. Given only the recorder object, return a scalar structure with values of all properties of recorder. The field names of the structure correspond to property names.
```

set (recorder, name, value)
set (recorder, properties)
properties = set (recorder)

```

Set the value of property specified by name to a given value.
If name and value are cell arrays of the same size, set each property to a corresponding value. Given a structure with fields corresponding to property names, set the value of those properties to the corresponding field values. Given only the recorder object, return a structure of settable properties.

\subsection*{33.5 Audio Data Processing}

Octave provides a few functions for dealing with audio data. An audio 'sample' is a single output value from an A/D converter, i.e., a small integer number (usually 8 or 16 bits), and audio data is just a series of such samples. It can be characterized by three parameters: the sampling rate (measured in samples per second or Hz, e.g., 8000 or 44100), the number of bits per sample (e.g., 8 or 16 ), and the number of channels ( 1 for mono, 2 for stereo, etc.).

There are many different formats for representing such data. Currently, only the two most popular, linear encoding and mu-law encoding, are supported by Octave. There is an excellent FAQ on audio formats by Guido van Rossum guido@cwi.nl which can be found at any FAQ ftp site, in particular in the directory/pub/usenet/news.answers/audio-fmts of the archive site rtfm.mit.edu.

Octave simply treats audio data as vectors of samples (non-mono data are not supported yet). It is assumed that audio files using linear encoding have one of the extensions lin or raw, and that files holding data in mu-law encoding end in \(\mathrm{au}, \mathrm{mu}\), or snd.
\(y=\operatorname{lin} 2 m u(x, n)\)
Convert audio data from linear to mu-law.
Mu-law values use 8 -bit unsigned integers. Linear values use \(n\)-bit signed integers or floating point values in the range \(-1 \leq x \leq 1\) if \(n\) is 0 .
If \(n\) is not specified it defaults to 0,8 , or 16 depending on the range of values in \(x\).
See also: [mu2lin], page 887.
\(y=m u 2 l i n(x, n)\)
Convert audio data from mu-law to linear.
Mu-law values are 8 -bit unsigned integers. Linear values use \(n\)-bit signed integers or floating point values in the range \(-1 \leq y \leq 1\) if \(n\) is 0 .
If \(n\) is not specified it defaults to 0 .
See also: [lin2mu], page 886.
record (sec)
record (sec, fs)
Record sec seconds of audio from the system's default audio input at a sampling rate of 8000 samples per second.
If the optional argument \(f s\) is given, it specifies the sampling rate for recording.
For more control over audio recording, use the audiorecorder class.
See also: [sound], page 887, [soundsc], page 887.
sound (y)
sound ( \(y, f s\) )
sound ( \(\mathrm{y}, \mathrm{fs}\), nbits)
Play audio data \(y\) at sample rate \(f s\) to the default audio device.
The audio signal y can be a vector or a two-column array, representing mono or stereo audio, respectively.
If \(f_{s}\) is not given, a default sample rate of 8000 samples per second is used.
The optional argument nbits specifies the bit depth to play to the audio device and defaults to 8 bits.
For more control over audio playback, use the audioplayer class.
See also: [soundsc], page 887, [record], page 887.
soundsc (y)
soundsc ( \(y\), fs)
soundsc (y, fs, nbits)
soundsc (..., [ymin, ymax])
Scale the audio data \(y\) and play it at sample rate \(f\) s to the default audio device.
The audio signal y can be a vector or a two-column array, representing mono or stereo audio, respectively.
If \(f s\) is not given, a default sample rate of 8000 samples per second is used.
The optional argument nbits specifies the bit depth to play to the audio device and defaults to 8 bits.
By default, \(y\) is automatically normalized to the range \([-1,1]\). If the range [ymin, \(y \max ]\) is given, then elements of \(y\) that fall within the range \(y\) min \(\leq y \leq y \max\) are scaled to the range \([-1,1]\) instead.
For more control over audio playback, use the audioplayer class.
See also: [sound], page 887, [record], page 887.

\section*{34 Object Oriented Programming}

Octave has the ability to create user-defined classes-including the capabilities of operator and function overloading. Classes can protect internal properties so that they may not be altered accidentally which facilitates data encapsulation. In addition, rules can be created to address the issue of class precedence in mixed class operations.

This chapter discusses the means of constructing a user class, how to query and set the properties of a class, and how to overload operators and functions. Throughout this chapter real code examples are given using a class designed for polynomials.

\subsection*{34.1 Creating a Class}

This chapter illustrates user-defined classes and object oriented programming through a custom class designed for polynomials. This class was chosen for its simplicity which does not distract unnecessarily from the discussion of the programming features of Octave. Even so, a bit of background on the goals of the polynomial class is necessary before the syntax and techniques of Octave object oriented programming are introduced.

The polynomial class is used to represent polynomials of the form
\[
a_{0}+a_{1} x+a_{2} x^{2}+\ldots a_{n} x^{n}
\]
where \(a_{0}, a_{1}\), etc. are elements of \(\Re\). Thus the polynomial can be represented by a vector
\[
a=[a 0, a 1, a 2, \ldots, a n] ;
\]

This is a sufficient specification to begin writing the constructor for the polynomial class. All object oriented classes in Octave must be located in a directory that is the name of the class prepended with the ' \(@\) ' symbol. For example, the polynomial class will have all of its methods defined in the @polynomial directory.

The constructor for the class must be the name of the class itself; in this example the constructor resides in the file @polynomial/polynomial.m. Ideally, even when the constructor is called with no arguments it should return a valid object. A constructor for the polynomial class might look like
```


## -*- texinfo -*-

## @deftypefn {} {} polynomial ()

## @deftypefnx {} {} polynomial (@var{a})

## Create a polynomial object representing the polynomial

## 

## @example

## a0 + a1 * x + a2 * x^2 + @dots{} + an * x^n

## @end example

## 

## @noindent

## from a vector of coefficients [a0 a1 a2 @dots{} an].

## @end deftypefn

function p = polynomial (a)

```
```

    if (nargin == 0)
    p.poly = 0;
    p = class (p, "polynomial");
    else
        if (isa (a, "polynomial"))
        p = a;
    elseif (isreal (a) && isvector (a))
        p.poly = a(:).'; # force row vector
        p = class (p, "polynomial");
    else
        error ("@polynomial: A must be a real vector");
    endif
    endif
    endfunction

```

Note that the return value of the constructor must be the output of the class function. The first argument to the class function is a structure and the second is the name of the class itself. An example of calling the class constructor to create an instance is
```

p = polynomial ([1, 0, 1]);

```

Methods are defined by m-files in the class directory and can have embedded documentation the same as any other m-file. The help for the constructor can be obtained by using the constructor name alone, that is, for the polynomial constructor help polynomial will return the help string. Help can be restricted to a particular class by using the class directory name followed by the method. For example, help @polynomial/polynomial is another way of displaying the help string for the polynomial constructor. This second means is the only way to obtain help for the overloaded methods and functions of a class.

The same specification mechanism can be used wherever Octave expects a function name. For example type @polynomial/disp will print the code of the disp method of the polynomial class to the screen, and dbstop @polynomial/disp will set a breakpoint at the first executable line of the disp method of the polynomial class.

To check whether a variable belongs to a user class, the isobject and isa functions can be used. For example:
```

$\mathrm{p}=$ polynomial ([1, 0, 1]);
isobject (p)
$\Rightarrow 1$
isa (p, "polynomial")
$\Rightarrow 1$

```
isobject (x)

Return true if x is a class object.
See also: [class], page 41, [typeinfo], page 41, [isa], page 41, [ismethod], page 891, [isprop], page 423.

The available methods of a class can be displayed with the methods function.
```

methods (obj)
methods ("classname")

```
```

methods (..., "-full")
mtds = methods (...)

```

List the names of the public methods for the object obj or the named class classname.
obj may be an Octave class object or a Java object. classname may be the name of an Octave class or a Java class.
If the optional argument "-full" is given then Octave returns full method signatures which include output type, name of method, and the number and type of inputs.
When called with no output arguments, methods prints the list of method names to the screen. Otherwise, the output argument mtds contains the list in a cell array of strings.

See also: [ismethod], page 891, [properties], page 909, [fieldnames], page 119.
To inquire whether a particular method exists for a user class, the ismethod function can be used.
```

ismethod (obj, method)
ismethod (class_name, method)

```

Return true if the string method is a valid method of the object obj or of the class clsname.

See also: [isprop], page 423, [isobject], page 890, [isjava], page 1055, [methods], page 890 .

For a polynomial class it makes sense to have a method to compute its roots.
```

function r = roots (p)
r = roots (fliplr (p.poly));
endfunction

```

We can check for the existence of the roots-method by calling:
```

p = polynomial ([1, 0, 1]);
ismethod (p, "roots")
\# 1

```

\subsection*{34.2 Class Methods}

There are a number of basic class methods that can (and should) be defined to allow the contents of the classes to be queried and set. The most basic of these is the disp method. The disp method is used by Octave whenever a class should be displayed on the screen. Usually this is the result of an Octave expression that doesn't end with a semicolon. If this method is not defined, then Octave won't print anything when displaying the contents of a class which can be confusing.
An example of a disp method for the polynomial class might be
```

function disp (p)
a = p.poly;
first = true;
for i = 1 : length (a);
if (a(i) != 0)

```
```

            if (first)
                        first = false;
            elseif (a(i) > 0 || isnan (a(i)))
                        printf (" +");
            endif
            if (a(i) < 0)
                        printf (" -");
            endif
            if (i == 1)
                        printf (" %.5g", abs (a(i)));
            elseif (abs (a(i)) != 1)
                    printf (" %.5g *", abs (a(i)));
            endif
            if (i > 1)
            printf (" X");
            endif
            if (i > 2)
                    printf (" ~ %d", i - 1);
            endif
            endif
    endfor
    if (first)
        printf (" 0");
    endif
    printf ("\n");
    endfunction

```

To be consistent with the Octave graphic handle classes, a class should also define the get and set methods. The get method accepts one or two arguments. The first argument is an object of the appropriate class. If no second argument is given then the method should return a structure with all the properties of the class. If the optional second argument is given it should be a property name and the specified property should be retrieved.
```

function val = get (p, prop)
if (nargin < 1)
print_usage ();
endif
if (nargin == 1)
val.poly = p.poly;
else
if (! ischar (prop))
error ("@polynomial/get: PROPERTY must be a string");
endif

```
```

        switch (prop)
        case "poly"
            val = p.poly;
        otherwise
            error ('@polynomial/get: invalid PROPERTY "%s"', prop);
        endswitch
    endif
    endfunction

```

Similarly, the first argument to the set method should be an object and any additional arguments should be property/value pairs.
```

function pout = set (p, varargin)
if (numel (varargin) < 2 || rem (numel (varargin), 2) != 0)
error ("@polynomial/set: expecting PROPERTY/VALUE pairs");
endif
pout = p;
while (numel (varargin) > 1)
prop = varargin{1};
val = varargin{2};
varargin(1:2) = [];
if (! ischar (prop) || ! strcmp (prop, "poly"))
error ("@polynomial/set: invalid PROPERTY for polynomial class");
elseif (! (isreal (val) \&\& isvector (val)))
error ("@polynomial/set: VALUE must be a real vector");
endif
pout.poly = val(:).'; \# force row vector
endwhile
endfunction

```

Note that Octave does not implement pass by reference; Therefore, to modify an object requires an assignment statement using the return value from the set method.
```

p = set (p, "poly", [1, 0, 0, 0, 1]);

```

The set method makes use of the subsasgn method of the class, and therefore this method must also be defined. The subsasgn method is discussed more thoroughly in the next section (see Section 34.3 [Indexing Objects], page 894).

Finally, user classes can be considered to be a special type of a structure, and they can be saved to a file in the same manner as a structure. For example:
```

p = polynomial ([1, 0, 1]);
save userclass.mat p
clear p
load userclass.mat

```

All of the file formats supported by save and load are supported. In certain circumstances a user class might contain a field that it doesn't make sense to save, or a field that needs to be initialized before it is saved. This can be done with the saveobj method of the class.
\(b=\) saveobj (a)
Method of a class to manipulate an object prior to saving it to a file.
The function saveobj is called when the object a is saved using the save function. An example of the use of saveobj might be to remove fields of the object that don't make sense to be saved or it might be used to ensure that certain fields of the object are initialized before the object is saved. For example:
```

function b = saveobj (a)
b = a;
if (isempty (b.field))
b.field = initfield (b);
endif
endfunction

```

See also: [loadobj], page 894, [class], page 41.
saveobj is called just prior to saving the class to a file. Similarly, the loadobj method is called just after a class is loaded from a file, and can be used to ensure that any removed fields are reinserted into the user object.
b = loadobj (a)
Method of a class to manipulate an object after loading it from a file.
The function loadobj is called when the object a is loaded using the load function. An example of the use of saveobj might be to add fields to an object that don't make sense to be saved. For example:
```

function b = loadobj (a)
b = a;
b.addmissingfield = addfield (b);
endfunction

```

See also: [saveobj], page 894, [class], page 41.

\subsection*{34.3 Indexing Objects}

\subsection*{34.3.1 Defining Indexing And Indexed Assignment}

Objects can be indexed with parentheses or braces, either like obj(idx) or like obj\{idx\}, or even like obj(idx).field. However, it is up to the programmer to decide what this indexing actually means. In the case of the polynomial class \(p(n)\) might mean either the coefficient of the \(n\)-th power of the polynomial, or it might be the evaluation of the polynomial at \(n\). The meaning of this subscripted referencing is determined by the subsref method.
subsref (val, idx)
Perform the subscripted element selection operation on val according to the subscript specified by idx.

The subscript idx must be a structure array with fields 'type' and 'subs'. Valid values for 'type’ are "()", "\{\}", and ".". The 'subs’ field may be either ":" or a cell array of index values.
The following example shows how to extract the first two columns of a matrix
```

val = magic (3)
\# val = [ % 1 % 6
3 5 7
4 9 2 ]
idx.type = "()";
idx.subs = {":", 1:2};
subsref (val, idx)
[ [ 8 1
3 5
4 1]

```

Note that this is the same as writing val(:, 1:2).
If idx is an empty structure array with fields 'type' and 'subs', return val.
See also: [subsasgn], page 896, [substruct], page 122.
For example, this class uses the convention that indexing with "()" evaluates the polynomial and indexing with "\{\}" returns the \(n\)-th coefficient (of the \(n\)-th power). The code for the subsref method looks like
```

function r = subsref (p, s)
if (isempty (s))
error ("@polynomial/subsref: missing index");
endif
switch (s(1).type)
case "()"
idx = s(1).subs;
if (numel (idx) != 1)
error ("@polynomial/subsref: need exactly one index");
endif
r = polyval (fliplr (p.poly), idx{1});
case "{}"
idx = s(1).subs;
if (numel (idx) != 1)
error ("@polynomial/subsref: need exactly one index");
endif
if (isnumeric (idx{1}))
r = p.poly(idx{1}+1);
else
r = p.poly(idx{1});

```
```

            endif
            case "."
                        fld = s.subs;
                        if (! strcmp (fld, "poly"))
                    error ('@polynomial/subsref: invalid property "%s"', fld);
                    endif
                        r = p.poly;
                otherwise
                        error ("@polynomial/subsref: invalid subscript type");
        endswitch
        if (numel (s) > 1)
        r = subsref (r, s(2:end));
        endif
    endfunction

```

The equivalent functionality for subscripted assignments uses the subsasgn method.
subsasgn (val, idx, rhs)
Perform the subscripted assignment operation according to the subscript specified by idx.
The subscript idx must be a structure array with fields 'type' and 'subs'. Valid values for 'type' are "()", "\{\}", and ".". The 'subs' field may be either ":" or a cell array of index values.
The following example shows how to set the two first columns of a 3 -by- 3 matrix to zero.
```

val = magic (3);
idx.type = "()";
idx.subs = {":", 1:2};
val = subsasgn (val, idx, 0)
[ [ 0 0 0 6
0

```

Note that this is the same as writing val (: , 1:2) \(=0\).
If idx is an empty structure array with fields 'type' and 'subs', return rhs.
See also: [subsref], page 894, [substruct], page 122, [optimize_subsasgn_calls], page 896 .
```

val = optimize_subsasgn_calls ()
old_val = optimize_subsasgn_calls (new_val)
optimize_subsasgn_calls (new_val, "local")

```

Query or set the internal flag for subsasgn method call optimizations.

If true, Octave will attempt to eliminate the redundant copying when calling the subsasgn method of a user-defined class.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [subsasgn], page 896.
Note that the subsref and subsasgn methods always receive the whole index chain, while they usually handle only the first element. It is the responsibility of these methods to handle the rest of the chain (if needed), usually by forwarding it again to subsref or subsasgn.

If you wish to use the end keyword in subscripted expressions of an object, then there must be an end method defined. For example, the end method for the polynomial class might look like
```

function r = end (obj, index_pos, num_indices)
if (num_indices != 1)
error ("polynomial object may only have one index");
endif
r = length (obj.poly) - 1;
endfunction

```
which is a fairly generic end method that has a behavior similar to the end keyword for Octave Array classes. An example using the polynomial class is then
```

p = polynomial ([1,2,3,4]);
p{end-1}
=> 3

```

Objects can also be used themselves as the index in a subscripted expression and this is controlled by the subsindex function.
\(i d x=\) subsindex (obj)
Convert an object to an index vector.
When obj is a class object defined with a class constructor, then subsindex is the overloading method that allows the conversion of this class object to a valid indexing vector. It is important to note that subsindex must return a zero-based real integer vector of the class "double". For example, if the class constructor were
```

function obj = myclass (a)
obj = class (struct ("a", a), "myclass");
endfunction

```
then the subsindex function
```

function idx = subsindex (obj)
idx = double (obj.a) - 1.0;
endfunction

```
could be used as follows
```

a = myclass (1:4);
b = 1:10;
b(a)

# 1 2 2 3 4

```

See also: [class], page 41, [subsref], page 894, [subsasgn], page 896.
Finally, objects can be used like ranges by providing a colon method.
```

r = colon (base, limit)
r = colon (base, increment, limit)

```

Return the result of the colon expression corresponding to base, limit, and optionally, increment.
This function is equivalent to the operator syntax base : limit or base : increment : limit.
See also: [linspace], page 530.

\subsection*{34.3.2 Indexed Assignment Optimization}

Octave's ubiquitous lazily-copied pass-by-value semantics implies a problem for performance of user-defined subsasgn methods. Imagine the following call to subsasgn
```

ss = substruct ("()", {1});
x = subsasgn (x, ss, 1);

```
where the corresponding method looking like this:
```

function x = subsasgn (x, ss, val)
x.myfield (ss.subs{1}) = val;
endfunction

```

The problem is that on entry to the subsasgn method, x is still referenced from the caller's scope, which means that the method will first need to unshare (copy) x and x .myfield before performing the assignment. Upon completing the call, unless an error occurs, the result is immediately assigned to x in the caller's scope, so that the previous value of x. myfield is forgotten. Hence, the Octave language implies a copy of N elements ( N being the size of x. myfield), where modifying just a single element would actually suffice. In other words, a constant-time operation is degraded to linear-time one. This may be a real problem for user classes that intrinsically store large arrays.

To partially solve the problem Octave uses a special optimization for user-defined subsasgn methods coded as m-files. When the method gets called as a result of the built-in assignment syntax (not a direct subsasgn call as shown above), i.e., \(x(1)=1\), AND if the subsasgn method is declared with identical input and output arguments, as in the example above, then Octave will ignore the copy of x inside the caller's scope; therefore, any changes made to x during the method execution will directly affect the caller's copy as well. This allows, for instance, defining a polynomial class where modifying a single element takes constant time.

It is important to understand the implications that this optimization brings. Since no extra copy of x will exist in the caller's scope, it is solely the callee's responsibility to not
leave x in an invalid state if an error occurs during the execution. Also, if the method partially changes x and then errors out, the changes will affect x in the caller's scope. Deleting or completely replacing x inside subsasgn will not do anything, however, only indexed assignments matter.

Since this optimization may change the way code works (especially if badly written), a function optimize_subsasgn_calls is provided to control it. This feature is enabled by default. Another way to avoid the optimization is to declare subsasgn methods with different output and input arguments like this:
```

function y = subsasgn (x, ss, val)
endfunction

```

\subsection*{34.4 Overloading Objects}

\subsection*{34.4.1 Function Overloading}

Any Octave function can be overloaded, and this allows an object-specific version of a function to be called as needed. A pertinent example for the polynomial class might be to overload the polyval function.
```

function [y, dy] = polyval (p, varargin)
if (nargout > 1)
[y, dy] = polyval (fliplr (p.poly), varargin{:});
else
y = polyval (fliplr (p.poly), varargin{:});
endif
endfunction

```

This function just hands off the work to the normal Octave polyval function. Another interesting example of an overloaded function for the polynomial class is the plot function.
```

function h = plot (p, varargin)
n = 128;
rmax = max (abs (roots (p.poly)));
x = [0 : (n - 1)] / (n - 1) * 2.2 * rmax - 1.1 * rmax;
if (nargout > 0)
h = plot (x, polyval (p, x), varargin{:});
else
plot (x, polyval (p, x), varargin{:});
endif
endfunction

```
which allows polynomials to be plotted in the domain near the region of the roots of the polynomial.

Functions that are of particular interest for overloading are the class conversion functions such as double. Overloading these functions allows the cast function to work with a user
class. It can also aid in the use of a class object with methods and functions from other classes since the object can be transformed to the requisite input form for the new function. An example double function for the polynomial class might look like
```

function a = double (p)
a = p.poly;
endfunction

```

\subsection*{34.4.2 Operator Overloading}

The following table shows, for each built-in numerical operation, the corresponding function name to use when providing an overloaded method for a user class.
\begin{tabular}{|c|c|c|}
\hline Operation & Method & Description \\
\hline \(a+b\) & plus (a, b) & Binary addition \\
\hline \(a-b\) & minus ( \(\mathrm{a}, \mathrm{b}\) ) & Binary subtraction \\
\hline +a & uplus (a) & Unary addition \\
\hline -a & uminus (a) & Unary subtraction \\
\hline \(\mathrm{a} . * \mathrm{~b}\) & times ( \(\mathrm{a}, \mathrm{b}\) ) & Element-wise multiplication \\
\hline \(\mathrm{a} * \mathrm{~b}\) & mtimes ( \(\mathrm{a}, \mathrm{b}\) ) & Matrix multiplication \\
\hline \(\mathrm{a} . / \mathrm{b}\) & rdivide (a, b) & Element-wise right division \\
\hline \(\mathrm{a} / \mathrm{b}\) & mrdivide (a, b) & Matrix right division \\
\hline \(\mathrm{a} . \backslash \mathrm{b}\) & ldivide ( \(\mathrm{a}, \mathrm{b}\) ) & Element-wise left division \\
\hline \(a \backslash b\) & mldivide (a, b) & Matrix left division \\
\hline \(\mathrm{a} \cdot{ }^{\text {- }} \mathrm{b}\) & power ( \(\mathrm{a}, \mathrm{b}\) ) & Element-wise power \\
\hline \(\mathrm{a}^{\wedge} \mathrm{b}\) & mpower (a, b) & Matrix power \\
\hline \(\mathrm{a}<\mathrm{b}\) & lt ( \(\mathrm{a}, \mathrm{b}\) ) & Less than \\
\hline \(\mathrm{a}<=\mathrm{b}\) & le ( \(\mathrm{a}, \mathrm{b}\) ) & Less than or equal to \\
\hline \(\mathrm{a}>\mathrm{b}\) & gt ( \(\mathrm{a}, \mathrm{b}\) ) & Greater than \\
\hline \(\mathrm{a}>=\mathrm{b}\) & ge ( \(\mathrm{a}, \mathrm{b}\) ) & Greater than or equal to \\
\hline \(\mathrm{a}==\mathrm{b}\) & eq ( \(\mathrm{a}, \mathrm{b}\) ) & Equal to \\
\hline a ! = b & ne ( \(\mathrm{a}, \mathrm{b}\) ) & Not equal to \\
\hline \(\mathrm{a} \& \mathrm{~b}\) & and ( \(\mathrm{a}, \mathrm{b}\) ) & Logical and \\
\hline \(\mathrm{a} \mid \mathrm{b}\) & or ( \(\mathrm{a}, \mathrm{b}\) ) & Logical or \\
\hline ! a & not (a) & Logical not \\
\hline \(a^{\prime}\) & ctranspose (a) & Complex conjugate transpose \\
\hline a.' & transpose (a) & Transpose \\
\hline \(a: b\) & colon ( \(\mathrm{a}, \mathrm{b}\) ) & Two element range \\
\hline a:b:c & colon ( \(\mathrm{a}, \mathrm{b}, \mathrm{c}\) ) & Three element range \\
\hline [a, b] & horzcat ( \(\mathrm{a}, \mathrm{b}\) ) & Horizontal concatenation \\
\hline [a; b] & vertcat (a, b) & Vertical concatenation \\
\hline \(\mathrm{a}\left(\mathrm{s}_{1}, \ldots, \mathrm{~s}_{n}\right)\) & subsref (a, s) & Subscripted reference \\
\hline \(\mathrm{a}\left(\mathrm{s}_{1}, \ldots, \mathrm{~s}_{n}\right)=\mathrm{b}\) & subsasgn ( \(\mathrm{a}, \mathrm{s}, \mathrm{b}\) ) & Subscripted assignment \\
\hline b (a) & subsindex (a) & Convert object to index \\
\hline disp & disp (a) & Object display \\
\hline
\end{tabular}

Table 34.1: Available overloaded operators and their corresponding class method
An example mtimes method for the polynomial class might look like
```

function p = mtimes (a, b)
p = polynomial (conv (double (a), double (b)));
endfunction

```

\subsection*{34.4.3 Precedence of Objects}

Many functions and operators take two or more arguments and the situation can easily arise where these functions are called with objects of different classes. It is therefore necessary to determine the precedence of which method from which class to call when there are mixed objects given to a function or operator. To do this the superiorto and inferiorto functions can be used
```

superiorto (class_name, ...)

```

When called from a class constructor, mark the object currently constructed as having a higher precedence than class_name.
More that one such class can be specified in a single call. This function may only be called from a class constructor.
See also: [inferiorto], page 901.
inferiorto (class_name, ...)
When called from a class constructor, mark the object currently constructed as having a lower precedence than class_name.
More that one such class can be specified in a single call. This function may only be called from a class constructor.
See also: [superiorto], page 901.
With the polynomial class, consider the case
```

2 * polynomial ([1, 0, 1]);

```
that mixes an object of the class "double" with an object of the class "polynomial". In this case the return type should be "polynomial" and so the superiorto function is used in the class constructor. In particular the polynomial class constructor would be modified to
```


## -*- texinfo -*-

## @deftypefn {} {} polynomial ()

## @deftypefnx {} {} polynomial (@var{a})

## Create a polynomial object representing the polynomial

## 

## @example

## a0 + a1 * x + a2 * x^2 + @dots{} + an * x^n

## @end example

## 

## @noindent

## from a vector of coefficients [a0 a1 a2 @dots{} an].

## @end deftypefn

function p = polynomial (a)

```
```

    if (nargin == 0)
    p.poly = [0];
    p = class (p, "polynomial");
    else
        if (strcmp (class (a), "polynomial"))
        p = a;
    elseif (isreal (a) && isvector (a))
        p.poly = a(:).'; # force row vector
        p = class (p, "polynomial");
    else
        error ("@polynomial: A must be a real vector");
    endif
    endif
    superiorto ("double");
    endfunction

```

Note that user classes always have higher precedence than built-in Octave types. Thus, marking the polynomial class higher than the "double" class is not actually necessary.

When confronted with two objects of equal precedence, Octave will use the method of the object that appears first in the list of arguments.

\subsection*{34.5 Inheritance and Aggregation}

Using classes to build new classes is supported by Octave through the use of both inheritance and aggregation.

Class inheritance is provided by Octave using the class function in the class constructor. As in the case of the polynomial class, the Octave programmer will create a structure that contains the data fields required by the class, and then call the class function to indicate that an object is to be created from the structure. Creating a child of an existing object is done by creating an object of the parent class and providing that object as the third argument of the class function.

This is most easily demonstrated by example. Suppose the programmer needs a FIR filter, i.e., a filter with a numerator polynomial but a denominator of 1 . In traditional Octave programming this would be performed as follows.
```

>> x = [some data vector];
>> n = [some coefficient vector];
>> y = filter (n, 1, x);

```

The equivalent behavior can be implemented as a class @FIRfilter. The constructor for this class is the file FIRfilter.m in the class directory @FIRfilter.
```


## -*- texinfo -*-

## @deftypefn {} {} FIRfilter ()

## @deftypefnx {} {} FIRfilter (@var{p})

## Create a FIR filter with polynomial @var{p} as coefficient vector.

## @end deftypefn

```
```

function f = FIRfilter (p)
if (nargin == 0)
p = @polynomial ([1]);
elseif (! isa (p, "polynomial"))
error ("@FIRfilter: P must be a polynomial object");
endif
f.polynomial = [];
f = class (f, "FIRfilter", p);
endfunction

```

As before, the leading comments provide documentation for the class constructor. This constructor is very similar to the polynomial class constructor, except that a polynomial object is passed as the third argument to the class function, telling Octave that the FIRfilter class will be derived from the polynomial class. The FIR filter class itself does not have any data fields, but it must provide a struct to the class function. Given that the @polynomial constructor will add an element named polynomial to the object struct, the @FIRfilter just initializes a struct with a dummy field polynomial which will later be overwritten.

Note that the sample code always provides for the case in which no arguments are supplied. This is important because Octave will call a constructor with no arguments when loading objects from saved files in order to determine the inheritance structure.

A class may be a child of more than one class (see [class], page 41), and inheritance may be nested. There is no limitation to the number of parents or the level of nesting other than memory or other physical issues.

For the FIRfilter class, more control about the object display is desired. Therefore, the display method rather than the disp method is overloaded (see Section 34.2 [Class Methods], page 891). A simple example might be
```

function display (f)
printf ("%s.polynomial", inputname (1));
disp (f.polynomial);
endfunction

```

Note that the FIRfilter's display method relies on the disp method from the polynomial class to actually display the filter coefficients. Furthermore, note that in the display method it makes sense to start the method with the line printf ( \(" \% \mathrm{~s}="\), inputname (1)) to be consistent with the rest of Octave which prints the variable name to be displayed followed by the value. In general it is not recommended to overload the display function.
display (obj)
Display the contents of the object obj prepended by its name.
The Octave interpreter calls the display function whenever it needs to present a class on-screen. Typically, this would be a statement which does not end in a semicolon to suppress output. For example:
myclass (...)

Or:
```

myobj = myclass (...)

```

In general, user-defined classes should overload the disp method to avoid the default output:
```

myobj = myclass (...)
=> myobj =
<class myclass>

```

When overloading the display method instead, one has to take care of properly displaying the object's name. This can be done by using the inputname function.
See also: [disp], page 271, [class], page 41, [subsref], page 894, [subsasgn], page 896.
Once a constructor and display method exist, it is possible to create an instance of the class. It is also possible to check the class type and examine the underlying structure.
```

octave:1> f = FIRfilter (polynomial ([11 1 1]/3))
f.polynomial = 0.33333 + 0.33333 * X + 0.33333* X ~ 2
octave:2> class (f)
ans = FIRfilter
octave:3> isa (f, "FIRfilter")
ans = 1
octave:4> isa (f, "polynomial")
ans = 1
octave:5> struct (f)
ans =
scalar structure containing the fields:
polynomial = 0.33333 + 0.33333 * X + 0.33333 * X ^ 2

```

The only thing remaining to make this class usable is a method for processing data. But before that, it is usually desirable to also have a way of changing the data stored in a class. Since the fields in the underlying struct are private by default, it is necessary to provide a mechanism to access the fields. The subsref method may be used for both tasks.
```

function r = subsref (f, x)
switch (x.type)
case "()"
n = f.polynomial;
r = filter (n.poly, 1, x.subs{1});
case "."
fld = x.subs;
if (! strcmp (fld, "polynomial"))
error ('@FIRfilter/subsref: invalid property "%s"', fld);
endif

```
```

        r = f.polynomial;
        otherwise
    error ("@FIRfilter/subsref: invalid subscript type for FIR filter");
        endswitch
    endfunction
    ```

The " ()" case allows us to filter data using the polynomial provided to the constructor.
```

octave:2> f = FIRfilter (polynomial ([ll 1 1 1]/3));
octave:3> x = ones (5,1);
octave:4> y = f(x)
y =

```
    0.33333
    0.66667
    1.00000
    1.00000
    1.00000

The "." case allows us to view the contents of the polynomial field.
```

octave:1> f = FIRfilter (polynomial ([11 1 1]/3));
octave:2> f.polynomial
ans = 0.33333 + 0.33333* X + 0.33333* X ~ 2

```

In order to change the contents of the object a subsasgn method is needed. For example, the following code makes the polynomial field publicly writable
```

function fout = subsasgn (f, index, val)
switch (index.type)
case "."
fld = index.subs;
if (! strcmp (fld, "polynomial"))
error ('@FIRfilter/subsasgn: invalid property "%s"', fld);
endif
fout = f;
fout.polynomial = val;
otherwise
error ("@FIRfilter/subsasgn: Invalid index type")
endswitch
endfunction

```
so that
```

octave:1> f = FIRfilter ();
octave:2> f.polynomial = polynomial ([lllll)
f.polynomial = 1 + 2 * X + 3 * X ^ 2

```

Defining the FIRfilter class as a child of the polynomial class implies that a FIRfilter object may be used any place that a polynomial object may be used. This is not a normal use of a filter. It may be a more sensible design approach to use aggregation rather than inheritance. In this case, the polynomial is simply a field in the class structure. A class constructor for the aggregation case might be
```


## -*- texinfo -*-

## @deftypefn {} {} FIRfilter ()

## @deftypefnx {} {} FIRfilter (@var{p})

## Create a FIR filter with polynomial @var{p} as coefficient vector.

## @end deftypefn

function f = FIRfilter (p)
if (nargin == 0)
f.polynomial = @polynomial ([1]);
else
if (! isa (p, "polynomial"))
error ("@FIRfilter: P must be a polynomial object");
endif
f.polynomial = p;
endif
f = class (f, "FIRfilter");
endfunction

```

For this example only the constructor needs changing, and all other class methods stay the same.

\section*{34.6 classdef Classes}

Since version 4.0, Octave has limited support for classdef classes. In contrast to the aforementioned classes, called old style classes in this section, classdef classes can be defined within a single m-file. Other innovations of classdef classes are:
- access rights for properties and methods,
- static methods, i.e., methods that are independent of an object, and
- the distinction between value and handle classes.

Several features have to be added in future versions of Octave to be fully compatible to matlab. An overview of what is missing can be found at https://wiki.octave.org/ Classdef.

\subsection*{34.6.1 Creating a classdef Class}

A very basic classdef value class (see Section 34.6.5 [Value Classes vs. Handle Classes], page 911) is defined by:
```

classdef some_class
properties
endproperties
methods
endmethods
endclassdef

```

In contrast to old style classes, the properties-endproperties block as well as the methods-endmethods block can be used to define properties and methods of the class. Because both blocks are empty, they can be omitted in this particular case.

For simplicity, a more advanced implementation of a classdef class is shown using the polynomial example again (see Section 34.1 [Creating a Class], page 889):
```

classdef polynomial2
properties
poly = 0;
endproperties
methods
function p = polynomial2 (a)
if (nargin == 1)
if (isa (a, "polynomial2"))
p.poly = a.poly;
elseif (isreal (a) \&\& isvector (a))
p.poly = a(:).'; \# force row vector
else
error ("polynomial2: A must be a real vector");
endif
endif
endfunction
function disp (p)
a = p.poly;
first = true;
for i = 1 : length (a);
if (a(i) != 0)
if (first)
first = false;
elseif (a(i) > 0 || isnan (a(i)))
printf (" +");
endif
if (a(i) < 0)
printf (" -");
endif
if (i == 1)
printf (" %.5g", abs (a(i)));
elseif (abs (a(i)) != 1)

```
```

                    printf (" %.5g *", abs (a(i)));
                        endif
                    if (i > 1)
                    printf (" X");
                    endif
                    if (i > 2)
                        printf (" ~ %d", i - 1);
                    endif
                endif
                endfor
                if (first)
                    printf (" 0");
            endif
            printf ("\n");
        endfunction
        endmethods
    endclassdef

```

An object of class polynomial2 is created by calling the class constructor:
```

>> p = polynomial2 ([1, 0, 1])

# p =

    1 + X - 2
    ```

\subsection*{34.6.2 Properties}

All class properties must be defined within properties blocks. The definition of a default value for a property is optional and can be omitted. The default initial value for each class property is [].

A properties block can have additional attributes to specify access rights or to define constants:
```

classdef some_class
properties (Access = mode)
prop1
endproperties
properties (SetAccess = mode, GetAccess = mode)
prop2
endproperties
properties (Constant = true)
prop3 = pi ()
endproperties
properties
prop4 = 1337
endproperties
endclassdef

```
where mode can be one of:
public The properties can be accessed from everywhere.
private The properties can only be accessed from class methods. Subclasses of that class cannot access them.
protected
The properties can only be accessed from class methods and from subclasses of that class.

When creating an object of some_class, prop1 has the default value [] and reading from and writing to prop1 is defined by a single mode. For prop2 the read and write access can be set differently. Finally, prop3 is a constant property which can only be initialized once within the properties block.

By default, in the example prop4, properties are not constant and have public read and write access.
```

properties (class_name)
properties (obj)
plist = properties (...)

```

Return or display the public properties for the named class class_name or classdef object obj.
If an output value is requested, return the list of property names in a cell array.
Programming Note: Property names are returned if the GetAccess attribute is public and if the Hidden attribute is false.

See also: [methods], page 890.

\subsection*{34.6.3 Methods}

All class methods must be defined within methods blocks. An exception to this rule is described at the end of this subsection. Those methods blocks can have additional attributes
specifying the access rights or whether the methods are static, i.e., methods that can be called without creating an object of that class.
```

classdef some_class
methods
function obj = some_class ()
disp ("New instance created.");
endfunction
function disp (obj)
disp ("Here is some_class.");
endfunction
endmethods
methods (Access = mode)
function $r=$ func (obj, r)
$r=2 * r ;$
endfunction
endmethods
methods (Static = true)
function $c=$ circumference (radius)
$c=2$ * pi () .* radius;
endfunction
endmethods
endclassdef

```

The constructor of the class is declared in the methods block and must have the same name as the class and exactly one output argument which is an object of its class.

It is also possible to overload built-in or inherited methods, like the disp function in the example above to tell Octave how objects of some_class should be displayed (see Section 34.2 [Class Methods], page 891).

In general, the first argument in a method definition is always the object that it is called from. Class methods can either be called by passing the object as the first argument to that method or by calling the object followed by a dot (".") and the method's name with subsequent arguments:
```

>> obj = some_class ();
New instance created.
>> disp (obj); \# both are
>> obj.disp (); \# equal

```

In some_class, the method func is defined within a methods block setting the Access attribute to mode, which is one of:
public The methods can be accessed from everywhere.
private The methods can only be accessed from other class methods. Subclasses of that class cannot access them.
protected
The methods can only be accessed from other class methods and from subclasses of that class.

The default access for methods is public.
Finally, the method circumference is defined in a static methods block and can be used without creating an object of some_class. This is useful for methods, that do not depend on any class properties. The class name and the name of the static method, separated by a dot ("."), call this static method. In contrast to non-static methods, the object is not passed as first argument even if called using an object of some_class.
```

>> some_class.circumference (3)
=> ans = 18.850
>> obj = some_class ();
New instance created.
>> obj.circumference (3)
=> ans = 18.850

```

Additionally, class methods can be defined as functions in a folder of the same name as the class prepended with the '@' symbol (see Section 34.1 [Creating a Class], page 889). The main classdef file has to be stored in this class folder as well.

\subsection*{34.6.4 Inheritance}

Classes can inherit from other classes. In this case all properties and methods of the superclass are inherited to the subclass, considering their access rights. Use this syntax to inherit from superclass:
```

classdef subclass < superclass
...
endclassdef

```

\subsection*{34.6.5 Value Classes vs. Handle Classes}

There are two intrinsically different types of classdef classes, whose major difference is the behavior regarding variable assignment. The first type are value classes:
```

classdef value_class
properties
prop1
endproperties
methods
function obj = set_prop1 (obj, val)
obj.prop1 = val;
endfunction
endmethods
endclassdef

```

Assigning an object of that class to another variable essentially creates a new object:
```

>> a = value_class ();
>> a.prop1 = 1;
>> b = a;
>> b.prop1 = 2;
>> b.prop1
=> ans = 2
>> a.prop1
=> ans = 1

```

But that also means that you might have to assign the output of a method that changes properties back to the object manually:
```

>> a = value_class ();
>> a.prop1 = 1;
>> a.set_prop1 (3);
=> ans =
<object value_class>
>> ans.prop1
=> ans = 3
>> a.prop1
mans = 1

```

The second type are handle classes. Those classes have to be derived from the abstract handle class:
```

classdef handle_class < handle
properties
prop1
endproperties
methods
function set_prop1 (obj, val)
obj.prop1 = val;
endfunction
endmethods
endclassdef

```

In the following example, the variables a and b refer to the very same object of class handle_class:
```

>> a = handle_class ();
>> a.prop1 = 1;
>> b = a;
>> b.prop1 = 2;
>> b.prop1
ans = 2
>> a.prop1
=>ans = 2

```

Object properties that are modified by a method of an handle class are changed persistently:
```

>> a.set_prop1 (3);
>> a.prop1
=> ans = 3

```

\section*{35 GUI Development}

Octave is principally a batch or command-line language. However, it does offer some features for constructing graphical interfaces that interact with users.

The GUI elements available are I/O dialogs, a progress bar, and UI elements for plot windows. For example, rather than hardcoding a filename for output results a script can open a dialog box and allow the user to choose a file. Similarly, if a calculation is expected to take a long time a script can display a progress bar. The various UI elements can be used to fully customize the plot window with menubars, toolbars, context menus, pushbuttons, sliders, etc.

Several utility functions make it possible to store private data for use with a GUI which will not pollute the user's variable space.

Finally, a program written in Octave might want to have long term storage of preferences or state variables. This can be done with user-defined preferences.

\subsection*{35.1 I/O Dialogs}

Simple dialog menus are available for choosing directories or files. They return a string variable which can then be used with any command requiring a filename.
```

dirname = uigetdir ()
dirname = uigetdir (init_path)
dirname = uigetdir (init_path, dialog_name)
Open a GUI dialog for selecting a directory.
If init_path is not given the current working directory is used.
dialog_name may be used to customize the dialog title.
The output dirname is a character string with the name of the selected directory. However, if the 'Cancel' button is clicked the output is of type double with the value 0.
See also: [uigetfile], page 915, [uiputfile], page 916.
[fname, fpath, fltidx] = uigetfile ()
[...] = uigetfile (flt)
[...] = uigetfile (flt, dialog_name)
[...] = uigetfile (flt, dialog_name, default_file)
[...] = uigetfile (..., "MultiSelect", mode)

```

Open a GUI dialog for selecting a file and return the filename fname, the path to this file fpath, and the filter index fltidx.
flt contains a (list of) file filter string(s) in one of the following formats:
"/path/to/filename.ext"
If a filename is given then the file extension is extracted and used as filter. In addition, the path is selected as current path in the dialog and the filename is selected as default file. Example: uigetfile ("myfun.m")
A single file extension "*.ext"
Example: uigetfile ("*.ext")

A 2-column cell array
containing a file extension in the first column and a brief description in the second column. Example: uigetfile (\{"*.ext", "My Description";"*.xyz", "XYZ-Format"\})
The filter string can also contain a semicolon separated list of filter extensions. Example: uigetfile (\{"*.gif;*.png;*.jpg", "Supported Picture Formats"\})

A directory name or path name
If the folder name of path name contains a trailing file separator, the contents of that folder will be displayed. If no trailing file separator is present the parent directory is listed. The substring to the right of the rightmost file separator (if any) will be interpreted as a file or directory name and if that file or directory exists it will be highlighted. If the path name or directory name is entirely or partly nonexistent, the current working directory will be displayed. No filter will be active.
dialog_name can be used to customize the dialog title.
If default_file is given then it will be selected in the GUI dialog. If, in addition, a path is given it is also used as current path.

Two or more files can be selected when setting the "MultiSelect" key to "on". In that case, fname is a cell array containing the files.
The outputs fname and fpath are strings returning the chosen name and path, respectively. However, if the 'Cancel' button is clicked the outputs are of type double with a value of 0 . fltidx is the index in the list of filter extensions flt that was selected.

See also: [uiputfile], page 916, [uigetdir], page 915.
[fname, fpath, fltidx] = uiputfile ()
[fname, fpath, fltidx] = uiputfile (flt)
[fname, fpath, fltidx] = uiputfile (flt, dialog_name)
[fname, fpath, fltidx] = uiputfile (flt, dialog_name, default_file)
Open a GUI dialog for selecting a file.
flt contains a (list of) file filter string(s) in one of the following formats:
"/path/to/filename.ext"
If a filename is given the file extension is extracted and used as filter. In addition the path is selected as current path in the dialog and the filename is selected as default file. Example: uiputfile ("myfun.m")
"*.ext" A single file extension. Example: uiputfile ("*.ext")
\{"*.ext", "My Description"\}
A 2-column cell array containing the file extension in the 1st column and a brief description in the 2nd column. Example: uiputfile (\{"*.ext","My Description";"*.xyz", "XYZ-Format"\})

The filter string can also contain a semicolon separated list of filter extensions. Example: uiputfile (\{"*.gif;*.png;*.jpg", "Supported Picture Formats"\})
dialog_name can be used to customize the dialog title. If default_file is given it is preselected in the GUI dialog. If, in addition, a path is given it is also used as current path.
fname and fpath return the chosen name and path, respectively. fltidx is the index in the list of filter extensions flt that was selected.
See also: [uigetfile], page 915, [uigetdir], page 915.
Additionally, there are dialog boxes for displaying help messages, warnings, or errors, and for getting text input from the user.
```

errordlg ()
errordlg (msg)
errordlg (msg, title)
errordlg (msg, title, opt)
h = errordlg (...)

```

Display an error dialog box with error message msg and caption title.
The default error message is "This is the default error string." and the default caption is "Error Dialog".
The error message may have multiple lines separated by newline characters (" n "), or it may be a cellstr array with one element for each line.
The third optional argument opt controls the behavior of the dialog. For details, see [msgbox], page 919.
The return value \(h\) is a handle to the figure object used for building the dialog.
Examples:
```

errordlg ("Some fancy error occurred.");
errordlg ("Some fancy error\nwith two lines.");
errordlg ({"Some fancy error", "with two lines."});
errordlg ("Some fancy error occurred.", "Fancy caption");

```

See also: [helpdlg], page 917, [warndlg], page 921, [msgbox], page 919, [inputdlg], page 918, [listdlg], page 918, [questdlg], page 920.
```

helpdlg ()
helpdlg (msg)
helpdlg (msg, title)
h = helpdlg (...)

```

Display a help dialog box with help message msg and caption title.
The default help message is "This is the default help string." and the default caption is "Help Dialog".
The help message may have multiple lines separated by newline characters (" n "), or it may be a cellstr array with one element for each line.
The return value \(h\) is a handle to the figure object used for building the dialog.
Examples:
```

helpdlg ("Some helpful text for the user.");
helpdlg ("Some helpful text\nwith two lines.");
helpdlg ({"Some helpful text", "with two lines."});
helpdlg ("Some helpful text for the user.", "Fancy caption");

```

See also: [errordlg], page 917, [warndlg], page 921, [msgbox], page 919, [inputdlg], page 918, [listdlg], page 918, [questdlg], page 920.
```

cstr = inputdlg (prompt)
cstr = inputdlg (prompt, title)
cstr = inputdlg (prompt, title, rowscols)
cstr = inputdlg (prompt, title, rowscols, defaults)
cstr = inputdlg (prompt, title, rowscols, defaults, options)

```

Return user input from a multi-textfield dialog box in a cell array of strings, or an empty cell array if the dialog is closed by the Cancel button.

Inputs:
prompt A cell array with strings labeling each text field. This input is required.
title \(\quad\) String to use for the caption of the dialog. The default is "Input Dialog".
rowscols Specifies the size of the text fields and can take three forms:
1. a scalar value which defines the number of rows used for each text field.
2. a vector which defines the individual number of rows used for each text field.
3. a matrix which defines the individual number of rows and columns used for each text field. In the matrix each row describes a single text field. The first column specifies the number of input rows to use and the second column specifies the text field width.
defaults A list of default values to place in each text field. It must be a cell array of strings with the same size as prompt.
options Not supported, only for MATLAB compatibility.
Example:
```

prompt = {"Width", "Height", "Depth"};
defaults = {"1.10", "2.20", "3.30"};
rowscols = [1,10; 2,20; 3,30];
dims = inputdlg (prompt, "Enter Box Dimensions", ...
rowscols, defaults);

```

See also: [errordlg], page 917, [helpdlg], page 917, [listdlg], page 918, [msgbox], page 919, [questdlg], page 920, [warndlg], page 921.
[sel, ok] = listdlg (key, value, ...)
Return user inputs from a list dialog box in a vector of selection indices (sel) and a flag indicating how the user closed the dialog box (ok).
The indices in sel are 1-based.
The value of ok is 1 if the user closed the box with the OK button, otherwise it is 0 and sel is empty.

Input arguments are specified in form of key, value pairs. The "ListString" argument pair must be specified.

Valid key and value pairs are:
"ListString"
a cell array of strings specifying the items to list in the dialog.
"SelectionMode"
can be either "Single" (only one item may be selected at a time) or "Multiple" (default).
"ListSize"
a two-element vector [width, height] specifying the size of the list field in pixels. The default is [160, 300].
"InitialValue"
a vector containing 1 -based indices of elements which will be pre-selected when the list dialog is first displayed. The default is 1 (first item).
"Name" a string to be used as the dialog caption. Default is "".
"PromptString"
a cell array of strings to be displayed above the list of items. Default is \{\}.
"OKString"
a string used to label the OK button. Default is "OK".
"CancelString"
a string used to label the Cancel button. Default is "Cancel".
Example:
```

my_options = {"An item", "another", "yet another"};
[sel, ok] = listdlg ("ListString", my_options,
"SelectionMode", "Multiple");
if (ok == 1)
disp ("You selected:");
for i = 1:numel (sel)
disp (sprintf ("\t%s", my_options{sel(i)}));
endfor
else
disp ("You cancelled.");
endif

```

See also: [menu], page 277, [errordlg], page 917, [helpdlg], page 917, [inputdlg], page 918, [msgbox], page 919, [questdlg], page 920, [warndlg], page 921.
\(h=\operatorname{msgbox}\) (msg)
\(h=\operatorname{msgbox}\) (msg, title)
\(h=\operatorname{msgbox}\) (msg, title, icon)
h = msgbox (msg, title, "custom", cdata)
\(h=\operatorname{msgbox}(m s g, t i t l e, ~ " c u s t o m ", ~ c d a t a, ~ c o l o r m a p) ~\)
\(h=\operatorname{msgbox}(\ldots, o p t)\)
Display msg using a message dialog box.

The message may have multiple lines separated by newline characters (" \(\backslash \mathrm{n}\) "), or it may be a cellstr array with one element for each line.
The optional input title (character string) can be used to decorate the dialog caption.
The optional argument icon selects a dialog icon. It can be one of "none" (default), "error", "help", "warn", or "custom". The latter must be followed by an image array cdata, and for indexed images the associated colormap.
The final optional argument opt controls the behavior of the dialog. If opt is a string, it may be one of
"non-modal" (default)
The dialog is normal.
"modal" If any dialogs already exist with the same title, the most recent is reused and all others are closed. The dialog is displayed "modal" which means it prevents users from interacting with any other GUI element until the dialog has been closed.
"replace"
If any dialogs already exist with the same title, the most recent is reused and all others are closed. The resulting dialog is set "non-modal".

If opt is a structure, it must contain fields "WindowStyle" and "Interpreter":
"WindowStyle"
The value must be "non-modal", "modal", or "replace". See above.
"Interpreter"
Controls the "interpreter" property of the text object used for displaying the message. The value must be "tex" (default), "none", or "latex".

The return value \(h\) is a handle to the figure object used for building the dialog.
Examples:
```

msgbox ("Some message for the user.");
msgbox ("Some message\nwith two lines.");
msgbox ({"Some message", "with two lines."});
msgbox ("Some message for the user.", "Fancy caption");

## A message dialog box with error icon

msgbox ("Some message for the user.", "Fancy caption", "error");

```

See also: [errordlg], page 917, [helpdlg], page 917, [inputdlg], page 918, [listdlg], page 918, [questdlg], page 920, [warndlg], page 921.
```

btn = questdlg (msg)
btn = questdlg (msg, title)
btn = questdlg (msg, title, default)
btn = questdlg (msg, title, btn1, btn2, default)
btn = questdlg (msg, title, btn1, btn2, btn3, default)

```

Display msg using a question dialog box and return the caption of the activated button.

The message may have multiple lines separated by newline characters (" n "), or it may be a cellstr array with one element for each line.
The optional title (character string) can be used to specify the dialog caption. It defaults to "Question Dialog".
The dialog may contain two or three buttons which will all close the dialog.
The string default identifies the default button, which is activated by pressing the ENTER key. It must match one of the strings given in btn1, btn2, or btn3.
If only msg and title are specified, three buttons with the default captions "Yes", "No", and "Cancel" are used.
If only two button captions, btn1 and btn2, are specified the dialog will have only these two buttons.

Examples:
```

btn = questdlg ("Close Octave?", "Some fancy title", ...
"Yes", "No", "No");
if (strcmp (btn, "Yes"))
exit ();
endif

```

See also: [errordlg], page 917, [helpdlg], page 917, [inputdlg], page 918, [listdlg], page 918, [msgbox], page 919, [warndlg], page 921.
warndlg ()
warndlg (msg)
warndlg (msg, title)
warndlg (msg, title, opt)
h = warndlg (...)
Display a warning dialog box with warning message msg and caption title.
The default warning message is "This is the default warning string." and the default caption is "Warning Dialog".
The warning message may have multiple lines separated by newline characters ("\n"), or it may be a cellstr array with one element for each line.
The third optional argument opt controls the behavior of the dialog. For details, see [msgbox], page 919.
The return value \(h\) is a handle to the figure object used for building the dialog.
Examples:
```

warndlg ("Some warning text for the user.");
warndlg ("Some warning text\nwith two lines.");
warndlg ({"Some warning text", "with two lines."});
warndlg ("Some warning text for the user.", "Fancy caption");

```

See also: [errordlg], page 917, [helpdlg], page 917, [msgbox], page 919, [inputdlg], page 918, [listdlg], page 918, [questdlg], page 920.
uisetfont ()
uisetfont (h)
uisetfont (fontstruct)
uisetfont (..., title)
fontstruct \(=\) uisetfont (...)
Open a font selection dialog.
If the first argument is a handle to a text, axes, or uicontrol object, pressing the OK button will change the font properties of the object.
The first argument may also be a structure with fields FontName, FontWeight, FontAngle, FontUnits, and FontSize, indicating the initially selected font.
The title of the dialog window can be specified by using the last argument title.
If an output argument fontstruct is requested, the selected font structure is returned. Otherwise, the font information is displayed onscreen.
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache -fv") manually after installing new fonts.

See also: [listfonts], page 934, [text], page 392, [axes], page 423, [uicontrol], page 925.
For creating new dialog types, there is a dialog function.
```

h = dialog ()
h = dialog ("property", value, ...)

```

Create an empty modal dialog window to which other uicontrols can be added.
The dialog box is a figure object with properties as recommended for a dialog box.
The default properties differing from a figure are:
buttondownfen
if isempty (allchild(gcbf)), close (gcbf), endif
colormap []
color defaultuicontrolbackgroundcolor
dockcontrols
off
handlevisibility
callback
integerhandle
off
inverthardcopy
off
menubar none
numbertitle
off
paperpositionmode
auto
resize off
windowstyle
modal
Multiple property-value pairs may be specified for the dialog object, but they must appear in pairs. The full list of properties is documented at Section 15.3.3.2 [Figure Properties], page 434.
The return value \(h\) is a graphics handle to the created figure.
Example:
```


## create an empty dialog window titled "Dialog Example"

h = dialog ("name", "Dialog Example");

## create a button (default style)

b = uicontrol (h, "string", "OK",
"position", [10 10 150 40],
"callback", "delete (gcf)");

## wait for dialog to resume or close

uiwait (h);

```

See also: [errordlg], page 917, [msgbox], page 919, [questdlg], page 920, [warndlg], page 921, [figure], page 399, [uiwait], page 936.

\subsection*{35.2 Progress Bar}
```

h = waitbar (frac)
h = waitbar (frac, msg)
h = waitbar (..., "createcancelbtn", fcn, ...)
h = waitbar (..., prop, val, ...)
waitbar (frac)
waitbar (frac, h)
waitbar (frac, h, msg)

```

Return a handle \(h\) to a new progress indicator ("waitbar") object.
The waitbar is filled to fraction frac which must be in the range \([0,1]\).
The optional message msg is centered and displayed above the waitbar.
A cancel button can be added to the bottom of the waitbar using the "createcancelbtn" property of waitbar figures. The action to be executed when the user presses the button is specified using a string or function handle fcn.
The appearance of the waitbar figure window can be configured by passing prop/val pairs to the function. The full list of properties is documented at Section 15.3.3.2 [Figure Properties], page 434.

When called with a single input the current waitbar, if it exists, is updated to the new value frac. If there are multiple outstanding waitbars they can be updated individually by passing the handle \(h\) of the specific waitbar to modify.

See also: [delete], page 405.

\subsection*{35.3 UI Elements}

The ui* series of functions work best with the qt graphics toolkit, although some functionality is available with the fltk toolkit. There is no support for the gnuplot toolkit.
```

uipanel (property, value,...)
uipanel (parent, property, value,...)
hui = uipanel (...)

```

Create a uipanel object.
uipanels are used as containers to group other uicontrol objects.
If parent is omitted then a uipanel for the current figure is created. If no figure is available, a new figure is created first.
If parent is given then a uipanel relative to parent is created.
Any provided property value pairs will override the default values of the created uipanel object.
The full list of properties is documented at Section 15.3.3.15 [Uipanel Properties], page 476.
The optional return value hui is a graphics handle to the created uipanel object.
Examples:
```


## create figure and panel on it

f = figure;
p = uipanel ("title", "Panel Title", "position", [. 25 . 25 . 5 . 5]);

## add two buttons to the panel

b1 = uicontrol ("parent", p, "string", "A Button", ...
"position", [18 10 150 36]);
b2 = uicontrol ("parent", p, "string", "Another Button", ...
"position",[18 60 150 36]);

```

See also: [figure], page 399, [uicontrol], page 925.
hui = uibuttongroup (property, value, ...)
hui = uibuttongroup (parent, property, value, ...)
uibuttongroup (h)
Create a uibuttongroup object and return a handle to it.
A uibuttongroup is used to group uicontrol objects.
If parent is omitted then a uibuttongroup for the current figure is created. If no figure is available, a new figure is created first.
If parent is given then a uibuttongroup relative to parent is created.
Any provided property value pairs will override the default values of the created uibuttongroup object.
The full list of properties is documented at Section 15.3.3.13 [Uibuttongroup Properties], page 472.

\section*{Examples:}
```


## Create figure and panel on it

f = figure;

## Create a button group

gp = uibuttongroup (f, "Position", [ 0 0.5 1 1])

## Create a buttons in the group

b1 = uicontrol (gp, "style", "radiobutton", ...
"string", "Choice 1", ...
"Position", [ 10 150 100 50 ]);
b2 = uicontrol (gp, "style", "radiobutton", ...
"string", "Choice 2", ...
"Position", [ 10 50 100 30 ]);

## Create a button not in the group

b3 = uicontrol (f, "style", "radiobutton", ...
"string", "Not in the group", ...
"Position", [ 10 50 100 50 ]);

```

When called with a single argument \(h\) which is a handle to an existing uibuttongroup object, switch the focus to the specified uibuttongroup. This functionality is not currently implemented.
See also: [figure], page 399, [uipanel], page 924.
hui \(=\) uicontrol (property, value, ...)
hui \(=\) uicontrol (parent, property, value,...)
uicontrol (h)
Create a uicontrol object and return a handle to it.
A uicontrol object is used to create simple interactive controls such as push buttons, checkboxes, edit and list controls.
If parent is omitted then a uicontrol for the current figure is created. If no figure is available, a new figure is created first.
If parent is given then a uicontrol relative to parent is created.
Any provided property value pairs will override the default values of the created uicontrol object.
The full list of properties is documented at Section 15.3.3.16 [Uicontrol Properties], page 478.
The type of uicontrol created is specified by the style property. If no style property is provided, a push button will be created.
Valid styles for uicontrol are:
"checkbox"
Create a checkbox control that allows user on/off selection.
"edit" Create an edit control that allows user input of single or multiple lines of text.
"listbox"
Create a listbox control that displays a list of items and allows user selection of single or multiple items.
"popupmenu"
Create a popupmenu control that displays a list of options that can be selected when the user clicks on the control.
"pushbutton"
Create a push button control that allows user to press to cause an action.
"radiobutton"
Create a radio button control intended to be used for mutually exclusive input in a group of radiobutton controls.
"slider" Create a slider control that allows user selection from a range of values by sliding knob on the control.
"text" Create a static text control to display single or multiple lines of text.
"togglebutton"
Create a toggle button control that appears like a push button but allows the user to select between two states.

Examples:
```


## Create figure and panel on it

f = figure;

## Create a button (default style)

b1 = uicontrol (f, "string", "A Button", ...
"position", [10 10 150 40]);

## Create an edit control

e1 = uicontrol (f, "style", "edit", "string", "editable text", ...
"position", [10 60 300 40]);

## Create a checkbox

c1 = uicontrol (f, "style", "checkbox", "string", "a checkbox", ...
"position", [10 120 150 40]);

```

When called with a single argument \(h\) which is a handle to an existing uicontrol object, switch the keyboard focus to the specified uicontrol. As a result, the uicontrol object will receive keyboard events that can be processed using the "keypressfcn" callback.

See also: [figure], page 399, [uipanel], page 924.
hui = uitable (property, value, ...)
hui \(=\) uitable (parent, property, value, ...)
Create a uitable object and return a handle to it.
A uitable object is used to show tables of data in a figure window.
If parent is omitted then a uitable for the current figure is created. If no figure is available, a new figure is created first.
If parent is given then a uitable relative to parent is created.
Any provided property value pairs will override the default values of the created uitable object.
The full list of properties is documented at Section 15.3.3.17 [Uitable Properties], page 481.

Examples:
```


## Create figure and place a table on it

f = figure ();
m = magic (8);
t = uitable (f, "Data", m, "ColumnWidth", { 40 });

## Create a table with labeled rows and columns

f = figure ();
d = reshape (1:9, [3, 3]);
row_names = { "Row1", "Row2", "Row3" };
col_names = { "Col1", "Col2", "Col3" };
t = uitable (f, "Data", d, ...
"RowName", row_names, "ColumnName", col_names);
p = get (t, "Position");
e = get (t, "Extent");
p(3:4) = e(3:4);
set (t, "Position", p);

## Long demo with callbacks

function uitable_demo ()
f = figure ("Name", "uitable Demo", "Menu", "none", ...
"Position", [10 10 1000 680]);
\#\# A basic example
d = { "char" , "A string";
"double" , 12.3456789;
"complex", 1+2i;
"bool" , true;
"single" , single(12.3456789);
"int8" , int8(-128);
"uint8" , uint8(128);
"int16" , int16(-32768);
"uint16" , uint16(32768);
"int32" , int32(-2147483648);
"uint32" , uint32(2147483648);
"int64" , int64(-2147483649);
"uint64" , uint64(2147843649)};
popup_options = {"A", "B", "C", "D", "E"};
columnformat_options = { "[]", "char", "pop-up", "numeric", ...
"short", "short e", "short eng", ...
"short g", "long", "long e", ...
"long eng", "long g", "bank", "+", ...
"rat", "logical"};

```
```

columnformat_values = columnformat_options;
columnformat_values{1} = "";
columnformat_values{3} = popup_options;
default_data = repmat (d(:,2), 1, columns (columnformat_options));
b_add = uicontrol (f, "Position", [285 630 600 50], ...
"UserData", [rows(d), 1], ...
"Style", "pushbutton", ...
"String", "Set data at selected point to selected datatype");
l_type_table = uicontrol (f, "Position", [ 0 603 120 25 ], ...
"String", "Datatype Table:", ...
"Style", "text");
t_type_table = uitable (f, "Position", [ 0 530 1000 70 ], ...
"Data", transpose (d(:, 2)), ...
"ColumnName", transpose (d(:, 1)), ...
"RowName", "Value", ...
"CellSelectionCallback", ...
@(x, y) set (b_add, "UserData", y.Indices ));
l_point_table = uicontrol (f, "Position", [ 0 640 60 25 ], ...
"String", "Point:", ...
"Style", "text");
t_point_table = uitable (f, "Position", [ 80 630 160 42 ], ...
"RowName", [], ...
"ColumnName", {"x", "y"}, ...
"Data", [ 1, 1 ], ...
"ColumnEditable", true);
l_editable_table = uicontrol (f, "Position", [ 0 502 200 25 ], ...
"Style", "text", ...
"String", "Set Data Columns Editable:");
t_editable_table = ...
uitable (f, "Position", [ 0 434 1000 65 ], ...
"Data", repmat (false, 1, columns (default_data)), ...
"ColumnEditable", true);
l_format_table = uicontrol (f, "Position", [ 0 406 200 25 ], ...
"Style", "text", ...
"String", "Set Data Column Format:");
t_format_table = ...
uitable (f, "Position", [ 0 338 1000 65 ], ...
"Data", columnformat_options, ...
"ColumnEditable", true, ...
"ColumnFormat", arrayfun (@(x) {columnformat_options}, ...|
1:columns (columnformat_options)));

```
```

l_data_table = uicontrol (f, "Style", "text", ...
"String", "Data:", ...
"Position", [ 0 310 60 25 ]);
t_data_table = uitable (f, "Position", [ 0 15 1000 290 ], ...
"Data", default_data, ...
"ColumnFormat", columnformat_values);
set (t_format_table, ...
"CellEditCallback", ...
@(x, y) update_column_format (y.NewData, y.Indices, ...
t_data_table, popup_options));
set (t_point_table, "CellEditCallback", ...
@(x, y) validate_point_table (x, y, t_data_table));
set (t_editable_table, "CellEditCallback", ...
@(x,y) set (t_data_table, ...
"ColumnEditable", get (t_editable_table, "Data")));
set (b_add, ...
"Callback", @(x, y) update_data (b_add, t_point_table, ...
t_type_table, t_data_table));
set (t_data_table, "CellSelectionCallback", ...
@(x, y) update_point_table (y.Indices, t_point_table));
endfunction
function validate_point_table (h, dat, t_data_table)
if (! (dat.NewData > 0 \&\& ...
dat.NewData < size (get (t_data_table, "Data"), dat.Indices(1, 1)) + 1))|
d = get (h, "Data");
d(dat.Indices) = 1;
set (h, "Data", d);
endif
endfunction
function update_column_format (format, indices, t_data_table, ...
popup_options)
cf = get (t_data_table, "ColumnFormat");
if (strcmp (format, "[]"))
format = "";
elseif (strcmp (format, "pop-up"))
format = popup_options;
endif
cf{indices(1,2)} = format;
set (t_data_table, "ColumnFormat", cf);
endfunction

```
```

function update_point_table (indices, t_point_table)
if (isempty (indices))
indices = [1, 1];
endif
set (t_point_table, "Data", indices(1,:));
endfunction
function update_data (b_add, t_point_table, t_type_table, ...
t_data_table)
indices = get (b_add, "UserData");
if (isempty (indices))
indices = [1, 1];
endif
d = get (t_data_table, "Data");
t_type_table_data = get (t_type_table, "Data");
p = get (t_point_table, "Data");
d(p(1,2), p(1,1)) = t_type_table_data(indices(1,2));
set (t_data_table, "Data", d);
endfunction

```

See also: [figure], page 399, [uicontrol], page 925.
hui \(=\) uimenu (property, value, ...)
hui = uimenu (h, property, value, ...)
Create a uimenu object and return a handle to it.
If \(h\) is omitted then a top-level menu for the current figure is created. If \(h\) is given then a submenu relative to \(h\) is created.
uimenu objects have the following specific properties:
"accelerator"
A string containing the key combination together with CTRL to execute this menu entry (e.g., "x" for CTRL+x).
"callback"
Is the function called when this menu entry is executed. It can be either a function string (e.g., "myfun"), a function handle (e.g., @myfun) or a cell array containing the function handle and arguments for the callback function (e.g., \{@myfun, arg1, arg2\}).
"checked"
Can be set "on" or "off". Sets a mark at this menu entry.
"enable" Can be set "on" or "off". If disabled the menu entry cannot be selected and it is grayed out.
"foregroundcolor"
A color value setting the text color for this menu entry.
"label" A string containing the label for this menu entry. A "\&"-symbol can be used to mark the "accelerator" character (e.g., "E\&xit")
```

"position"

```

An scalar value containing the relative menu position. The entry with the lowest value is at the first position starting from left or top.
"separator"
Can be set "on" or "off". If enabled it draws a separator line above the current position. It is ignored for top level entries.

The full list of properties is documented at Section 15.3.3.12 [Uimenu Properties], page 470.
Examples:
```

f = uimenu ("label", "\&File", "accelerator", "f");
e = uimenu ("label", "\&Edit", "accelerator", "e");
uimenu (f, "label", "Close", "accelerator", "q", ...
"callback", "close (gcf)");
uimenu (e, "label", "Toggle \&Grid", "accelerator", "g", ...
"callback", "grid (gca)");

```

See also: [figure], page 399.
hui \(=\) uicontextmenu (property, value, ...)
hui = uicontextmenu (h, property, value, ...)
Create a uicontextmenu object and return a handle to it.
If \(h\) is omitted then a uicontextmenu for the current figure is created. If no figure is available, a new figure is created first.
If \(h\) is given then a uicontextmenu relative to \(h\) is created.
Any provided property value pairs will override the default values of the created uicontextmenu object.
The full list of properties is documented at Section 15.3.3.14 [Uicontextmenu Properties], page 474.
Examples:
```


## create figure and uicontextmenu

f = figure ();
c = uicontextmenu (f);

## create menus in the context menu

m1 = uimenu ("parent", c, "label", "Menu item 1", ...
"callback", "disp('menu item 1')");
m2 = uimenu ("parent", c, "label", "Menu item 2", ...
"callback", "disp('menu item 2')");

## set the context menu for the figure

set (f, "uicontextmenu", c);

```

See also: [figure], page 399, [uimenu], page 930.
uitoolbar (property, value, ...)
uitoolbar (parent, property, value, ...)
hui = uitoolbar (...)
Create a uitoolbar object. A uitoolbar displays uitoggletool and uipushtool buttons. If parent is omitted then a uitoolbar for the current figure is created. If no figure is available, a new figure is created first.
If parent is given then a uitoolbar relative to parent is created.
Any provided property value pairs will override the default values of the created uitoolbar object.
The full list of properties is documented at Section 15.3.3.18 [Uitoolbar Properties], page 483.
The optional return value hui is a graphics handle to the created uitoolbar object.
Examples:
\% create figure without a default toolbar
f = figure ("toolbar", "none");
\% create empty toolbar
t = uitoolbar (f);
See also: [figure], page 399, [uitoggletool], page 933, [uipushtool], page 932.
uipushtool (property, value, ...)
uipushtool (parent, property, value, ...)
hui \(=\) uipushtool (...)
Create a uipushtool object.
uipushtools are buttons that appear on a figure toolbar. The button is created with a border that is shown when the user hovers over the button. An image can be set using the cdata property.
If parent is omitted then a uipushtool for the current figure is created. If no figure is available, a new figure is created first. If a figure is available, but does not contain a uitoolbar, a uitoolbar will be created.

If parent is given then a uipushtool is created on the parent uitoolbar.
Any provided property value pairs will override the default values of the created uipushtool object.
The full list of properties is documented at Section 15.3.3.19 [Uipushtool Properties], page 485.

The optional return value hui is a graphics handle to the created uipushtool object.
Examples:
```

% create figure without a default toolbar
f = figure ("toolbar", "none");
% create empty toolbar
t = uitoolbar (f);
% create a 19x19x3 black square
img=zeros(19,19,3);
% add pushtool button to toolbar
b = uipushtool (t, "cdata", img);

```

See also: [figure], page 399, [uitoolbar], page 931, [uitoggletool], page 933.
```

uitoggletool (property, value,...)
uitoggletool (parent, property, value,...)
hui = uitoggletool (...)
Create a uitoggletool object.

```
uitoggletool are togglebuttons that appear on a figure toolbar. The button is created with a border that is shown when the user hovers over the button. An image can be set using the cdata property.

If parent is omitted then a uitoggletool for the current figure is created. If no figure is available, a new figure is created first. If a figure is available, but does not contain a uitoolbar, a uitoolbar will be created.

If parent is given then a uitoggletool is created on the parent uitoolbar.
Any provided property value pairs will override the default values of the created uitoggletool object.
The full list of properties is documented at Section 15.3.3.20 [Uitoggletool Properties], page 486.

The optional return value hui is a graphics handle to the created uitoggletool object.
Examples:
```

% create figure without a default toolbar
f = figure ("toolbar", "none");
% create empty toolbar
t = uitoolbar (f);
% create a 19x19x3 black square
img=zeros(19,19,3);
% add uitoggletool button to toolbar
b = uitoggletool (t, "cdata", img);

```

See also: [figure], page 399, [uitoolbar], page 931, [uipushtool], page 932.

\subsection*{35.4 GUI Utility Functions}

These functions do not implement a GUI element but are useful when developing programs that do. The functions uiwait, uiresume, and waitfor are only available with the qt or fltk toolkits.
data \(=\) guidata ( \(h\) )
guidata (h, data)
Query or set user-custom GUI data.
The GUI data is stored in the figure handle \(h\). If \(h\) is not a figure handle then it's parent figure will be used for storage.
data must be a single object which means it is usually preferable for it to be a data container such as a cell array or struct so that additional data items can be added easily.

See also: [getappdata], page 495, [setappdata], page 494, [get], page 430, [set], page 430, [getpref], page 937, [setpref], page 937.
hdata = guihandles (h)
hdata = guihandles
Return a structure of object handles for the figure associated with handle \(h\).
If no handle is specified the current figure returned by gcf is used.
The fieldname for each entry of hdata is taken from the "tag" property of the graphic object. If the tag is empty then the handle is not returned. If there are multiple graphic objects with the same tag then the entry in hdata will be a vector of handles. guihandles includes all possible handles, including those for which "HandleVisibility" is "off".

See also: [guidata], page 933, [findobj], page 488, [findall], page 489, [allchild], page 431.
have_window_system ()
Return true if a window system is available (X11, Windows, or Apple OS X) and false otherwise.

See also: [isguirunning], page 934.
isguirunning ()
Return true if Octave is running in GUI mode and false otherwise.
See also: [have_window_system], page 934.
pos = getpixelposition (h)
pos = getpixelposition (h, rel_to_fig)
Return the position of a user interface component in pixel units.
The first argument \(h\) must be a handle to a valid graphics object of type uibuttongroup, uicontrol, uipanel, uitable, axes, or figure. For other object types, the function returns zeros.
By default, the position is returned relative to the object's parent. If the second argument rel_to_fig is logically true, the position is computed relative to the enclosing figure object.
The return value pos is a 4 -element vector with values [ lower_left_X, lower_left_ Y, width, height ].
See also: [get], page 430.
fonts = listfonts ()
fonts = listfonts (h)
List system fonts.
If a handle to a graphics object \(h\) is provided, also include the font from the object's "FontName" property in the list.
Programming Note: On systems that don't use FontConfig natively (all but Linux), the font cache is built when Octave is installed. You will need to run system ("fc-cache \(-f v "\) ) manually after installing new fonts.

See also: [uisetfont], page 921, [text], page 392, [axes], page 423, [uicontrol], page 925.
```

movegui
movegui (h)
movegui (pos)
movegui (h, pos)
movegui (h, event)
movegui (h, event, pos)

```

Move a figure specified by figure handle \(h\) to a position on the screen defined by pos.
\(h\) is a figure handle, or a handle to a graphics object. In the latter case, its parent figure will be used. If unspecified, \(h\) will be set to the handle of the relevant figure if a callback is being executed (gcbf), otherwise it will be set to the handle of the current figure (gcf).
pos is either a two-value numeric vector or a string. If pos is numeric then it must be of the form [h, v] specifying the horizontal and vertical offsets of the figure with respect to the screen. A positive value indicates the offset between the left (or bottom for the vertical component) of the screen, and the left (or bottom) of the figure. A negative value indicates the offset between the right (or top) of the screen and the right (or top) of the figure.
Possible values for pos as a string are
north Top center of the screen.
south Bottom center of the screen.
east Right center of the screen.
west Left center of the screen.
northeast
Top right of the screen.
northwest
Top left of the screen.
southeast
Bottom right of the screen.
southwest
Bottom left of the screen.
center Center of the screen.
onscreen (default)
The figure will be minimally moved to be entirely visible on the screen, with a 30 pixel extra padding from the sides of the screen. This is the default value if none is provided.
event contains event data that will be ignored. This construct facilitates a call to movegui from a callback.
openvar (name)
Open the variable name in the graphical Variable Editor.
uiwait
uiwait (h)
uiwait (h, timeout)
Suspend program execution until the figure with handle \(h\) is deleted or uiresume is called.

When no figure handle is specified this function uses the current figure. If the figure handle is invalid or there is no current figure, this functions returns immediately.

When specified, timeout defines the number of seconds to wait for the figure deletion or the uiresume call. The timeout value must be at least 1. If a smaller value is specified, a warning is issued and a timeout value of 1 is used instead. If a noninteger value is specified, it is truncated towards 0 . If timeout is not specified, the program execution is suspended indefinitely.

See also: [uiresume], page 936, [waitfor], page 936.
uiresume (h)
Resume program execution suspended with uiwait.
The handle \(h\) must be the same as the on specified in uiwait. If the handle is invalid or there is no uiwait call pending for the figure with handle \(h\), this function does nothing.
See also: [uiwait], page 936.

\section*{waitfor (h)}
waitfor (h, prop)
waitfor (h, prop, value)
waitfor (..., "timeout", timeout)
Suspend the execution of the current program until a condition is satisfied on the graphics handle \(h\).
While the program is suspended graphics events are still processed normally, allowing callbacks to modify the state of graphics objects. This function is reentrant and can be called from a callback, while another waitfor call is pending at the top-level.
In the first form, program execution is suspended until the graphics object \(h\) is destroyed. If the graphics handle is invalid or if \(h\) is the root graphics handle and no property prop was provided, the function returns immediately.
In the second form, execution is suspended until the graphics object is destroyed or the property named prop is modified. If the graphics handle is invalid or the property does not exist, the function returns immediately.
In the third form, execution is suspended until the graphics object is destroyed or the property named prop is set to value. The function isequal is used to compare property values. If the graphics handle is invalid, the property does not exist or the property is already set to value, the function returns immediately.
An optional timeout can be specified using the property "timeout". This timeout value is the number of seconds to wait for the condition to be true. timeout must be at least 1. If a smaller value is specified, a warning is issued and a value of 1 is used instead. If the timeout value is not an integer, it is truncated towards 0 .

To define a condition on a property named "timeout", use the string '\timeout' instead.
In all cases, typing CTRL-C stops program execution immediately.
See also: [waitforbuttonpress], page 419, [isequal], page 165.

\subsection*{35.5 User-Defined Preferences}
```

val = getpref ("group", "pref")
val = getpref ("group", "pref", default)
{val1, val2, ...} = getpref ("group", {"pref1", "pref2", ...})
prefstruct = getpref ("group")
prefstruct = getpref ()

```

Return the preference value corresponding to the named preference pref in the preference group group.
The named preference group must be a string.
If pref does not exist in group and default is specified, create the preference with value default and return default.
The preference pref may be a string or cell array of strings. If it is a cell array of strings then a cell array of preferences is returned.
The corresponding default value default may be any Octave value, .e.g., double, struct, cell array, object, etc. Or, if pref is a cell array of strings then default must be a cell array of values with the same size as pref.
If neither pref nor default are specified, return a structure of preferences for the preference group group.
If no arguments are specified, return a structure containing all groups of preferences and their values.
See also: [addpref], page 937, [setpref], page 937, [ispref], page 938, [rmpref], page 938.
setpref ("group", "pref", val)
setpref ("group", \{"pref1", "pref2", ...\}, \{val1, val2, ...\})
Set the preference pref to the given val in the named preference group group.
The named preference group must be a string.
The preference pref may be a string or a cell array of strings.
The corresponding value val may be any Octave value, .e.g., double, struct, cell array, object, etc. Or, if pref is a cell array of strings then val must be a cell array of values with the same size as pref.
If the named preference or group does not exist, it is added.
See also: [addpref], page 937, [getpref], page 937, [ispref], page 938, [rmpref], page 938.
addpref ("group", "pref", val)
addpref ("group", \{"pref1", "pref2", ...\}, \{val1, val2, ...\})
Add the preference pref and associated value val to the named preference group group.

The named preference group must be a string.
The preference pref may be a string or a cell array of strings. An error will be issued if the preference already exists.
The corresponding value val may be any Octave value, .e.g., double, struct, cell array, object, etc. Or, if pref is a cell array of strings then val must be a cell array of values with the same size as pref.
See also: [setpref], page 937, [getpref], page 937, [ispref], page 938, [rmpref], page 938.
rmpref ("group", "pref")
rmpref ("group", \{"pref1", "pref2", ...\})
rmpref ("group")
Remove the named preference pref from the preference group group.
The named preference group must be a string.
The preference pref may be a string or cell array of strings.
If pref is not specified, remove the preference group group.
It is an error to remove a nonexistent preference or group.
See also: [addpref], page 937, [ispref], page 938, [setpref], page 937, [getpref], page 937.
ispref ("group", "pref")
ispref ("group", \{"pref1", "pref2", ...\})
ispref ("group")
Return true if the named preference pref exists in the preference group group.
The named preference group must be a string.
The preference pref may be a string or a cell array of strings.
If pref is not specified, return true if the preference group group exists.
See also: [getpref], page 937, [addpref], page 937, [setpref], page 937, [rmpref], page 938.
prefdir
prefdir (1)
dir = prefdir
Return the directory that holds the preferences for Octave.
Examples:
Display the preferences directory
prefdir
Change to the preferences folder
cd (prefdir)
If called with an argument, the preferences directory is created if it doesn't already exist.

See also: [getpref], page 937, [setpref], page 937, [addpref], page 937, [rmpref], page 938, [ispref], page 938.
preferences
Display the GUI preferences dialog window for Octave.

\subsection*{35.6 Octave Workspace Windows}

The functions below make windows that are a part of Octave's own GUI interface visible, and move the keyboard focus to the selected window. Their utility lies in the ability to call these functions from a script and highlight a window without using a mouse for selection.
commandhistory ()
Show the GUI command history window and give it the keyboard focus.
See also: [commandwindow], page 939, [filebrowser], page 939, [workspace], page 939.
commandwindow ()
Show the GUI command window and give it the keyboard focus.
See also: [commandhistory], page 939, [filebrowser], page 939, [workspace], page 939.
filebrowser ()
Show the GUI file browser window and give it the keyboard focus.
See also: [commandwindow], page 939, [commandhistory], page 939, [workspace], page 939.
workspace ()
Show the GUI workspace window and give it the keyboard focus.
See also: [commandwindow], page 939, [commandhistory], page 939, [filebrowser], page 939 .

\section*{36 System Utilities}

This chapter describes the functions that are available to allow you to get information about what is happening outside of Octave, while it is still running, and use this information in your program. For example, you can get information about environment variables, the current time, and even start other programs from the Octave prompt.

\subsection*{36.1 Timing Utilities}

Octave's core set of functions for manipulating time values are patterned after the corresponding functions from the standard C library. Several of these functions use a data structure for time that includes the following elements:
usec Microseconds after the second (0-999999).
sec Seconds after the minute (0-60). This number can be 60 to account for leap seconds.
min Minutes after the hour (0-59).
hour Hours since midnight (0-23).
mday \(\quad\) Day of the month (1-31).
mon Months since January (0-11).
year Years since 1900.
wday Days since Sunday (0-6).
yday Days since January 1 (0-365).
isdst Daylight saving time flag.
gmtoff Seconds offset from UTC.
zone Time zone.
In the descriptions of the following functions, this structure is referred to as a tm_struct.
seconds = time ()
Return the current time as the number of seconds since the epoch.
The epoch is referenced to 00:00:00 UTC (Coordinated Universal Time) 1 Jan 1970. For example, on Monday February 17, 1997 at 07:15:06 UTC, the value returned by time was 856163706 .
See also: [strftime], page 943, [strptime], page 945, [localtime], page 942, [gmtime], page 942, [mktime], page 943, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.
\(\mathrm{t}=\) now ()
Return the current local date/time as a serial day number (see [datenum], page 948).
The integral part, floor (now) corresponds to the number of days between today and Jan 1, 0000.
The fractional part, rem (now, 1) corresponds to the current time.
See also: [clock], page 945, [date], page 946, [datenum], page 948.
ctime ( \(t\) )
Convert a value returned from time (or any other non-negative integer), to the local time and return a string of the same form as asctime.

The function ctime (time) is equivalent to asctime (localtime (time)). For example:
```

ctime (time ())
=> "Mon Feb 17 01:15:06 1997\n"

```

See also: [asctime], page 943, [time], page 941, [localtime], page 942.

\section*{tm_struct \(=\) gmtime \((t)\)}

Given a value returned from time, or any non-negative integer, return a time structure corresponding to UTC (Coordinated Universal Time).

For example:
```

gmtime (time ())
=> {
usec = 0
sec = 6
min = 15
hour = 7
mday = 17
mon = 1
year = 97
wday = 1
yday = 47
isdst = 0
gmtoff = 0
zone = GMT
}

```

See also: [strftime], page 943, [strptime], page 945, [localtime], page 942, [mktime], page 943 , [time], page 941, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.

Given a value returned from time, or any non-negative integer, return a time structure corresponding to the local time zone.
```

localtime (time ())
=> {
usec = 0
sec = 6
min = 15
hour = 1
mday = 17
mon = 1
year = 97
wday = 1
yday = 47
isdst = 0
gmtoff = -21600
zone = CST
}

```

See also: [strftime], page 943, [strptime], page 945, [gmtime], page 942, [mktime], page 943, [time], page 941, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.
seconds \(=\) mktime (tm_struct)
Convert a time structure corresponding to the local time to the number of seconds since the epoch.
For example:
```

mktime (localtime (time ()))
\# 856163706

```

See also: [strftime], page 943, [strptime], page 945, [localtime], page 942, [gmtime], page 942, [time], page 941, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.
asctime (tm_struct)
Convert a time structure to a string using the following format: "ddd mmm m HH:MM:SS yyyy \(\backslash n "\).
For example:
```

asctime (localtime (time ()))
=> "Mon Feb 17 01:15:06 1997\n"

```

This is equivalent to ctime (time ()).
See also: [ctime], page 942, [localtime], page 942, [time], page 941.
strftime (fmt, tm_struct)
Format the time structure tm_struct in a flexible way using the format string fmt that contains ' \(\%\) ' substitutions similar to those in printf.
Except where noted, substituted fields have a fixed size; numeric fields are padded if necessary. Padding is with zeros by default; for fields that display a single number,
padding can be changed or inhibited by following the \(\% \%\) with one of the modifiers described below. Unknown field specifiers are copied as normal characters. All other characters are copied to the output without change. For example:
```

strftime ("%r (%Z) %A %e %B %Y", localtime (time ()))
=> "01:15:06 AM (CST) Monday 17 February 1997"

```

Octave's strftime function supports a superset of the ANSI C field specifiers.
Literal character fields:
\%\% \% character.
\%n Newline character.
\(\% \mathrm{t}\) Tab character.
Numeric modifiers (a nonstandard extension):
- (dash) Do not pad the field.
_ (underscore)
Pad the field with spaces.
Time fields:
\%H Hour (00-23).
\%I Hour (01-12).
\(\% \mathrm{k} \quad\) Hour (0-23).
\(\% 1 \quad\) Hour (1-12).
\(\%\) M Minute (00-59).
\(\%\) Locale's AM or PM.
\(\%\) Time, 12-hour (hh:mm:ss [AP]M).
\%R Time, 24-hour (hh:mm).
\(\%\) Time in seconds since 00:00:00, Jan 1, 1970 (a nonstandard extension).
\(\%\) S Second (00-61).
\%T Time, 24-hour (hh:mm:ss).
\(\% \mathrm{X} \quad\) Locale's time representation ( \(\% \mathrm{H}: \% \mathrm{M}: \% \mathrm{~S}\) ).
\(\%\) Offset from UTC ( \(\pm \mathrm{hhmm}\) ), or nothing if no time zone is determinable.
\(\%\) Time zone (EDT), or nothing if no time zone is determinable.
Date fields:
\(\%\) Locale's abbreviated weekday name (Sun-Sat).
\(\%\) Locale's full weekday name, variable length (Sunday-Saturday).
\(\%\) Locale's abbreviated month name (Jan-Dec).
\%B Locale's full month name, variable length (January-December).
\%c Locale's date and time (Sat Nov 04 12:02:33 EST 1989).
\(\%\) Century (00-99).
\(\%\) d Day of month (01-31).
\(\%\) Day of month (1-31).
\%D Date (mm/dd/yy).
\%h Same as \%b.
\%j Day of year (001-366).
\(\% m \quad\) Month (01-12).
\(\% \quad\) Ueek number of year with Sunday as first day of week (00-53).
\%w Day of week (0-6).
\%W Week number of year with Monday as first day of week (00-53).
\(\%\) Locale's date representation (mm/dd/yy).
\(\%\) Last two digits of year (00-99).
\(\% Y \quad\) Year (1970-).
See also: [strptime], page 945, [localtime], page 942, [gmtime], page 942, [mktime], page 943, [time], page 941, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.
[tm_struct, nchars] = strptime (str, fmt)
Convert the string str to the time structure tm_struct under the control of the format string fint.
If fmt fails to match, nchars is 0 ; otherwise, it is set to the position of last matched character plus 1. Always check for this unless you're absolutely sure the date string will be parsed correctly.
See also: [strftime], page 943, [localtime], page 942, [gmtime], page 942, [mktime], page 943, [time], page 941, [now], page 941, [date], page 946, [clock], page 945, [datenum], page 948, [datestr], page 949, [datevec], page 951, [calendar], page 951, [weekday], page 952.

Most of the remaining functions described in this section are not patterned after the standard C library. Some are available for compatibility with matlab and others are provided because they are useful.
clock ()
Return the current local date and time as a date vector.
The date vector contains the following fields: current year, month (1-12), day (1-31), hour ( \(0-23\) ), minute ( \(0-59\) ), and second ( \(0-61\) ). The seconds field has a fractional part after the decimal point for extended accuracy.

For example:
```

fix (clock ())
\#}\begin{array}{lllllll}{1993}\&{8}\&{20}\&{4}\&{56}\&{1}

```
clock is more accurate on systems that have the gettimeofday function.
See also: [now], page 941, [date], page 946, [datevec], page 951.
date ()
Return the current date as a character string in the form DD-MMM-YYYY.
For example:
```

date ()
=> 20-Aug-1993

```

See also: [now], page 941, [clock], page 945, [datestr], page 949, [localtime], page 942.
etime ( \(t 2\), \(t 1\) )
Return the difference in seconds between two time values returned from clock ( \(t 2\) t1).
For example:
```

t0 = clock ();

# many computations later...

elapsed_time = etime (clock (), t0);

```
will set the variable elapsed_time to the number of seconds since the variable to was set.

See also: [tic], page 947, [toc], page 947, [clock], page 945, [cputime], page 946, [addtodate], page 951.
[total, user, system] = cputime ();
Return the CPU time used by your Octave session.
The first output is the total time spent executing your process and is equal to the sum of second and third outputs, which are the number of CPU seconds spent executing in user mode and the number of CPU seconds spent executing in system mode, respectively.
If your system does not have a way to report CPU time usage, cputime returns 0 for each of its output values.
Note that because Octave used some CPU time to start, it is reasonable to check to see if cputime works by checking to see if the total CPU time used is nonzero.

See also: [tic], page 947, [toc], page 947.
is_leap_year ()
is_leap_year (year)
Return true if year is a leap year and false otherwise.
If no year is specified, is_leap_year uses the current year.
For example:
```

is_leap_year (2000)
\# 1

```

See also: [weekday], page 952, [eomday], page 952, [calendar], page 951.
tic ()
id = tic ()
Initialize a wall-clock timer.
Calling tic without an output argument resets the internal timer. Subsequent calls to toc return the number of seconds since the timer was set.
If called with one output argument, tic creates a new timer instance and returns a timer identifier id. The id is a scalar of type uint64 that may be passed to toc to check elapsed time on this timer, rather than the default internal timer.
Example 1 : benchmarking code with internal timer
```

tic;

# many computations later...

elapsed_time = toc;

```

Example 2 : mixed timer id and internal timer
```

tic;
pause (1);
toc

# Elapsed time is 1.0089 seconds.

id = tic;
pause (2);
toc (id)

# Elapsed time is 2.01142 seconds.

toc
Elapsed time is 3.02308 seconds.

```

Calling tic and toc in this way allows nested timing calls.
If you are more interested in the CPU time that your process used, you should use the cputime function instead. The tic and toc functions report the actual wall clock time that elapsed between the calls. This may include time spent processing other jobs or doing nothing at all.
See also: [toc], page 947, [cputime], page 946.
toc ()
toc (id)
elapsed_time \(=\) toc (...)
Measure elapsed time on a wall-clock timer.
With no arguments, return the number of seconds elapsed on the internal timer since the last call to tic.
When given the identifier id of a specific timer, return the number of seconds elapsed since the timer id was initialized.
See [tic], page 947, for examples of the use of tic/toc.
See also: [tic], page 947, [cputime], page 946.
pause ()
pause (n)
old_state = pause ("on")
```

old_state = pause ("off")

```
old_state = pause ("query")

Suspend the execution of the program or change the state of the pause function.
If invoked without an input arguments then the program is suspended until a character is typed. If argument \(n\) is a positive real value, it indicates the number of seconds the program shall be suspended, for example:
```

tic; pause (0.05); toc
\dashv Elapsed time is 0.05039 seconds.

```

The following example prints a message and then waits 5 seconds before clearing the screen.
```

disp ("wait please...");
pause (5);
clc;

```

If invoked with a string argument "on", "off", or "query", the state of the pause function is changed or queried. When the state is "off", the pause function returns immediately. The optional return value contains the previous state of the pause function. In the following example pause is disabled locally:
```

old_state = pause ("off");
tic; pause (0.05); toc
Elapsed time is 3.00407e-05 seconds.
pause (old_state);

```

While the program is suspended Octave still handles figures painting and graphics callbacks execution.

See also: [kbhit], page 278 .
```

days = datenum (datevec)
days = datenum (year, month, day)
days = datenum (year, month, day, hour)
days = datenum (year, month, day, hour, minute)
days = datenum (year, month, day, hour, minute, second)
days = datenum ("datestr")
days = datenum ("datestr", f)
days = datenum ("datestr", p)
[days, secs] = datenum (...)

```

Return the date/time input as a serial day number, with Jan 1, 0000 defined as day 1.

The integer part, floor (days) counts the number of complete days in the date input.
The fractional part, rem (days, 1) corresponds to the time on the given day.
The input may be a date vector (see [datevec], page 951), date string (see [datestr], page 949 ), or directly specified as input.
When processing input datestrings, \(f\) is the format string used to interpret date strings (see [datestr], page 949). If no format \(f\) is specified, then a relatively slow search is performed through various formats. It is always preferable to specify the format string \(f\) if it is known. Formats which do not specify a particular time component
will have the value set to zero. Formats which do not specify a date will default to January 1st of the current year.
\(p\) is the year at the start of the century to which two-digit years will be referenced. If not specified, it defaults to the current year minus 50 .
The optional output secs holds the time on the specified day with greater precision than days.
Notes:
- Years can be negative and/or fractional.
- Months below 1 are considered to be January.
- Days of the month start at 1 .
- Days beyond the end of the month go into subsequent months.
- Days before the beginning of the month go to the previous month.
- Days can be fractional.

Caution: datenums represent a specific time for the Earth as a whole. They do not take in to account time zones (shifts in time based on location), nor seasonal changes due to Daylight Savings Time (shifts in time based on local regulation). Be aware that it is possible to create datenums that, when interpreted by a function which accounts for time zone and DST shifts such as datestr, are nonexistent or ambiguous.
Caution: this function does not attempt to handle Julian calendars so dates before October 15, 1582 are wrong by as much as eleven days. Also, be aware that only Roman Catholic countries adopted the calendar in 1582. It took until 1924 for it to be adopted everywhere. See the Wikipedia entry on the Gregorian calendar for more details.
Warning: leap seconds are ignored. A table of leap seconds is available on the Wikipedia entry for leap seconds.
See also: [datestr], page 949, [datevec], page 951, [now], page 941, [clock], page 945, [date], page 946.
str = datestr (date)
str = datestr (date, f)
str = datestr (date, f, p)
Format the given date/time according to the format \(f\) and return the result in str.
date is a serial date number (see [datenum], page 948), a date vector (see [datevec], page 951), or a string or cell array of strings. In the latter case, it is passed to datevec to guess the input date format.
\(f\) can be an integer which corresponds to one of the codes in the table below, or a date format string.
\(p\) is the year at the start of the century in which two-digit years are to be interpreted in. If not specified, it defaults to the current year minus 50 .
For example, the date 730736.65149 (2000-09-07 15:38:09.0934) would be formatted as follows:
\begin{tabular}{lll} 
Code & Format & Example \\
0 & dd-mmm-yyyy HH:MM:SS & \(07-\) Sep-2000 15:38:09 \\
1 & dd-mmm-yyyy & \(07-\) Sep-2000 \\
2 & \(\mathrm{~mm} / \mathrm{dd} /\) yy & \(09 / 07 / 00\) \\
3 & mmm & Sep \\
4 & m & S \\
5 & mm & 09 \\
6 & \(\mathrm{~mm} / \mathrm{dd}\) & \(09 / 07\) \\
7 & dd & 07 \\
8 & ddd & Thu \\
9 & d & T \\
10 & yyyy & 2000 \\
11 & yy & 00 \\
12 & mmmyy & Sep00 \\
13 & HH:MM:SS & \(15: 38: 09\) \\
14 & HH:MM:SS PM & \(3: 38: 09 \mathrm{PM}\) \\
15 & HH:MM & \(15: 38\) \\
16 & HH:MM PM & \(3: 38\) PM \\
17 & QQ-YY & Q3-00 \\
18 & QQ & Q3 \\
19 & dd/mm & \(07 / 09\) \\
20 & dd/mm/yy & \(07 / 09 / 00\) \\
21 & mmm.dd,yyyy HH:MM:SS & Sep.07,2000 15:38:08 \\
22 & mmm.dd,yyyy & Sep.07,2000 \\
23 & mm/dd/yyyy & \(09 / 07 / 2000\) \\
24 & dd/mm/yyyy & \(07 / 09 / 2000\) \\
25 & yy/mm/dd & \(00 / 09 / 07\) \\
26 & yyyy/mm/dd & \(2000 / 09 / 07\) \\
27 & QQ-YYYY & Q3-2000 \\
28 & mmmyyyy & Sep2000 \\
29 & yyyy-mm-dd & \(2000-09-07\) \\
30 & yyyymmddTHHMMSS & \(20000907 \mathrm{T153808}\) \\
31 & yyyy-mm-dd HH:MM:SS & \(2000-09-0715: 38: 08\) \\
1 & &
\end{tabular}

If \(f\) is a format string, the following symbols are recognized:
\begin{tabular}{lll} 
Symbol & Meaning & Example \\
yyyy & Full year & 2005 \\
yy & Two-digit year & 05 \\
mmmm & Full month name & December \\
mmm & Abbreviated month name & Dec \\
mm & Numeric month number (padded with zeros) & \(01,08,12\) \\
m & First letter of month name (capitalized) & D \\
dddd & Full weekday name & Sunday \\
ddd & Abbreviated weekday name & Sun \\
dd & Numeric day of month (padded with zeros) & 11 \\
d & First letter of weekday name (capitalized) & S
\end{tabular}
\begin{tabular}{lll} 
HH & \begin{tabular}{l} 
Hour of day, padded with zeros, \\
or padded with spaces if PM is set
\end{tabular} & \(09: 00\) \\
MM & Minute of hour (padded with zeros) & \(9: 00 \mathrm{AM}\) \\
SS & Second of minute (padded with zeros) & \(10: 05\) \\
FFF & Milliseconds of second (padded with zeros) & \(10: 05: 03\) \\
AM & Use 12-hour time format & 10:05:03.012 \\
PM & Use 12-hour time format & 11:30 AM \\
11:30 PM
\end{tabular}

If \(f\) is not specified or is -1 , then use 0,1 or 16 , depending on whether the date portion or the time portion of date is empty.
If \(p\) is not specified, it defaults to the current year minus 50 .
If a matrix or cell array of dates is given, a column vector of date strings is returned.
See also: [datenum], page 948, [datevec], page 951, [date], page 946, [now], page 941, [clock], page 945.
v = datevec (date)
\(\mathrm{v}=\) datevec (date, f)
\(v=\) datevec (date, \(p\) )
\(v=\) datevec (date, \(f, p\) )
[y, m, d, h, mi, s] = datevec (...)
Convert a serial date number (see [datenum], page 948) or date string (see [datestr], page 949) into a date vector.
A date vector is a row vector with six members, representing the year, month, day, hour, minute, and seconds respectively.
\(f\) is the format string used to interpret date strings (see [datestr], page 949). If date is a string, but no format is specified, then a relatively slow search is performed through various formats. It is always preferable to specify the format string \(f\) if it is known. Formats which do not specify a particular time component will have the value set to zero. Formats which do not specify a date will default to January 1st of the current year.
\(p\) is the year at the start of the century to which two-digit years will be referenced. If not specified, it defaults to the current year minus 50 .
See also: [datenum], page 948, [datestr], page 949, [clock], page 945, [now], page 941, [date], page 946.
\(d=\) addtodate (d, \(q, f\) )
Add \(q\) amount of time (with units \(f\) ) to the serial datenum, \(d\).
\(f\) must be one of "year", "month", "day", "hour", "minute", "second", or "millisecond".

See also: [datenum], page 948, [datevec], page 951, [etime], page 946.
\(c=\) calendar ()
\(c=\) calendar (d)
\(c=c a l e n d a r(y, m)\)
calendar (...)
Return the current monthly calendar in a \(6 \times 7\) matrix.

If \(d\) is specified, return the calendar for the month containing the date \(d\), which must be a serial date number or a date string.
If \(y\) and \(m\) are specified, return the calendar for year \(y\) and month \(m\).
If no output arguments are specified, print the calendar on the screen instead of returning a matrix.

See also: [datenum], page 948, [datestr], page 949.
\([n, s]=\) weekday (d)
\([n, s]=\) weekday (d, format)
Return the day of the week as a number in \(n\) and as a string in \(s\).
The days of the week are numbered \(1-7\) with the first day being Sunday.
\(d\) is a serial date number or a date string.
If the string format is not present or is equal to "short" then \(s\) will contain the abbreviated name of the weekday. If format is "long" then \(s\) will contain the full name.
Table of return values based on format:
\begin{tabular}{lll}
\(n\) & "short" & "long" \\
1 & Sun & Sunday \\
2 & Mon & Monday \\
3 & Tue & Tuesday \\
4 & Wed & Wednesday \\
5 & Thu & Thursday \\
6 & Fri & Friday \\
7 & Sat & Saturday
\end{tabular}

See also: [eomday], page 952, [is_leap_year], page 946, [calendar], page 951, [datenum], page 948, [datevec], page 951.
\(e=\) eomday ( \(y, m\) )
Return the last day of the month \(m\) for the year \(y\).
See also: [weekday], page 952, [datenum], page 948, [datevec], page 951, [is_leap_year], page 946, [calendar], page 951.
datetick ()
datetick (axis_str)
datetick (date_format)
datetick (axis_str, date_format)
datetick (..., "keeplimits")
datetick (..., "keepticks")
datetick (hax, ...)
Add date-formatted tick labels to an axis.
The axis to apply the ticks to is determined by axis_str which can take the values "x", "y", or "z". The default value is "x".
The formatting of the labels is determined by the variable date_format, which can either be a string or positive integer that datestr accepts.

If the first argument hax is an axes handle, then plot into this axes, rather than the current axes returned by gca.
See also: [datenum], page 948, [datestr], page 949.

\subsection*{36.2 Filesystem Utilities}

Octave includes many utility functions for copying, moving, renaming, and deleting files; for creating, reading, and deleting directories; for retrieving status information on files; and for manipulating file and path names.
```

movefile f1
movefile f1 f2
movefile f1 f2 f
movefile (f1)
movefile (f1, f2)
movefile (f1, f2, 'f')
[status, msg, msgid] = movefile (...)

```

Move the source file or directory \(f 1\) to the destination \(f 2\).
The name f1 may contain globbing patterns, or may be a cell array of strings. If f1 expands to multiple filenames, \(f 2\) must be a directory.
If no destination \(f 2\) is specified then the destination is the present working directory. If \(f 2\) is a filename then \(f 1\) is renamed to \(f 2\).
When the force flag ' \(f\) ' is given any existing files will be overwritten without prompting.
If successful, status is logical 1, and msg, msgid are empty character strings (""). Otherwise, status is logical 0, msg contains a system-dependent error message, and msgid contains a unique message identifier. Note that the status code is exactly opposite that of the system command.
See also: [rename], page 953, [copyfile], page 953, [unlink], page 954, [delete], page 405, [glob], page 959.
rename old new
[status, msg] = rename (old, new)
Change the name of file old to new.
If successful, status is 0 and msg is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.
See also: [movefile], page 953, [copyfile], page 953, [ls], page 982, [dir], page 983.
copyfile f1 f2
copyfile \(f 1 f 2 f\)
copyfile (f1, f2)
copyfile (f1, f2, 'f')
[status, msg, msgid] = copyfile (...)
Copy the source file(s) or directory \(f 1\) to the destination \(f 2\).
The name f1 may contain globbing patterns, or may be a cell array of strings. If f1 expands to multiple filenames, \(f 2\) must be a directory.

When the force flag ' f ' is given any existing files will be overwritten without prompting.
If successful, status is logical 1 , and msg, msgid are empty character strings (""). Otherwise, status is logical 0, msg contains a system-dependent error message, and msgid contains a unique message identifier. Note that the status code is exactly opposite that of the system command.
See also: [movefile], page 953, [rename], page 953, [unlink], page 954, [delete], page 405, [glob], page 959.
unlink (file)
[status, msg] = unlink (file)
Delete the file named file.
If successful, status is 0 and msg is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.
See also: [delete], page 405, [rmdir], page 955.
link old new
[status, msg] = link (old, new)
Create a new link (also known as a hard link) to an existing file.
If successful, status is 0 and msg is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.

See also: [symlink], page 954, [unlink], page 954, [readlink], page 954, [lstat], page 956.
symlink old new
[status, msg] = symlink (old, new)
Create a symbolic link new which contains the string old.
If successful, status is 0 and msg is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.
See also: [link], page 954, [unlink], page 954, [readlink], page 954, [lstat], page 956.
readlink symlink
[result, err, msg] = readlink (symlink)
Read the value of the symbolic link symlink.
If successful, result contains the contents of the symbolic link symlink, err is 0 , and msg is an empty string. Otherwise, err is nonzero and msg contains a systemdependent error message.
See also: [lstat], page 956, [symlink], page 954, [link], page 954, [unlink], page 954, [delete], page 405.
mkdir dirname
mkdir parent dirname
mkdir (dirname)
mkdir (parent, dirname)
[status, msg, msgid] = mkdir (...)
Create a directory named dirname in the directory parent, creating any intermediate directories if necessary.

If dirname is a relative path, and no parent directory is specified, then the present working directory is used.
If successful, status is logical 1 , and msg, msgid are empty character strings (""). Otherwise, status is logical \(0, \mathrm{msg}\) contains a system-dependent error message, and msgid contains a unique message identifier. Note that the status code is exactly opposite that of the system command.
When creating a directory permissions will be set to 0777 - UMASK.
See also: [rmdir], page 955, [pwd], page 983, [cd], page 982, [umask], page 955.
rmdir dir
rmdir (dir, "s")
[status, msg, msgid] = rmdir (...)
Remove the directory named dir.
If the optional second parameter is supplied with value "s", recursively remove all subdirectories as well.

If successful, status is logical 1 , and msg, msgid are empty character strings (""). Otherwise, status is logical 0 , msg contains a system-dependent error message, and msgid contains a unique message identifier.
See also: [mkdir], page 954, [confirm_recursive_rmdir], page 955, [pwd], page 983.
val = confirm_recursive_rmdir ()
old_val = confirm_recursive_rmdir (new_val)
confirm_recursive_rmdir (new_val, "local")
Query or set the internal variable that controls whether Octave will ask for confirmation before recursively removing a directory tree.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [rmdir], page 955.
mkfifo (name, mode)
[status, msg] = mkfifo (name, mode)
Create a FIFO special file named name with file mode mode.
mode is interpreted as an octal number and is subject to umask processing. The final calculated mode is mode - umask.
If successful, status is 0 and msg is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.
See also: [pipe], page 976, [umask], page 955.
umask (mask)
Set the permission mask for file creation.
The parameter mask is an integer, interpreted as an octal number.
If successful, returns the previous value of the mask (as an integer to be interpreted as an octal number); otherwise an error message is printed.

The permission mask is a UNIX concept used when creating new objects on a file system such as files, directories, or named FIFOs. The object to be created has base permissions in an octal number mode which are modified according to the octal value of mask. The final permissions for the new object are mode - mask.

See also: [fopen], page 295, [mkdir], page 954, [mkfifo], page 955.
```

[info, err, msg] = stat (file)
[info, err, msg] = stat (fid)
[info, err, msg] = lstat (file)
[info, err, msg] = lstat (fid)

```

Return a structure info containing the following information about file or file identifier fid.
dev ID of device containing a directory entry for this file.
ino File number of the file.
mode File mode, as an integer. Use the functions S_ISREG, S_ISDIR, S_ISCHR, S_ISBLK, S_ISFIFO, S_ISLNK, or S_ISSOCK to extract information from this value.
modestr File mode, as a string of ten letters or dashes as would be returned by ls -1 .
nlink Number of links.
uid User ID of file's owner.
gid Group ID of file's group.
rdev ID of device for block or character special files.
size Size in bytes.
atime Time of last access in the same form as time values returned from time. See Section 36.1 [Timing Utilities], page 941.
mtime Time of last modification in the same form as time values returned from time. See Section 36.1 [Timing Utilities], page 941.
ctime Time of last file status change in the same form as time values returned from time. See Section 36.1 [Timing Utilities], page 941.
blksize Size of blocks in the file.
blocks Number of blocks allocated for file.
If the call is successful err is 0 and msg is an empty string. If the file does not exist, or some other error occurs, info is an empty matrix, err is -1 , and msg contains the corresponding system error message.
If file is a symbolic link, stat will return information about the actual file that is referenced by the link. Use lstat if you want information about the symbolic link itself.
For example:
```

[info, err, msg] = stat ("/vmlinuz")

```
```


# info =

    {
        atime = 855399756
        rdev = 0
        ctime = 847219094
        uid = 0
        size = 389218
        blksize = 4096
        mtime = 847219094
        gid = 6
        nlink = 1
        blocks = 768
        mode = -rw-r--r--
        modestr = -rw-r--r--
        ino = 9316
        dev = 2049
    }
    merr = 0
msg =

```

See also: [lstat], page 956, [ls], page 982, [dir], page 983, [isfile], page 959, [isfolder], page 959.

S_ISBLK (mode)
Return true if mode corresponds to a block device.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
S_ISCHR (mode)
Return true if mode corresponds to a character device.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
S_ISDIR (mode)
Return true if mode corresponds to a directory.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
S_ISFIFO (mode)
Return true if mode corresponds to a fifo.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
S_ISLNK (mode)
Return true if mode corresponds to a symbolic link.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.

S_ISREG (mode)
Return true if mode corresponds to a regular file.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
S_ISSOCK (mode)
Return true if mode corresponds to a socket.
The value of mode is assumed to be returned from a call to stat.
See also: [stat], page 956, [lstat], page 956.
fileattrib ()
fileattrib (file)
[status, attrib] = fileattrib (...)
[status, msg, msgid] = fileattrib (...)
Report attribute information about file.
If no file or directory is specified, report information about the present working directory.
If successful, the output is a structure with the following fields:
Name Full name of file.
archive True if file is an archive (Windows).
system True if file is a system file (Windows).
hidden True if file is a hidden file (Windows).
directory
True if file is a directory.

\section*{UserRead}

GroupRead
OtherRead
True if the user (group; other users) has read permission for file.
UserWrite
GroupWrite
OtherWrite
True if the user (group; other users) has write permission for file.

\section*{UserExecute}

GroupExecute
OtherExecute
True if the user (group; other users) has execute permission for file.
If an attribute does not apply (e.g., archive on a Unix system) then the field is set to NaN .

If file contains globbing characters, information about all matching files is returned in a structure array.
If outputs are requested, the first is status which takes the value 1 when the operation was successful, and 0 otherwise. The second output contains the structure
described above (attrib) if the operation was successful; otherwise, the second output is a system-dependent error message ( msg ). The third output is an empty string (" ") when the operation was successful, or a unique message identifier (msgid) in the case of failure.

See also: [stat], page 956, [glob], page 959.
\(t f=\) isfile (f)
Return true if \(f\) is a regular file and false otherwise.
If \(f\) is a cell array of strings, \(t f\) is a logical array of the same size.
See also: [isfolder], page 959, [exist], page 144, [stat], page 956, [is_absolute_filename], page 962, [is_rooted_relative_filename], page 962.
isdir (f)
This function is not recommended. Use isfolder or file_in_loadpath instead.
Return true if \(f\) is a directory and false otherwise.
Compatibility Note: The matlab function of the same name will also search for \(f\) in the load path directories. To emulate this behavior use
```

tf = ! isempty (file_in_loadpath (f))

```

See also: [isfolder], page 959, [file_in_loadpath], page 215, [exist], page 144, [stat], page 956, [is_absolute_filename], page 962, [is_rooted_relative_filename], page 962.
\(t f=\) isfolder (f)
Return true if \(f\) is a directory and false otherwise.
If \(f\) is a cell array of strings, \(t f\) is a logical array of the same size.
See also: [isfile], page 959, [exist], page 144, [stat], page 956, [is_absolute_filename], page 962, [is_rooted_relative_filename], page 962.
files \(=\) readdir (dir)
[files, err, msg] = readdir (dir)
Return the names of files in the directory dir as a cell array of strings.
If an error occurs, return an empty cell array in files. If successful, err is 0 and msg is an empty string. Otherwise, err is nonzero and msg contains a system-dependent error message.
See also: [1s], page 982, [dir], page 983, [glob], page 959, [what], page 148.
glob (pattern)
Given an array of pattern strings (as a char array or a cell array) in pattern, return a cell array of filenames that match any of them, or an empty cell array if no patterns match.
The pattern strings are interpreted as filename globbing patterns (as they are used by Unix shells).
Within a pattern
* matches any string, including the null string,
? matches any single character, and
[...] matches any of the enclosed characters.
Tilde expansion is performed on each of the patterns before looking for matching filenames. For example:
```

ls
\#
file1 file2 file3 myfile1 myfile1b
glob ("*file1")
\#
{
[1,1] = file1
[2,1] = myfile1
}
glob ("myfile?")
\#
{
[1,1] = myfile1
}
glob ("file[12]")
\#
{
[1,1] = file1
[2,1] = file2
}

```

Note: On Windows, patterns that contain non-ASCII characters are not supported.
See also: [ls], page 982, [dir], page 983, [readdir], page 959, [what], page 148.
file_in_path (path, file)
file_in_path (path, file, "all")
Return the absolute name of file if it can be found in path.
The value of path should be a colon-separated list of directories in the format described for path. If no file is found, return an empty character string. For example:
```

file_in_path (EXEC_PATH, "sh")
=> "/bin/sh"

```

If the second argument is a cell array of strings, search each directory of the path for element of the cell array and return the first that matches.
If the third optional argument "all" is supplied, return a cell array containing the list of all files that have the same name in the path. If no files are found, return an empty cell array.
See also: [file_in_loadpath], page 215, [dir_in_loadpath], page 216, [path], page 214.
filesep ()
filesep ("all")
Return the system-dependent character used to separate directory names.
If "all" is given, the function returns all valid file separators in the form of a string. The list of file separators is system-dependent. It is ' \(/\) ' (forward slash) under UNIX or Mac OS X, '/' and ' \(\\) ' (forward and backward slashes) under Windows.

See also: [pathsep], page 215.
[dir, name, ext] = fileparts (filename)
Return the directory, name, and extension components of filename.
The input filename is a string which is parsed. There is no attempt to check whether the filename or directory specified actually exists.

See also: [fullfile], page 961, [filesep], page 960.
filename \(=\) fullfile (dir1, dir2, ..., file)
Build complete filename from separate parts.
The function joins any number of path components intelligently. The return value is the concatenation of each component with exactly one file separator between each part of the path and at most one leading and/or trailing file separator.
The input arguments might be strings or cell strings. Any input arguments that are cell strings must contain one single string or must be equal in size. In that case, the function returns a cell string of filepaths of the same dimensions as the input cell strings, e.g.:
```

fullfile ("/home/username", "data", {"f1.csv", "f2.csv", "f3.csv"})
\#
{
[1,1] = /home/username/data/f1.csv
[1,2] = /home/username/data/f2.csv
[1,3] = /home/username/data/f3.csv
}

```

On Windows systems, while forward slash file separators do work, they are replaced by backslashes. In addition, drive letters are stripped of leading file separators to obtain a valid file path.
Note: fullfile does not perform any validation of the resulting full filename.
See also: [fileparts], page 961, [filesep], page 960.
tilde_expand (string)
tilde_expand (cellstr)
Perform tilde expansion on string.
If string begins with a tilde character, \(\left({ }^{(\sim)}\right)\), all of the characters preceding the first slash (or all characters, if there is no slash) are treated as a possible user name, and the tilde and the following characters up to the slash are replaced by the home directory of the named user. If the tilde is followed immediately by a slash, the tilde is replaced by the home directory of the user running Octave.
If the input is a cell array of strings cellstr then tilde expansion is performed on each string element.
Examples:
```

tilde_expand ("~joeuser/bin")
=> "/home/joeuser/bin"
tilde_expand ("~/bin")
=> "/home/jwe/bin"

```
[cname, status, msg] = canonicalize_file_name (fname)
Return the canonical name of file fname.
If the file does not exist the empty string ("") is returned. No tilde expansion of fname is performed.

See also: [make_absolute_filename], page 962, [is_absolute_filename], page 962, [is_rooted_relative_filename], page 962, [is_same_file], page 962, [tilde_expand], page 961.
make_absolute_filename (file)
Return the full name of file beginning from the root of the file system.
No check is done for the existence of file. No tilde expansion of file is performed.
See also: [canonicalize_file_name], page 962, [is_absolute_filename], page 962, [is_rooted_relative_filename], page 962, [isfolder], page 959, [tilde_expand], page 961.
is_absolute_filename (file)
Return true if file is an absolute filename.
See also: [is_rooted_relative_filename], page 962, [make_absolute_filename], page 962, [isfolder], page 959.
same = is_same_file (filepath1, filepath2)
Return true if filepath1 and filepath2 refer to the same file.
If either filepath1 or filepath2 is a cell array of strings, then an array of the same size is returned, containing the values described above for every member of the cell array. The other argument may also be a cell array of strings (of the same size) or a string.
Programming Notes: Depending on the operating system and file system, the same file or folder can be referred to with different paths. In particular, paths on the Windows platform may differ in case (file1 vs. FILE1), file separator ('\' vs. '/'), and format (A~spaces.txt ( 8.3 convention) vs. A filename with spaces.txt). This function returns true if the paths in filepath1 and filepath2 actually refer to the same file or folder, and false otherwise.
Note that unlike strcmp, this function requires that filepath1 and filepath2 exist, as well as point to the same location, in order to return true.

See also: [canonicalize_file_name], page 962, [strcmp], page 84.
is_rooted_relative_filename (file)
Return true if file is a rooted-relative filename.
See also: [is_absolute_filename], page 962, [make_absolute_filename], page 962, [isfolder], page 959.
val = recycle ()
old_val = recycle (new_val)
Query or set the preference for recycling deleted files.
When recycling is enabled, commands which would permanently erase files instead move them to a temporary location (such as the directory labeled Trash).

Programming Note: This function is provided for matlab compatibility, but recycling is not implemented in Octave. To help avoid accidental data loss an error will be raised if an attempt is made to enable file recycling.
See also: [delete], page 405, [rmdir], page 955.

\subsection*{36.3 File Archiving Utilities}
bunzip2 (bzfile)
bunzip2 (bzfile, dir)
filelist = bunzip2 (...)
Unpack the bzip2 archive bzfile.
If dir is specified the files are unpacked in this directory rather than the one where bzfile is located.
The optional output filelist is a list of the uncompressed files.
See also: [bzip2], page 965, [unpack], page 965, [gunzip], page 963, [unzip], page 964, [untar], page 964.
filelist = gzip (files)
filelist = gzip (files, dir)
Compress the list of files and directories specified in files.
files is a character array or cell array of strings. Shell wildcards in the filename such as ' \(*\) ' or '?' are accepted and expanded. Each file is compressed separately and a new file with a ".gz" extension is created. The original files are not modified, but existing compressed files will be silently overwritten. If a directory is specified then gzip recursively compresses all files in the directory.
If dir is defined the compressed files are placed in this directory, rather than the original directory where the uncompressed file resides. Note that this does not replicate a directory tree in dir which may lead to files overwriting each other if there are multiple files with the same name.
If dir does not exist it is created.
The optional output filelist is a list of the compressed files.
See also: [gunzip], page 963, [unpack], page 965, [bzip2], page 965, [zip], page 964, [tar], page 964.
gunzip (gzfile)
gunzip (gzfile, dir)
filelist = gunzip (...)
Unpack the gzip archive gzfile.
If gzfile is a directory, all gzfiles in the directory will be recursively unpacked.
If dir is specified the files are unpacked in this directory rather than the one where gzfile is located.
The optional output filelist is a list of the uncompressed files.
See also: [gzip], page 963, [unpack], page 965, [bunzip2], page 963, [unzip], page 964, [untar], page 964.
```

filelist = tar (tarfile, files)
filelist = tar (tarfile, files, rootdir)

```

Pack the list of files and directories specified in files into the TAR archive tarfile.
files is a character array or cell array of strings. Shell wildcards in the filename such as '*' or '?' are accepted and expanded. Directories are recursively traversed and all files are added to the archive.

If rootdir is defined then any files without absolute pathnames are located relative to rootdir rather than the current directory.
The optional output filelist is a list of the files that were included in the archive.
See also: [untar], page 964, [unpack], page 965, [bzip2], page 965, [gzip], page 963, [zip], page 964 .
untar (tarfile)
untar (tarfile, dir)
filelist \(=\) untar (...)
Unpack the TAR archive tarfile.
If dir is specified the files are unpacked in this directory rather than the current directory.
The optional output filelist is a list of the uncompressed files.
See also: [tar], page 964, [unpack], page 965, [bunzip2], page 963, [gunzip], page 963, [unzip], page 964.
filelist \(=\) zip (zipfile, files)
filelist \(=\) zip (zipfile, files, rootdir)
Compress the list of files and directories specified in files into the ZIP archive zipfile. files is a character array or cell array of strings. Shell wildcards in the filename such as ' \(*\) ' or '?' are accepted and expanded. Directories are recursively traversed and all files are compressed and added to the archive.

If rootdir is defined then any files without absolute pathnames are located relative to rootdir rather than the current directory.

The optional output filelist is a list of the files that were included in the archive.
See also: [unzip], page 964, [unpack], page 965, [bzip2], page 965, [gzip], page 963, [tar], page 964.
unzip (zipfile)
unzip (zipfile, dir)
filelist = unzip (...)
Unpack the ZIP archive zipfile.
If dir is specified the files are unpacked in this directory rather than the current directory.
The optional output filelist is a list of the uncompressed files.
See also: [zip], page 964, [unpack], page 965, [bunzip2], page 963, [gunzip], page 963, [untar], page 964.
```

files = unpack (file)
files = unpack (file, dir)
files = unpack (file, dir, filetype) Valid filetype extensions:
bz
bz2 bzip archive
gz gzip archive
tar tar archive
tarbz
tarbz2
tbz
tbz2 tar + bzip archive
targz
tgz tar + gzip archive
z compress archive
zip zip archive

```

Unpack the archive file based on its extension to the directory dir.
If file is a list of strings, then each file is unpacked individually. Shell wildcards in the filename such as '*' or '?' are accepted and expanded.
If dir is not specified or is empty ([]), it defaults to the current directory. If a directory is in the file list, then filetype must also be specified.
The specific archive filetype is inferred from the extension of the file. The filetype may also be specified directly using a string which corresponds to a known extension.

The optional return value is a list of files unpacked.
See also: [bunzip2], page 963, [gunzip], page 963, [unzip], page 964, [untar], page 964, [bzip2], page 965, [gzip], page 963, [zip], page 964, [tar], page 964.
filelist = bzip2 (files)
filelist = bzip2 (files, dir)
Compress the list of files specified in files.
files is a character array or cell array of strings. Shell wildcards in the filename such as ' \(*\) ' or '?' are accepted and expanded. Each file is compressed separately and a new file with a ".bz2" extension is created. The original files are not modified, but existing compressed files will be silently overwritten.
If dir is defined the compressed files are placed in this directory, rather than the original directory where the uncompressed file resides. Note that this does not replicate a directory tree in dir which may lead to files overwriting each other if there are multiple files with the same name.
If dir does not exist it is created.
The optional output filelist is a list of the compressed files.
See also: [bunzip2], page 963, [unpack], page 965, [gzip], page 963, [zip], page 964, [tar], page 964.

\subsection*{36.4 Networking Utilities}
gethostname ()
Return the hostname of the system where Octave is running.

\subsection*{36.4.1 FTP Objects}

Octave supports the FTP protocol through an object-oriented interface. Use the function ftp to create an FTP object which represents the connection. All FTP functions take an FTP object as the first argument.
\(f=\mathrm{ftp}(\) host \()\)
\(f=\) ftp (host, username, password)
Connect to the FTP server host with username and password.
If username and password are not specified, user "anonymous" with no password is used. The returned FTP object \(f\) represents the established FTP connection.

The list of actions for an FTP object are shown below. All functions require an FTP object as the first argument.

\section*{Method Description}
ascii Set transfer type to ascii
binary Set transfer type to binary
cd Change remote working directory
close Close FTP connection
delete Delete remote file
dir List remote directory contents
mget Download remote files
mkdir Create remote directory
mput Upload local files
rename \(\quad\) Rename remote file or directory
rmdir Remove remote directory
close (f)
Close the FTP connection represented by the FTP object \(f\).
\(f\) is an FTP object returned by the ftp function.
mget (f, file)
mget (f, dir)
mget (f, remote_name, target)
Download a remote file file or directory dir to the local directory on the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
The arguments file and dir can include wildcards and any files or directories on the remote server that match will be downloaded.
If a third string argument target is given, then it must indicate the path to the local destination directory. target may be a relative or absolute path.
mput (f, file)
Upload the local file file into the current remote directory on the FTP connection \(f\). \(f\) is an FTP object returned by the ftp function.
The argument file is passed through the glob function and any files that match the wildcards in file will be uploaded.
cd (f)
cd (f, path)
Get or set the remote directory on the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
If path is not specified, return the remote current working directory. Otherwise, set the remote directory to path and return the new remote working directory.
If the directory does not exist, an error message is printed and the working directory is not changed.
```

lst = dir (f)

```

List the current directory in verbose form for the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
ascii (f)
Set the FTP connection \(f\) to use ASCII mode for transfers.
ASCII mode is only appropriate for text files as it will convert the remote host's newline representation to the local host's newline representation.
\(f\) is an FTP object returned by the ftp function.

\section*{binary (f)}

Set the FTP connection \(f\) to use binary mode for transfers.
In binary mode there is no conversion of newlines from the remote representation to the local representation.
\(f\) is an FTP object returned by the ftp function.
delete (f, file)
Delete the remote file file over the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
rename (f, oldname, newname)
Rename or move the remote file or directory oldname to newname, over the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
mkdir (f, path)
Create the remote directory path, over the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.
rmdir (f, path)
Remove the remote directory path, over the FTP connection \(f\).
\(f\) is an FTP object returned by the ftp function.

\subsection*{36.4.2 WWW Access}

Octave can communicate with websites across the Internet. The web function will launch an external web browser to interactively view a site. The remaining functions-urlread, urlwrite, webread, webwrite-are internal Octave functions which can import or export data to/from Octave and a website identified by a URL (Uniform Resource Locator).
```

status = web ()
status = web (url)
status = web (url, option)
status = web (url, option_1,..., option_N)
[status,h, url] = web (...)

```

Open url in the default system web browser.
With no arguments given, the address https://www.octave.org is opened.
Additional options can be passed for MATLAB compatibility, but are ignored.
- '-browser' Open url in the default system browser.
- '-new' No effect on the system browser.
- '-noaddressbox' No effect on the system browser.
- '-notoolbar' No effect on the system browser.

The return value status has one of the values:
- ' 0 ' Found and opened system browser successfully.
- '1' Cannot find the system browser.
- '2' System browser found, but an error occurred.

The return values handle and url are currently unimplemented but given for compatibility.
See also: [weboptions], page 970, [webread], page 969, [webwrite], page 970, [urlread], page 968, [urlwrite], page 969.
\(s=u r l r e a d\) (url)
[s, success] = urlread (url)
[s, success, message] = urlread (url)
[...] = urlread (url, method, param)
Download a remote file specified by its url and return its content in string \(s\).
For example:
```

s = urlread ("ftp://ftp.octave.org/pub/README");

```

The variable success is 1 if the download was successful, otherwise it is 0 in which case message contains an error message.

If no output argument is specified and an error occurs, then the error is signaled through Octave's error handling mechanism.

This function uses libcurl. Curl supports, among others, the HTTP, FTP, and FILE protocols. Username and password may be specified in the URL. For example:
```

s = urlread ("http://user:password@example.com/file.txt");

```

GET and POST requests can be specified by method and param. The parameter method is either 'get' or 'post' and param is a cell array of parameter and value pairs. For example:
```

s = urlread ("http://www.google.com/search", "get",
{"query", "octave"});

```

See also: [urlwrite], page 969.
urlwrite (url, localfile)
\(f=\) urlwrite (url, localfile)
[f, success] = urlwrite (url, localfile)
[f, success, message] = urlwrite (url, localfile)
Download a remote file specified by its url and save it as localfile.
For example:
```

urlwrite ("ftp://ftp.octave.org/pub/README",
"README.txt");

```

The full path of the downloaded file is returned in \(f\).
The variable success is 1 if the download was successful, otherwise it is 0 in which case message contains an error message.
If no output argument is specified and an error occurs, then the error is signaled through Octave's error handling mechanism.
This function uses libcurl. Curl supports, among others, the HTTP, FTP, and FILE protocols. Username and password may be specified in the URL, for example:
```

urlwrite ("http://username:password@example.com/file.txt",
"file.txt");

```

GET and POST requests can be specified by method and param. The parameter method is either 'get' or 'post' and param is a cell array of parameter and value pairs. For example:
```

urlwrite ("http://www.google.com/search", "search.html",
"get", {"query", "octave"});

```

See also: [urlread], page 968.
```

response = webread (url)
response = webread (url, name1, value1,...)
response = webread (..., options)

```

Read content from RESTful web service.
Read content from the web service specified by url and return the content in response.
All key-value pairs given (name1, value1, ...) are appended as query parameters to url. To place a query in the body of the message, use webwrite. The web service defines the acceptable query parameters.
options is a weboptions object that may be used to add other HTTP request options. This argument can be used with either calling form. See help weboptions for a complete list of supported HTTP options.
See also: [weboptions], page 970, [webwrite], page 970.
response \(=\) webwrite (url, name1, value1, ...)
response = webwrite (url, data)
response \(=\) webwrite (..., options)
Write data to RESTful web services.
Write content to the web service specified by url and return the response in response.
All key-value pairs given (name1, value1, . . .) are added as pairs of query parameters to the body of request method (get, post, put, etc.).
options is a weboptions object that may be used to add other HTTP request options. This argument can be used with either calling form. See help weboptions for a complete list of supported HTTP options.

See also: [weboptions], page 970, [webread], page 969.
output = weboptions ()
output = weboptions (name1, value1, ...)
Specify parameters for RESTful web services.
weboptions with no inputs returns a default weboptions object to specify parameters for a request to a web service. A weboptions object can be an optional input argument to the webread and webwrite functions.
Multiple name and value pair arguments may be specified in any order as name1, value1, name2, value2, etc.
The option names must match exactly one of those specified in the table below.
The following options are available:
- 'CharacterEncoding' - Specify the character encoding of the data:
'auto' (default), 'UTF-8', 'US-ASCII' 'auto' chooses an encoding based on the content-type of the data.
- 'UserAgent' - Specify the User Agent for the connection.

Default value is 'GNU Octave/version', where 'version' is the current version of Octave as returned by version.
- 'Timeout' - Specify the timeout value for the connection in seconds.

Default is 10 seconds. 'Inf' is not currently supported.
- 'Username' - User identifier for a basic HTTP connection.

Default is NULL. It must be a string.
- 'Password' - User authentication password for HTTP connection.

Default is NULL. It must be a string or character vector. Programming Note: If you display a weboption object with the Password property set, the value is displayed as a string containing ' \(*\) '. However, the object stores the value of the Password property as plain text.
- 'KeyName' - Specify the name of an additional key to be added to the HTTP request header. It should be coupled with 'KeyValue'. It must be a string or character vector.
- 'KeyValue' - Specify the value of the key 'KeyName'.
'KeyName' must be present in order to assign to this field.
- 'HeaderFields' - Specify the header fields for the connection.

Names and values of header fields, specified as an m-by-2 array of strings or cell array of character vectors to add to the HTTP request header. HeaderFields \(\{\mathrm{i}, 1\}\) is the name of a field and HeaderFields \(\{i, 2\}\) is its value.
```

weboptions ("HeaderFields", {"Content-Length" "78";"Content-Type" "application/
Creates a weboptions object that contains two header fields:
Content-Length with value 78 and Content-Type with value application/json.|

```
- 'ContentType' - Specify the content type of the data.

The following values are available: 'auto', 'text', 'json'
Default is 'auto'. It automatically determines the content type. All other formats like 'audio', 'binary', etc. available in matlab are not currently supported.
- 'ContentReader' - Not yet implemented. Only for matlab compatibility.
- 'MediaType' - Not yet implemented. Only for matlab compatibility.
- 'RequestMethod' - Specifies the type of request to be made.

The following methods are available: 'get', 'put', 'post', 'delete', 'patch' webread uses the HTTP GET method. webwrite uses the HTTP POST method as default.
- 'ArrayFormat' - Not yet implemented. Only for matlab compatibility.
- 'CertificateFilename' - Not yet implemented. Only for matlab compatibility.
See also: [webread], page 969, [webwrite], page 970.

\subsection*{36.4.3 Base64 and Binary Data Transmission}

Some transmission channels can not accept binary data. It is customary to encode binary data in Base64 for transmission and to decode the data upon reception.
\(s=\) base64_encode ( \(x\) )
Encode a double matrix or array \(x\) into the base64 format string \(s\).
See also: [base64_decode], page 971, [matlab.net.base64decode], page 972, [matlab.net.base64encode], page 971.
x = base64_decode (s)
\(x\) = base64_decode (s, dims)
Decode the double matrix or array \(x\) from the base64 encoded string \(s\).
The optional input parameter dims should be a vector containing the dimensions of the decoded array.
See also: [base64_encode], page 971, [matlab.net.base64decode], page 972, [matlab.net.base64encode], page 971.
b64_str = matlab.net.base64encode (in)
Convert in to a base64 encoded string b64_str.
The input in can be a string or numeric vector. The output b64_str will be encoded according to RFC 4648.
See also: [matlab.net.base64decode], page 972, [base64_decode], page 971, [base64_encode], page 971, [unicode2native], page 94.
```

out_vec = matlab.net.base64decode (b64_str)

```

Convert base64 encoded b64_str to uint8 vector out_vec.
The input b64_str must be a string vector. The output out_vec will be a uint8 vector that is decoded according to RFC 4648.

See also: [matlab.net.base64encode], page 971, [base64_decode], page 971, [base64_encode], page 971, [native2unicode], page 95.

\subsection*{36.5 Controlling Subprocesses}

Octave includes some high-level commands like system and popen for starting subprocesses. If you want to run another program to perform some task and then look at its output, you will probably want to use these functions.

Octave also provides several very low-level Unix-like functions which can also be used for starting subprocesses, but you should probably only use them if you can't find any way to do what you need with the higher-level functions.
```

system ("string")
system ("string", return_output)
system ("string", return_output, type)
[status, output] = system (...)

```

Execute a shell command specified by string.
If system is called with one or more output arguments, or if the optional argument return_output is true and the subprocess is started synchronously, then the output from the command is returned as a variable. Otherwise, if the subprocess is executed synchronously, its output is sent to the standard output. To send the output of a command executed with system through the pager, use a command like
```

[~, text] = system ("cmd");
more on;
disp (text);
more on;
printf ("%s\n", nthargout (2, "system", "cmd"));

```
    or

If the optional argument type is "async", the process is started in the background and the process ID of the child process is returned immediately. Otherwise, the child process is started and Octave waits until it exits. If the type argument is omitted, it defaults to the value "sync".
The system function can return two values. The first is the exit status of the command and the second is any output from the command that was written to the standard output stream. For example,
```

[status, output] = system ("echo foo \& exit 2");

```
will set the variable output to the string 'foo', and the variable status to the integer '2'.
For commands run asynchronously, status is the process id of the command shell that is started to run the command.

The shell used for executing commands varies with operating system and is typically /bin/sh for UNIX systems and cmd.exe for Windows systems.
See also: [unix], page 973, [dos], page 973.
unix ("command")
status = unix ("command")
[status, text] = unix ("command")
[...] = unix ("command", "-echo")
Execute a system command if running under a Unix-like operating system, otherwise do nothing.
Octave waits for the external command to finish before returning the exit status of the program in status and any output in text.
When called with no output argument, or the "-echo" argument is given, then text is also sent to standard output.

See also: [dos], page 973, [system], page 972, [isunix], page 986, [ismac], page 986, [ispc], page 986.
dos ("command")
status = dos ("command")
[status, text] = dos ("command")
[...] = dos ("command", "-echo")
Execute a system command if running under a Windows-like operating system, otherwise do nothing.
Octave waits for the external command to finish before returning the exit status of the program in status and any output in text.
When called with no output argument, or the "-echo" argument is given, then text is also sent to standard output.
See also: [unix], page 973, [system], page 972, [isunix], page 986, [ismac], page 986, [ispc], page 986.
open file
output \(=\) open (file)
Open the file file in Octave or in an external application based on the file type as determined by the filename extension.
By default, recognized file types are
.m Open file in the editor. No output value is returned.
.mat
octave-workspace
Open the data file with load. If no return value output is requested, variables are loaded in the base workspace. Otherwise output will be a structure containing loaded data. See [load function], page 281.
.ofig Open the figure with hgload. See [hgload function], page 418.
.fig, .ofig
Load the figure returned.

Custom file extensions may also be handled if a function openxxx, where xxx is the extension, is found in the load path. The function must accept the file name as input. For example, in order to load ". dat" data files in the base workspace, as is done by default for ".mat" files, one may define "opendat. m " with the following contents:
```

function retval = opendat (fname)
evalin ("base", sprintf ("load ('%s');", fname));
endfunction

```

Other file types are opened in the appropriate external application.
```

output = perl (scriptfile)
output = perl (scriptfile, argument1, argument2,...)
[output, status] = perl (...)

```

Invoke Perl script scriptfile, possibly with a list of command line arguments.
Return output in output and optional status in status. If scriptfile is not an absolute filename it is searched for in the current directory and then in the Octave loadpath.

See also: [system], page 972, [python], page 974.
```

output = python (scriptfile)
output = python (scriptfile, argument1, argument2,...)
[output, status] = python (...)

```

Invoke Python script scriptfile, possibly with a list of command line arguments.
Return output in output and optional status in status. If scriptfile is not an absolute filename it is searched for in the current directory and then in the Octave loadpath.

See also: [system], page 972, [perl], page 974 .
```

fid = popen (command, mode)

```

Start a process and create a pipe.
The name of the command to run is given by command. The argument mode may be
"r" The pipe will be connected to the standard output of the process, and open for reading.
"W" The pipe will be connected to the standard input of the process, and open for writing.

The file identifier corresponding to the input or output stream of the process is returned in fid.

For example:
```

fid = popen ("ls -ltr / | tail -3", "r");
while (ischar (s = fgets (fid)))
fputs (stdout, s);
endwhile
\dashv drwxr-xr-x 33 root root 3072 Feb 15 13:28 etc
\dashv drwxr-xr-x 3 root root 1024 Feb 15 13:28 lib
\dashv drwxrwxrwt 15 root root 2048 Feb 17 14:53 tmp

```

See also: [popen2], page 975.
pclose (fid)
Close a file identifier that was opened by popen.
The function fclose may also be used for the same purpose.
See also: [fclose], page 296, [popen], page 974.
[in, out, pid] = popen2 (command, args)
Start a subprocess with two-way communication.
The name of the process is given by command, and args is an array or cell array of strings containing options for the command.
The file identifiers for the input and output streams of the subprocess are returned in in and out. If execution of the command is successful, pid contains the process ID of the subprocess. Otherwise, pid is -1 .
For example:
```

[in, out, pid] = popen2 ("sort", "-r");
fputs (in, "these\nare\nsome\nstrings\n");
fclose (in);
EAGAIN = errno ("EAGAIN");
done = false;
do
s = fgets (out);
if (ischar (s))
fputs (stdout, s);
elseif (errno () == EAGAIN)
pause (0.1);
fclear (out);
else
done = true;
endif
until (done)
fclose (out);
waitpid (pid);
these
| strings
\dashv some
\dashv are

```

Note that popen2, unlike popen, will not "reap" the child process. If you don't use waitpid to check the child's exit status, it will linger until Octave exits.
See also: [popen], page 974, [waitpid], page 977.
val = EXEC_PATH ()
old_val = EXEC_PATH (new_val)
EXEC_PATH (new_val, "local")
Query or set the internal variable that specifies a colon separated list of directories to append to the shell PATH when executing external programs.
The initial value of is taken from the environment variable OCTAVE_EXEC_PATH, but that value can be overridden by the command line argument --exec-path PATH.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [IMAGE_PATH], page 861, [OCTAVE_HOME], page 987, [OCTAVE_EXEC_HOME], page 987.

In most cases, the following functions simply decode their arguments and make the corresponding Unix system calls. For a complete example of how they can be used, look at the definition of the function popen2.
[pid, msg] = fork ()
Create a copy of the current process.
Fork can return one of the following values:
\(>0 \quad\) You are in the parent process. The value returned from fork is the process id of the child process. You should probably arrange to wait for any child processes to exit.
\(0 \quad\) You are in the child process. You can call exec to start another process. If that fails, you should probably call exit.
\(<0 \quad\) The call to fork failed for some reason. You must take evasive action. A system dependent error message will be waiting in msg.
[err, msg] = exec (file, args)
Replace current process with a new process.
Calling exec without first calling fork will terminate your current Octave process and replace it with the program named by file. For example,
```

exec ("ls", "-1")

```
will run ls and return you to your shell prompt.
If successful, exec does not return. If exec does return, err will be nonzero, and msg will contain a system-dependent error message.
[read_fd, write_fd, err, msg] = pipe ()
Create a pipe and return the reading and writing ends of the pipe into read_fd and write_fd respectively.

If successful, err is 0 and \(m s g\) is an empty string. Otherwise, err is nonzero and msg contains a system-dependent error message.
See also: [mkfifo], page 955.
[fid, msg] = dup2 (old, new)
Duplicate a file descriptor.
If successful, fid is greater than zero and contains the new file ID. Otherwise, fid is negative and msg contains a system-dependent error message.
See also: [fopen], page 295, [fclose], page 296, [fcntl], page 979.
[pid, status, msg] = waitpid (pid, options)
Wait for process pid to terminate.
The pid argument can be:
\(-1 \quad\) Wait for any child process.
\(0 \quad\) Wait for any child process whose process group ID is equal to that of the Octave interpreter process.
\(>0 \quad\) Wait for termination of the child process with ID pid.
The options argument can be a bitwise OR of zero or more of the following constants:
\(0 \quad\) Wait until signal is received or a child process exits (this is the default if the options argument is missing).
WNOHANG Do not hang if status is not immediately available.
wUNTRACED
Report the status of any child processes that are stopped, and whose status has not yet been reported since they stopped.

WCONTINUE
Return if a stopped child has been resumed by delivery of SIGCONT. This value may not be meaningful on all systems.

If the returned value of pid is greater than 0 , it is the process ID of the child process that exited. If an error occurs, pid will be less than zero and msg will contain a system-dependent error message. The value of status contains additional systemdependent information about the subprocess that exited.

See also: [WCONTINUE], page 977, [WCOREDUMP], page 978, [WEXITSTATUS], page 978, [WIFCONTINUED], page 978, [WIFSIGNALED], page 978, [WIFSTOPPED], page 978, [WNOHANG], page 979, [WSTOPSIG], page 979, [WTERMSIG], page 979, [WUNTRACED], page 979.

WCONTINUE ()
Return the numerical value of the WCONTINUE macro.
WCONTINUE is the option argument that may be passed to waitpid to indicate that it should also return if a stopped child has been resumed by delivery of a SIGCONT signal.
See also: [waitpid], page 977, [WNOHANG], page 979, [WUNTRACED], page 979.

WCOREDUMP (status)
Given status from a call to waitpid, return true if the child produced a core dump.
This function should only be employed if WIFSIGNALED returned true. The macro used to implement this function is not specified in POSIX.1-2001 and is not available on some Unix implementations (e.g., AIX, SunOS).

See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WIFSTOPPED], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WEXITSTATUS (status)
Given status from a call to waitpid, return the exit status of the child.
This function should only be employed if WIFEXITED returned true.
See also: [waitpid], page 977, [WIFEXITED], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WIFCONTINUED (status)
Given status from a call to waitpid, return true if the child process was resumed by delivery of SIGCONT.
See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WSTOPSIG], page 979.

WIFSIGNALED (status)
Given status from a call to waitpid, return true if the child process was terminated by a signal.

See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WIFSTOPPED (status)
Given status from a call to waitpid, return true if the child process was stopped by delivery of a signal.
This is only possible if the call was done using WUNTRACED or when the child is being traced (see ptrace(2)).

See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WIFEXITED (status)
Given status from a call to waitpid, return true if the child terminated normally.
See also: [waitpid], page 977, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WNOHANG ()
Return the numerical value of the WNOHANG macro.
WNOHANG is the option argument that may be passed to waitpid to indicate that it should return its status immediately instead of waiting for a process to exit.
See also: [waitpid], page 977, [WUNTRACED], page 979, [WCONTINUE], page 977.
WSTOPSIG (status)
Given status from a call to waitpid, return the number of the signal which caused the child to stop.
This function should only be employed if WIFSTOPPED returned true.
See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WTERMSIG], page 979, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WIFCONTINUED], page 978.

WTERMSIG (status)
Given status from a call to waitpid, return the number of the signal that caused the child process to terminate.
This function should only be employed if WIFSIGNALED returned true.
See also: [waitpid], page 977, [WIFEXITED], page 978, [WEXITSTATUS], page 978, [WIFSIGNALED], page 978, [WCOREDUMP], page 978, [WIFSTOPPED], page 978, [WSTOPSIG], page 979, [WIFCONTINUED], page 978.

WUNTRACED ()
Return the numerical value of the WUNTRACED macro.
WUNTRACED is the option argument that may be passed to waitpid to indicate that it should also return if the child process has stopped but is not traced via the ptrace system call
See also: [waitpid], page 977, [WNOHANG], page 979, [WCONTINUE], page 977.
fcntl (fid, request, arg)
[status, msg] = fcntl (fid, request, arg)
Change the properties of the open file fid.
The following values may be passed as request:
F_DUPFD Return a duplicate file descriptor.
F_GETFD Return the file descriptor flags for fid.
F_SETFD Set the file descriptor flags for fid.
F_GETFL Return the file status flags for fid. The following codes may be returned (some of the flags may be undefined on some systems).
O_RDONLY Open for reading only.
O_WRONLY Open for writing only.
O_RDWR Open for reading and writing.
O_APPEND Append on each write.
```

O_CREAT Create the file if it does not exist.
O_NONBLOCK

```
                    Non-blocking mode.
0_SYNC Wait for writes to complete.
O_ASYNC Asynchronous I/O.

F_SETFL Set the file status flags for fid to the value specified by arg. The only flags that can be changed are 0_APPEND and 0_NONBLOCK.

If successful, status is 0 and \(m s g\) is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.

See also: [fopen], page 295, [dup2], page 977.
kill (pid, sig)
[status, msg] = kill (pid, sig)
Send signal sig to process pid.
If pid is positive, then signal sig is sent to pid.
If pid is 0 , then signal sig is sent to every process in the process group of the current process.
If pid is -1 , then signal sig is sent to every process except process 1 .
If pid is less than -1 , then signal sig is sent to every process in the process group -pid.
If sig is 0 , then no signal is sent, but error checking is still performed.
If successful, status is 0 and \(m s g\) is an empty string. Otherwise, status is -1 and msg contains a system-dependent error message.

SIG ()
Return a structure containing Unix signal names and their defined values.

\subsection*{36.6 Process, Group, and User IDs}
pgid = getpgrp ()
Return the process group id of the current process.
pid \(=\) getpid ()
Return the process id of the current process.
See also: [getppid], page 980.
pid = getppid ()
Return the process id of the parent process.
See also: [getpid], page 980.
euid \(=\) geteuid ()
Return the effective user id of the current process.
See also: [getuid], page 981.
uid = getuid ()
Return the real user id of the current process.
See also: [geteuid], page 980.
egid = getegid ()
Return the effective group id of the current process.
See also: [getgid], page 981.
gid \(=\) getgid ()
Return the real group id of the current process.
See also: [getegid], page 981.

\subsection*{36.7 Environment Variables}
getenv (var)
Return the value of the environment variable var.
For example,
getenv ("PATH")
returns a string containing the value of your path.
See also: [setenv], page 981, [unsetenv], page 981.
setenv (var, value)
setenv (var)
putenv (...)
Set the value of the environment variable var to value.
If no value is specified then the variable will be assigned the null string.
See also: [unsetenv], page 981, [getenv], page 981.
status = unsetenv (var)
Delete the environment variable var.
Return 0 if the variable was deleted, or did not exist, and -1 if an error occurred.
See also: [setenv], page 981, [getenv], page 981.
homedir = get_home_directory ()
Return the current home directory.
On most systems, this is equivalent to getenv ("HOME"). On Windows systems, if the environment variable HOME is not set then it is equivalent to fullfile (getenv ("HOMEDRIVE"), getenv ("HOMEPATH"))

See also: [getenv], page 981.

\subsection*{36.8 Current Working Directory}
cd dir
cd
old_dir = cd
old_dir = cd (dir)
chdir ...
Change the current working directory to dir.
If called with no input or output arguments, the current directory is changed to the user's home directory ("~").
For example,
```

cd ~/octave

```
changes the current working directory to ~/octave. If the directory does not exist, an error message is printed and the working directory is not changed.
chdir is an alias for cd and can be used in all of the same calling formats.
Compatibility Note: When called with no arguments, matlab prints the present working directory rather than changing to the user's home directory.
See also: [pwd], page 983, [mkdir], page 954, [rmdir], page 955, [dir], page 983, [1s], page 982.

1s
ls filenames
ls options
ls options filenames
list = ls (...)
List directory contents.
The ls function forwards to the ls command if it is available. It falls back to calling the native operating system's directory listing command. Available options may vary from system to system.
Filenames are subject to shell expansion if they contain any wildcard characters ' \(*\) ', '?', ' []'. If these wildcard characters are escaped with a backslash ' \(\backslash\) ' (e.g., ' \(\backslash *\) ') then they are not treated as wildcards, but instead as the corresponding literal character. If the optional output list is requested then ls returns a character array with one row for each file/directory name.
Example usage on a UNIX-like system:
```

ls -l
\dashvt total 12

```

See also: [dir], page 983, [readdir], page 959, [glob], page 959, [what], page 148, [stat], page 956, [filesep], page 960, [ls_command], page 982.
val = ls_command ()
old_val = ls_command (new_val)
Query or set the shell command used by Octave's ls command.

See also: [ls], page 982.

\section*{dir}
dir (directory)
[list] = dir (directory)
Display file listing for directory directory.
If directory is not specified then list the present working directory.
If a return value is requested, return a structure array with the fields
name File or directory name.
folder Location of file or directory
date Timestamp of file modification (string value).
bytes File size in bytes.
isdir True if name is a directory.
datenum Timestamp of file modification as serial date number (double).
statinfo Information structure returned from stat.
If directory is a filename, rather than a directory, then return information about the named file. directory may also be a list rather than a single directory or file.
directory is subject to shell expansion if it contains any wildcard characters '*', '?', ' []'. If these wildcard characters are escaped with a backslash ' \(\backslash\) ' (e.g., ' \(\backslash *\) ') on a POSIX platform, then they are not treated as wildcards, but as the corresponding literal character. On Windows, it is not possible to escape wildcard characters because backslash ' \(\backslash\) ' is treated as a file separator. On Windows, use ls for file or folder names that contain characters that would be treated as wildcards by dir.
Note that for symbolic links, dir returns information about the file that the symbolic link points to rather than the link itself. However, if the link points to a nonexistent file, dir returns information about the link.
See also: [ls], page 982, [readdir], page 959, [glob], page 959, [what], page 148, [stat], page 956, [lstat], page 956.
pwd ()
dir \(=\) pwd ()
Return the current working directory.
See also: [cd], page 982, [dir], page 983, [ls], page 982, [mkdir], page 954, [rmdir], page 955 .

\subsection*{36.9 Password Database Functions}

Octave's password database functions return information in a structure with the following fields.
name The user name.
passwd The encrypted password, if available.
uid The numeric user id.
gid The numeric group id.
gecos The GECOS field.
dir The home directory.
shell The initial shell.
In the descriptions of the following functions, this data structure is referred to as a pw_struct.
pw_struct = getpwent ()
Return a structure containing an entry from the password database, opening it if necessary.
Once the end of the data has been reached, getpwent returns 0 .
See also: [setpwent], page 984, [endpwent], page 984.
pw_struct = getpwuid (uid).
Return a structure containing the first entry from the password database with the user ID uid.
If the user ID does not exist in the database, getpwuid returns 0 .
See also: [getpwnam], page 984.
pw_struct = getpwnam (name)
Return a structure containing the first entry from the password database with the user name name.
If the user name does not exist in the database, getpwname returns 0 .
See also: [getpwuid], page 984.
setpwent ()
Return the internal pointer to the beginning of the password database.
See also: [getpwent], page 984, [endpwent], page 984.
endpwent ()
Close the password database.
See also: [getpwent], page 984, [setpwent], page 984.

\subsection*{36.10 Group Database Functions}

Octave's group database functions return information in a structure with the following fields.
\begin{tabular}{ll} 
name & The user name. \\
passwd & The encrypted password, if available. \\
gid & The numeric group id. \\
mem & The members of the group.
\end{tabular}

In the descriptions of the following functions, this data structure is referred to as a grp_struct.
grp_struct = getgrent ()
Return an entry from the group database, opening it if necessary.
Once the end of data has been reached, getgrent returns 0 .
See also: [setgrent], page 985, [endgrent], page 985.
grp_struct = getgrgid (gid).
Return the first entry from the group database with the group ID gid.
If the group ID does not exist in the database, getgrgid returns 0 .
See also: [getgrnam], page 985.
grp_struct \(=\) getgrnam (name)
Return the first entry from the group database with the group name name.
If the group name does not exist in the database, getgrnam returns 0 .
See also: [getgrgid], page 985.
setgrent ()
Return the internal pointer to the beginning of the group database.
See also: [getgrent], page 984, [endgrent], page 985.
endgrent ()
Close the group database.
See also: [getgrent], page 984, [setgrent], page 985.

\subsection*{36.11 System Information}
```

computer ()
comp = computer ()
[comp, maxsize] = computer ()
[comp, maxsize, endian] = computer ()
arch = computer ("arch") that Octave is running on. example:

```
```

computer ()

```
computer ()
    \dashv x86_64-pc-linux-gnu
    \dashv x86_64-pc-linux-gnu
mycomp = computer ()
mycomp = computer ()
    => mycomp = x86_64-pc-linux-gnu
```

    => mycomp = x86_64-pc-linux-gnu
    ```

Print or return a string of the form cpu-vendor-os that identifies the type of computer

If invoked with an output argument, the value is returned instead of printed. For

If two output arguments are requested, also return the maximum number of elements for an array. This will depend on whether Octave has been compiled with 32-bit or 64 -bit index vectors.
If three output arguments are requested, also return the byte order of the current system as a character ("B" for big-endian or "L" for little-endian).

If the argument "arch" is specified, return a string indicating the architecture of the computer on which Octave is running.
Results may be different if Octave was invoked with the -traditional option.
See also: [isunix], page 986, [ismac], page 986, [ispc], page 986.
[uts, err, msg] = uname ()
Return system information in the structure.
For example:
```

uname ()
=> {
sysname = x86_64
nodename = segfault
release = 2.6.15-1-amd64-k8-smp
version = Linux
machine = \#2 SMP Thu Feb 23 04:57:49 UTC 2006
}

```

If successful, err is 0 and \(m s g\) is an empty string. Otherwise, err is nonzero and msg contains a system-dependent error message.
nproc ()
nproc (query)
Return the current number of available (logical) processors.
This returns the number of logical processors. For processors with hyperthreading, this is larger than the number of physical cores.
If called with the optional argument query, modify how processors are counted as follows:
all total number of processors.
current processors available to the current process.
overridable
same as current, but overridable through the OMP_NUM_THREADS environment variable.
ispc ()
Return true if Octave is running on a Windows system and false otherwise.
See also: [isunix], page 986, [ismac], page 986.
isunix ()
Return true if Octave is running on a Unix-like system and false otherwise.
See also: [ismac], page 986, [ispc], page 986.
ismac ()
Return true if Octave is running on a Mac OS X system and false otherwise.
See also: [isunix], page 986, [ispc], page 986.
isieee ()
Return true if your computer claims to conform to the IEEE standard for floating point calculations.
No actual tests are performed.
isdeployed ()
Return true if the current program has been compiled and is running separately from the Octave interpreter and false if it is running in the Octave interpreter.
Currently, this function always returns false in Octave.
isstudent ()
Return true if running in the student edition of matlab.
isstudent always returns false in Octave.
See also: [false], page 65.
OCTAVE_HOME ()
Return the name of the top-level Octave installation directory. OCTAVE_HOME corresponds to the configuration variable prefix.
See also: [EXEC_PATH], page 976, [IMAGE_PATH], page 861, [OCTAVE_EXEC_HOME], page 987.

OCTAVE_EXEC_HOME ()
Return the name of the top-level Octave installation directory for architecturedependent files. If not specified separately, the value is the same as OCTAVE_HOME. OCTAVE_EXEC_HOME corresponds to the configuration variable exec_prefix.
See also: [EXEC_PATH], page 976, [IMAGE_PATH], page 861, [OCTAVE_HOME], page 987.
matlabroot ()
Return the name of the top-level Octave installation directory.
This is an alias for the function OCTAVE_HOME provided for compatibility.
See also: [OCTAVE_HOME], page 987.
cfg_dir = user_config_dir ()
Return the (platform-specific) directory for user configuration.
See also: [user_data_dir], page 987.
data_dir = user_data_dir ()
Return the (platform-specific) directory for user data.
See also: [user_config_dir], page 987.
OCTAVE_VERSION ()
Return the version number of Octave as a string.
See also: [ver], page 988, [version], page 988.
```

v = version ()
[v, d] = version ()
v = version (feature)

```

Get version information for Octave.
If called without input argument, the first return value \(v\) gives the version number of Octave as a string. The second return value \(d\) holds the release date as a string.
The following options can be passed for feature:
"-date" for the release date of the running build,
"-description"
for a description of the release (always an empty string),
"-release"
for the name of the running build (always an empty string),
"-java" for version information of the Java VM,
"-fftw" for version information for the linked FFTW,
"-blas" for version information for the linked BLAS,
"-lapack"
for version information for the linked LAPACK.
"-hgid" the mercurial ID of the sources used to build Octave.
The information returned for the "-blas" and "-lapack" options might be unreliable. It might report which library was linked in when Octave was built instead of which library is currently used.
The variant with no input and output argument is an alias for the function OCTAVE_VERSION provided for compatibility.

See also: [OCTAVE_VERSION], page 987, [ver], page 988.

\section*{ver}
ver Octave
ver package
v = ver (...)
Display a header containing the current Octave version number, license string, and operating system. The header is followed by a list of installed packages, versions, and installation directories.

Use the package name package or Octave to query a specific component.
When called with an output argument, return a vector of structures describing Octave and each installed package. Each structure includes the following fields.

Name Package name.
Version Version of the package.
Release Release of the package (currently unused, defaults to []).
Date Date that the version was released.
See also: [version], page 988, [usejava], page 1058, [pkg], page 997.
compare_versions (v1, v2, operator)
Compare two version strings using the given operator.
This function assumes that versions v1 and v2 are arbitrarily long strings made of numeric and period characters possibly followed by an arbitrary string (e.g., "1.2.3", "0.3", "0.1.2+", or "1.2.3.4-test1").
The version is first split into numeric and character portions and then the parts are padded to be the same length (i.e., "1.1" would be padded to be "1.1.0" when being compared with "1.1.1", and separately, the character parts of the strings are padded with nulls).
The operator can be any logical operator from the set
- "==" equal
- "<" less than
- "<=" less than or equal to
- ">" greater than
- ">=" greater than or equal to
- "! =", "~ = " not equal

Note that version "1.1-test2" will compare as greater than "1.1-test10". Also, since the numeric part is compared first, "a" compares less than "1a" because the second string starts with a numeric part even though double ("a") is greater than double ("1").
verLessThan (package, version)
Return true if the installed version of the package is less than version.
package is the name of the package to check. Use "Octave" as the package to check the version of Octave itself.
version is the version to compare it to. A version is a string in the format accepted by compare_versions: an arbitrarily long string composed of numeric and period characters, possibly followed by an arbitrary string (e.g., "1.2.3", "0.3", "0.1.2+", or "1.2.3.4-test1").
Examples:
```

tf = verLessThan ("Octave", "5")
tf = 0
tf = verLessThan ("io", "2.4.12")

# ...

if (! verLessThan ("Octave", "5"))
\#\# ... use new Octave 5 features ...
endif

```

See also: [compare_versions], page 989, [version], page 988, [ver], page 988, [pkg], page 997.
license
license inuse
```

license inuse feature
license ("inuse")
retval = license ("inuse")
retval = license ("test", feature)
retval = license ("checkout", feature)
[retval, errmsg] = license ("checkout", feature)

```

Get license information for Octave and Octave packages.
GNU Octave is free software distributed under the GNU General Public License (GPL), and a license manager makes no sense. This function is provided only for mATLAB compatibility.
When called with no extra input arguments, it returns the Octave license, otherwise the first input defines the operation mode and must be one of the following strings: inuse, test, and checkout. The optional feature argument can either be "octave" (core), or an Octave package.
"inuse" Returns a list of loaded features, i.e., octave and the list of loaded packages. If an output is requested, it returns a struct array with the fields "feature", and "user".
"test" Return true if the specified feature is installed, false otherwise.
An optional third argument "enable" or "disable" is accepted but ignored.
"checkout"
Return true if the specified feature is installed, false otherwise. An optional second output will have an error message if a package is not installed.

See also: [pkg], page 997, [ver], page 988, [version], page 988.
memory ()
[userdata, systemdata] = memory ()
Display or return information about the memory usage of Octave.
If the function is called without output arguments, a table with an overview of the current memory consumption is displayed.
The output argument userdata is a structure with the following fields containing data for the Octave process:

MaxPossibleArrayBytes
Maximum size for an array to be allocated. Be aware that this includes all physical memory and swap space. Allocating that amount of memory might result in system instability, data corruption, and/or file system corruption. Note that depending on the platform (32-bit systems), the largest contiguous memory block might further limit the maximum possible allocatable array. This check is not currently implemented.

\section*{MemAvailableAllArrays}

The total size of available memory in bytes.
ram_available_all_arrays
The maximum size for an array that can be allocated in physical memory (excluding swap space). Note that depending on the platform (32-bit systems), the largest contiguous memory block might further limit the maximum possible allocatable array. This check is not currently implemented.

MemUsedMATLAB
mem_used_octave
The memory (including swap space) currently used by Octave in bytes.
ram_used_octave
The physical memory (excluding swap space) currently used by Octave in bytes.

The output argument systemdata is a nested structure with the following fields containing information about the system's memory:

\section*{PhysicalMemory.Available}

The currently available physical memory in bytes.
PhysicalMemory.Total
The total physical memory in bytes.
SystemMemory.Available
The currently available memory (including swap space) in bytes.

SystemMemory.Total
The total memory (including swap space) in bytes.

VirtualAddressSpace.Available
The currently available virtual address space in bytes.

VirtualAddressSpace.Total
The total virtual address space in bytes.

Example \#1 : print formatted table of memory usage
```

memory ()

# 

System RAM: 3934008 KiB, swap: 4087804 KiB
Octave RAM: 170596 KiB, virt: 1347944 KiB
Free RAM: 1954940 KiB, swap: 4087804 KiB
Available RAM: 2451948 KiB, total: 6042744 KiB

```

Example \#2 : return structs with memory usage information
```

[userdata, systemdata] = memory ()

# 

    userdata =
        scalar structure containing the fields:
        MaxPossibleArrayBytes = 6.1622e+09
        MemAvailableAllArrays = 6.1622e+09
        ram_available_all_arrays = 2.4883e+09
        MemUsedMATLAB = 1.3825e+09
        mem_used_octave = 1.3825e+09
        ram_used_octave = 1.7824e+08
    systemdata =
scalar structure containing the fields:
PhysicalMemory =
scalar structure containing the fields:
Available = 2.4954e+09
Total = 4.0284e+09
SystemMemory =
scalar structure containing the fields:
Available = 6.6813e+09
Total = 8.2143e+09
VirtualAddressSpace =
scalar structure containing the fields:
Available = 2.8147e+14
Total = 2.8147e+14

```

Programming Note: This function is implemented for Linux and Windows only.
See also: [computer], page 985, [getpid], page 980, [getrusage], page 992, [nproc], page 986 , [uname], page 986.
getrusage ()
Return a structure containing a number of statistics about the current Octave process.
Not all fields are available on all systems. If it is not possible to get CPU time statistics, the CPU time slots are set to zero. Other missing data are replaced by NaN. The list of possible fields is:
idrss Unshared data size.
inblock Number of block input operations.
isrss Unshared stack size.
ixrss \(\quad\) Shared memory size.
majflt Number of major page faults.
maxrss Maximum data size.
minflt Number of minor page faults.
msgrcv Number of messages received.
msgsnd Number of messages sent.
nivcsw Number of involuntary context switches.
nsignals Number of signals received.
nswap Number of swaps.
nvcsw \(\quad\) Number of voluntary context switches.
oublock Number of block output operations.
stime A structure containing the system CPU time used. The structure has the elements sec (seconds) usec (microseconds).
utime A structure containing the user CPU time used. The structure has the elements sec (seconds) usec (microseconds).
```

value = winqueryreg (rootkey, subkey, valuename)
value = winqueryreg (rootkey, subkey)
names = winqueryreg ("name", rootkey, subkey)

```

Query names or value from the Windows registry.
On Windows, return the value of the registry key subkey from the root key rootkey. You can specify the name of the queried registry value with the optional argument valuename. Otherwise, if called with only two arguments or valuename is empty, then the default value of subkey is returned. If the registry value is of type "REG_DWORD" then value is of class int32. If the value is of the type "REG_SZ" or "REG_EXPAND_SZ" a string is returned.

If the first argument is "name", a cell array of strings with the names of the values at that key is returned.

The variable rootkey must be a string with a valid root key identifier:

HKCR
HKEY_CLASSES_ROOT
HKEY_CURRENT_CONFIG
HKCU
HKEY_CURRENT_USER
HKLM
HKEY_LOCAL_MACHINE
HKU
HKEY_USERS
HKEY_PERFORMANCE_DATA
Examples:
Get a list of value names at the key 'HKCU\Environment':
```

valuenames = winqueryreg ("name", "HKEY_CURRENT_USER", ...
"Environment");

```

For each valuenames, display the value:
```

for k = 1:numel (valuenames)
val = winqueryreg ("HKEY_CURRENT_USER", "Environment", ...
valuenames{k});
str = sprintf ("%s = %s", valuenames{k}, num2str (val));
disp (str);
endfor

```

On non-Windows platforms this function fails with an error.

\subsection*{36.12 Hashing Functions}

It is often necessary to find if two strings or files are identical. This might be done by comparing them character by character and looking for differences. However, this can be slow, and so comparing a hash of the string or file can be a rapid way of finding if the files differ.

Another use of the hashing function is to check for file integrity. The user can check the hash of the file against a known value and find if the file they have is the same as the one that the original hash was produced with.

Octave supplies the hash function to calculate hash values of strings and files, the latter in combination with the fileread function. The hash function supports many commonly used hash methods.
hash (hfun, str)
Calculate the hash value of the string str using the hash function hfun.
The available hash functions are given in the table below.
'MD2' Message-Digest Algorithm 2 (RFC 1319).
'MD4' Message-Digest Algorithm 4 (RFC 1320).
'MD5' Message-Digest Algorithm 5 (RFC 1321).
'SHA1' Secure Hash Algorithm 1 (RFC 3174)
'SHA224' Secure Hash Algorithm 2 (224 Bits, RFC 3874)
'SHA256' Secure Hash Algorithm 2 (256 Bits, RFC 6234)
‘SHA384' Secure Hash Algorithm 2 (384 Bits, RFC 6234)
'SHA512' Secure Hash Algorithm 2 (512 Bits, RFC 6234)
To calculate for example the MD5 hash value of the string "abc" the hash function is called as follows:
```

hash ("md5", "abc")
\dashv ans = 900150983cd24fb0d6963f7d28e17f72

```

For the same string, the SHA-1 hash value is calculated with:
```

hash ("sha1", "abc")
\dashv ans = a9993e364706816aba3e25717850c26c9cd0d89d

```

And to compute the hash value of a file, e.g., file = "file.txt", call hash in combination with the fileread:
```

hash ("md5", fileread (file));

```

\section*{37 Packages}

Since Octave is Free Software users are encouraged to share their programs with others. To aid this sharing Octave supports the installation of extra packages. The 'Octave Forge' project is a community-maintained set of packages that can be downloaded and installed in Octave. At the time of writing the 'Octave Forge' project can be found online at https:// octave.sourceforge.io, but since the Internet is an ever-changing place this may not be true at the time of reading. Therefore it is recommended to see the Octave websitefor an updated reference.

\subsection*{37.1 Installing and Removing Packages}

Assuming a package is available in the file image-1.0.0.tar.gz it can be installed from the Octave prompt with the command
```

pkg install image-1.0.0.tar.gz

```

If the package is installed successfully nothing will be printed on the prompt, but if a warning or error occurred during installation it will be reported. It is possible to install several packages at once by writing several package file names after the pkg install command. If a different version of the package is already installed it will be removed prior to installing the new package. This makes it easy to upgrade and downgrade the version of a package, but makes it impossible to have several versions of the same package installed at once.

To see which packages are installed type
```

pkg list
\Package Name | Version | Installation directory

- ---------------+---------+--------------------------
\dashv image *| 1.0.0 | /home/jwe/octave/image-1.0.0

```

In this case, version 1.0.0 of the image package is installed. The ' \(*\) ' character next to the package name shows that the image package is loaded and ready for use.

It is possible to remove a package from the system using the pkg uninstall command like this
```

pkg uninstall image

```

If the package is removed successfully nothing will be printed in the prompt, but if a warning or error occurred it will be reported. It should be noted that the package file used for installation is not needed for removal, and that only the package name as reported by pkg list should be used when removing a package. It is possible to remove several packages at once by writing several package names after the pkg uninstall command.

To minimize the amount of code duplication between packages, it is possible that one package depends on another one. If a package depends on another, it will check if that package is installed during installation. If it is not, an error will be reported and the package will not be installed. This behavior can be disabled by passing the -nodeps flag to the pkg install command
```

pkg install -nodeps my_package_with_dependencies.tar.gz

```

Since the installed package expects its dependencies to be installed it may not function correctly. Because of this it is not recommended to disable dependency checking.

\section*{pkg command pkg_name}
pkg command option pkg_name
[out1, ...] = pkg (command, ...)
Manage or query packages (groups of add-on functions) for Octave.
Packages can be installed globally (i.e., for all users of the system) or locally (i.e., for the current user only).

Global packages are installed by default in a system-wide location. This is usually a subdirectory of the folder where Octave itself is installed. Therefore, Octave needs write access to this folder to install global packages, which is usually only available when Octave is run with administrative privileges, such as when run as root (or superuser) on Unix-like systems, or run with elevated privileges ("Run as administrator") on Windows.

In contrast, local packages are installed by default in the user's home directory (or user profile on Windows) and are only available to that specific user. Usually, they can be installed without administrative privileges.
When Octave is running with administrative privileges, pkg will install packages to the global package location by default. Otherwise, packages will be installed to the local location by default. The user can override this default installation location with optional arguments (-local or -global) as described below. The currently used default package installation location can be queried with pkg prefix.
For global and local packages, there are separate databases holding the information about the installed packages. If some package is installed globally as well as locally, the local installation takes precedence over ("shadows") the global one. Which (global or local) package installation is used can also be manipulated by using prefixes and/or using the 'local_list' input argument. Using these mechanisms, several different releases of the same package can be installed side by side as well (but cannot be loaded simultaneously).

Packages might depend on external software and/or other packages. To be able to install such packages, these dependencies should be installed beforehand. A package that depends on other package(s) can still be installed using the -nodeps flag. The effects of unsatisfied dependencies on external software - like libraries-depends on the individual package.
Packages must be loaded before they can be used. When loading a package, Octave performs the following tasks:
1. If the package depends on other packages (and pkg load is called without the -nodeps option), the package is not loaded immediately. Instead, those dependencies are loaded first (recursively if needed).
2. When all dependencies are satisfied, the package's subdirectories are added to the search path.

This load order leads to functions that are provided by dependencies being potentially shadowed by functions of the same name that are provided by top-level packages.
Each time, a package is added to the search path, initialization script(s) for the package are automatically executed if they are provided by the package.

Depending on the value of command and on the number of requested return arguments, pkg can be used to perform several tasks. Possible values for command are:
'install' Install named packages. For example,
```

pkg install image-1.0.0.tar.gz

```
installs the package found in the file image-1.0.0.tar.gz. The file containing the package can be a URL, e.g.,
```

pkg install 'http://somewebsite.org/image-1.0.0.tar.gz'

```
installs the package found in the given URL. This requires an internet connection and the cURL library.
Security risk: no verification of the package is performed before the installation. It has the same security issues as manually downloading the package from the given URL and installing it.
No support: the GNU Octave community is not responsible for packages installed from foreign sites. For support or for reporting bugs you need to contact the maintainers of the installed package directly (see the DESCRIPTION file of the package)
The option variable can contain options that affect the manner in which a package is installed. These options can be one or more of
-nodeps The package manager will disable dependency checking. With this option it is possible to install a package even when it depends on another package which is not installed on the system. Use this option with care.
-local A local installation (package available only to current user) is forced, even if Octave is being run with administrative privileges.
-global A global installation (package available to all users) is forced, even if Octave is not being run with administrative privileges. The user must have write access to the global package store.
-forge Install a package directly from the Octave Forge repository. This requires an internet connection and the cURL library.
Security risk: no verification of the package is performed before the installation. There are no signatures for packages, or checksums to confirm the correct file was downloaded. It has the same security issues as manually downloading the package from the Octave Forge repository and installing it.
-verbose The package manager will print the output of all commands as they are performed.
'update' Check installed Octave Forge packages against repository and update any outdated items. Updated packages are installed either globally or locally depending on whether Octave is running with elevated privileges. This requires an internet connection and the cURL library.

Options for the install command and the names of individual packages to be checked for updates may be specified as a list following the update command. If the -local or -global option is specified, pkg update limits the update check to the local or global installed packages, and installs updates in that same context. For example,
Update all packages:
pkg update
Update all local packages:
pkg update -local
Update certain packages, ignore dependencies, max verbosity:
```

pkg update -verbose -nodeps image signal geometry

```

Updates for multiple packages are sorted alphabetically and not checked for dependencies affected by installation order. If dependency order related pkg update failure occurs, use pkg update -nodeps to ignore dependencies, or pkg install -forge <package_name> to update individual packages manually.
'uninstall'
Uninstall named packages. For example, pkg uninstall image
removes the image package from the system. If another installed package depends on the image package an error will be issued. The package can be uninstalled anyway by using the -nodeps option.
'load' Add named packages to the path. After loading a package it is possible to use the functions provided by the package. For example,
```

pkg load image

```
adds the image package to the path.
Note: When loading a package, pkg will automatically try to load any unloaded dependencies as well, unless the -nodeps flag has been specified. For example,
```

pkg load signal

```
adds the signal package and also tries to load its dependency: the control package. Be aware that the functionality of package(s) loaded will probably be impacted by use of the -nodeps flag. Even if necessary dependencies are loaded later, the functionality of top-level packages can still be affected because the optimal loading order may not have been followed.
'unload' Remove named packages from the path. After unloading a package it is no longer possible to use the functions provided by the package. Trying to unload a package that other loaded packages still depend on will result in an error; no packages will be unloaded in this case. A package can be forcibly removed with the -nodeps flag, but be aware that the functionality of dependent packages will likely be affected. As when loading
packages, reloading dependencies after having unloaded them with the -nodeps flag may not restore all functionality of the dependent packages as the required loading order may be incorrect.
'list' Show the list of currently installed packages. For example,
pkg list
will produce a short report with the package name, version, and installation directory for each installed package. Supply a package name to limit reporting to a particular package. For example:
```

pkg list image

```

If a single return argument is requested then pkg returns a cell array where each element is a structure with information on a single package.
```

installed_packages = pkg ("list")

```

If two output arguments are requested pkg splits the list of installed packages into those which were installed by the current user, and those which were installed by the system administrator.
```

[user_packages, system_packages] = pkg ("list")

```

The "-forge" option lists packages available at the Octave Forge repository. This requires an internet connection and the cURL library. For example:
```

oct_forge_pkgs = pkg ("list", "-forge")

```
'describe'

Show a short description of installed packages. With the option "-verbose" also list functions provided by the package. For example,
```

pkg describe -verbose

```
will describe all installed packages and the functions they provide. Display can be limited to a set of packages:
```


## describe control and signal packages

pkg describe control signal

```

If one output is requested a cell of structure containing the description and list of functions of each package is returned as output rather than printed on screen:
```

desc = pkg ("describe", "secs1d", "image")

```

If any of the requested packages is not installed, pkg returns an error, unless a second output is requested:
[desc, flag] = pkg ("describe", "secs1d", "image")
flag will take one of the values "Not installed", "Loaded", or "Not loaded" for each of the named packages.
'prefix' Set the installation prefix directory. For example,
pkg prefix ~/my_octave_packages
sets the installation prefix to ~/my_octave_packages. Packages will be installed in this directory.

It is possible to get the current installation prefix by requesting an output argument. For example:
```

pfx = pkg ("prefix")

```

The location in which to install the architecture dependent files can be independently specified with an addition argument. For example:
pkg prefix ~/my_octave_packages ~/my_arch_dep_pkgs
'local_list'
Set the file in which to look for information on locally installed packages. Locally installed packages are those that are available only to the current user. For example:
```

pkg local_list ~/.octave_packages

```

It is possible to get the current value of local_list with the following
pkg local_list
'global_list'
Set the file in which to look for information on globally installed packages. Globally installed packages are those that are available to all users. For example:
pkg global_list /usr/share/octave/octave_packages
It is possible to get the current value of global_list with the following
pkg global_list
'build' Build a binary form of a package or packages. The binary file produced will itself be an Octave package that can be installed normally with pkg. The form of the command to build a binary package is
pkg build builddir image-1.0.0.tar.gz ...
where builddir is the name of a directory where the temporary installation will be produced and the binary packages will be found. The options -verbose and -nodeps are respected, while all other options are ignored.
'rebuild' Rebuild the package database from the installed directories. This can be used in cases where the package database has been corrupted.
'test' Perform the built-in self tests contained in all functions provided by the named packages. For example:
pkg test image
See also: [ver], page 988, [news], page 22.

\subsection*{37.2 Using Packages}

By default installed packages are not available from the Octave prompt, but it is possible to control this using the pkg load and pkg unload commands. The functions from a package can be added to the Octave path by typing
```

pkg load package_name

```
where package_name is the name of the package to be added to the path.
In much the same way a package can be removed from the Octave path by typing pkg unload package_name

\subsection*{37.3 Administrating Packages}

It is possible to make both per-user (local) and system-wide (global) installations of a package. If the user performing the installation is root (or Administrator with elevated privileges on Windows), the packages by default install in a system-wide directory that defaults to OCTAVE_HOME/share/octave/packages/. If the user is not root (or Octave is running without elevated privileges), packages are installed locally. The default installation directory for local packages is user_data_dir/octave/OCTAVE_API_VERSION/packages. Packages will be installed in a subdirectory of the installation directory that will be named after the package. It is possible to change the installation directory by using the pkg prefix command:
pkg prefix new_installation_directory
The current installation directory can be retrieved by typing
current_installation_directory = pkg ("prefix")
The package manager stores some information about the installed packages in configuration files. For per-user (local) packages, this information is stored in the file user_config_ dir/octave/OCTAVE_API_VERSION/octave_packages by default. For system-wide (global) installations, it is stored in OCTAVE_HOME/share/octave/octave_packages. The path to the per-user file can be changed with the pkg local_list command:
pkg local_list /path/to/new_file
For system-wide installations, this can be changed in the same way using the pkg global_ list command. If these commands are called without a new path, the current path will be returned. To retain these settings between sessions, they can be set in one of the startup files, see Section 2.1.2 [Startup Files], page 18.

\subsection*{37.4 Creating Packages}

Internally a package is simply a gzipped tar file that contains a top level directory of any given name. This directory will in the following be referred to as package and may contain the following files:
```

package/CITATION

```

This is am optional file describing instructions on how to cite the package for publication. It will be displayed verbatim by the function citation.
package/COPYING
This is a required file containing the license of the package. No restrictions is made on the license in general. If however the package contains dynamically linked functions the license must be compatible with the GNU General Public License.
package/DESCRIPTION
This is a required file containing information about the package. See Section 37.4.1 [The DESCRIPTION File], page 1005, for details on this file.
package/ChangeLog
This is an optional file describing all the changes made to the package source files.
package/INDEX
This is an optional file describing the functions provided by the package. If this file is not given then one with be created automatically from the functions in the package and the Categories keyword in the DESCRIPTION file. See Section 37.4.2 [The INDEX File], page 1007, for details on this file.
package/NEWS
This is an optional file describing all user-visible changes worth mentioning. As this file increases on size, old entries can be moved into package/ONEWS.
package/ONEWS
This is an optional file describing old entries from the NEWS file.
package/PKG_ADD
An optional file that includes commands that are run when the package is added to the users path. Note that PKG_ADD directives in the source code of the package will also be added to this file by the Octave package manager. Note that symbolic links are to be avoided in packages, as symbolic links do not exist on some file systems, and so a typical use for this file is the replacement of the symbolic link
```

ln -s foo.oct bar.oct

```
with an autoload directive like
```

autoload ('bar', which ('foo'));

```

See Section 37.4.3 [PKG_ADD and PKG_DEL Directives], page 1008, for details on PKG_ADD directives.
package/PKG_DEL
An optional file that includes commands that are run when the package is removed from the users path. Note that PKG_DEL directives in the source code of the package will also be added to this file by the Octave package manager. See Section 37.4.3 [PKG_ADD and PKG_DEL Directives], page 1008, for details on PKG_DEL directives.
package/pre_install.m
This is an optional function that is run prior to the installation of a package. This function is called with a single argument, a struct with fields names after the data in the DESCRIPTION, and the paths where the package functions will be installed.
package/post_install.m
This is an optional function that is run after the installation of a package. This function is called with a single argument, a struct with fields names after the data in the DESCRIPTION, and the paths where the package functions were installed.
package/on_uninstall.m
This is an optional function that is run prior to the removal of a package. This function is called with a single argument, a struct with fields names after the data in the DESCRIPTION, the paths where the package functions are installed, and whether the package is currently loaded.

Besides the above mentioned files, a package can also contain one or more of the following directories:
package/inst
An optional directory containing any files that are directly installed by the package. Typically this will include any m-files.
package/src
An optional directory containing code that must be built prior to the packages installation. The Octave package manager will execute ./configure in this directory if this script exists, and will then call make if a file Makefile exists in this directory. make install will however not be called. The environment variables MKOCTFILE, OCTAVE_CONFIG, and OCTAVE will be set to the full paths of the programs mkoctfile, octave-config, and octave, respectively, of the correct version when configure and make are called. If a file called FILES exists all files listed there will be copied to the inst directory, so they also will be installed. If the FILES file doesn't exist, src/*.m and src/*. oct will be copied to the inst directory.
package/doc
An optional directory containing documentation for the package. The files in this directory will be directly installed in a sub-directory of the installed package for future reference.
package/bin
An optional directory containing files that will be added to the Octave EXEC_PATH when the package is loaded. This might contain external scripts, etc., called by functions within the package.

\subsection*{37.4.1 The DESCRIPTION File}

The DESCRIPTION file contains various information about the package, such as its name, author, and version. This file has a very simple format
- Lines starting with '\#' are comments.
- Lines starting with a blank character are continuations from the previous line.
- Everything else is of the form NameOfOption: ValueOfOption.

The following is a simple example of a DESCRIPTION file
```

Name: The name of my package
Version: 1.0.0
Date: 2007-18-04
Author: The name (and possibly email) of the package author.
Maintainer: The name (and possibly email) of the current
package maintainer.
Title: The title of the package
Description: A short description of the package. If this
description gets too long for one line it can continue
on the next by adding a space to the beginning of the
following lines.
License: GPLv3+

```

The package manager currently recognizes the following keywords
Name Name of the package.
Version Version of the package. A package version is typically digits separated by dots but may also contain '+', '-', '~', and alphanumeric characters (in the "C" locale). For example, " \(2.1 .0+\) " could indicate a development version of a package. Versions are compared using [compare_versions], page 989.
Date Date of last update.
Author Original author of the package.
Maintainer
Maintainer of the package.
Title A one line description of the package.
Description
A one paragraph description of the package.
Categories
Optional keyword describing the package (if no INDEX file is given this is mandatory).
Problems Optional list of known problems.
Url Optional list of homepages related to the package.
Depends A list of other Octave packages that this package depends on. This can include dependencies on particular versions, with the following format:

Depends: package (>= 1.0.0)
Possible operators are \(<,<=,==,>=\) or \(>\). If the part of the dependency in () is missing, any version of the package is acceptable. Multiple dependencies can be defined as a comma separated list. This can be used to define a range of versions of a particular package:
```

Depends: package (>= 1.0.0), package (< 1.5.0)

```

It is also possible to depend on particular versions of Octave core:
Depends: octave (>= 3.8.0)
License An optional short description of the used license (e.g., GPL version 3 or newer). This is optional since the file COPYING is mandatory.

SystemRequirements
These are the external install dependencies of the package and are not checked by the package manager. This is here as a hint to the distribution packager. They follow the same conventions as the Depends keyword.

BuildRequires
These are the external build dependencies of the package and are not checked by the package manager. This is here as a hint to the distribution packager. They follow the same conventions as the Depends keyword. Note that in general, packaging systems such as rpm or deb autoprobe the install dependencies from the build dependencies, and therefore a BuildRequires dependency usually removes the need for a SystemRequirements dependency.

The developer is free to add additional arguments to the DESCRIPTION file for their own purposes. One further detail to aid the packager is that the SystemRequirements and BuildRequires keywords can have a distribution dependent section, and the automatic build process will use these. An example of the format of this is

BuildRequires: libtermcap-devel [Mandriva] libtermcap2-devel
where the first package name will be used as a default and if the RPMs are built on a Mandriva distribution, then the second package name will be used instead.

\subsection*{37.4.2 The INDEX File}

The optional INDEX file provides a categorical view of the functions in the package. This file has a very simple format
- Lines beginning with ' \(\#\) ' are comments.
- The first non-comment line should look like this
```

toolbox >> Toolbox name

```
- Lines beginning with an alphabetical character indicates a new category of functions.
- Lines starting with a white space character indicate that the function names on the line belong to the last mentioned category.
The format can be summarized with the following example:
```


# A comment

toolbox >> Toolbox name
Category Name 1
function1 function2 function3
function4
Category Name 2
function2 function5

```

If you wish to refer to a function that users might expect to find in your package but is not there, providing a work around or pointing out that the function is available elsewhere, you can use:
```

fn = workaround description

```

This workaround description will not appear when listing functions in the package with pkg describe but they will be published in the HTML documentation online. Workaround descriptions can use any HTML markup, but keep in mind that it will be enclosed in a bold-italic environment. For the special case of:
```

fn = use <code>alternate expression</code>

```
the bold-italic is automatically suppressed. You will need to use <code> even in references:
```

fn = use <a href="someothersite.html"><code>fn</code></a>

```

Sometimes functions are only partially compatible, in which case you can list the noncompatible cases separately. To refer to another function in the package, use <f>fn</f>. For example:
```

eig (a, b) = use <f>qz</f>

```

Since sites may have many missing functions, you can define a macro rather than typing the same link over and again.
```

\$id = expansion

```
defines the macro id. You can use \(\$\) id anywhere in the description and it will be expanded. For example:
```

\$TSA = see <a href="link_to_spctools">SPC Tools</a>
arcov = \$TSA <code>armcv</code>

```
id is any string of letters, numbers and _.

\subsection*{37.4.3 PKG_ADD and PKG_DEL Directives}

If the package contains files called PKG_ADD or PKG_DEL the commands in these files will be executed when the package is added or removed from the users path. In some situations such files are a bit cumbersome to maintain, so the package manager supports automatic creation of such files. If a source file in the package contains a PKG_ADD or PKG_DEL directive they will be added to either the PKG_ADD or PKG_DEL files.

In m-files a PKG_ADD directive looks like this
```


## PKG_ADD: some_octave_command

```

Such lines should be added before the function keyword. In C++ files a PKG_ADD directive looks like this
```

// PKG_ADD: some_octave_command

```

In both cases some_octave_command should be replaced by the command that should be placed in the PKG_ADD file. PKG_DEL directives work in the same way, except the PKG_ADD keyword is replaced with PKG_DEL and the commands get added to the PKG_DEL file.

\subsection*{37.4.4 Missing Components}

If a package relies on a component, such as another Octave package, that may not be present it may be useful to install a function which informs users what to do when a particular component is missing. The function must be written by the package maintainer and registered with Octave using missing_component_hook.
```

val = missing_component_hook ()
old_val = missing_component_hook (new_val)
missing_component_hook (new_val, "local")

```

Query or set the internal variable that specifies the function to call when a component of Octave is missing.
This can be useful for packagers that may split the Octave installation into multiple sub-packages, for example, to provide a hint to users for how to install the missing components.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
The hook function is expected to be of the form
```

fcn (component)

```

Octave will call fcn with the name of the function that requires the component and a string describing the missing component. The hook function should return an error message to be displayed.
See also: [missing_function_hook], page 1100.

\section*{Appendix A External Code Interface}
"The sum of human wisdom is not contained in any one language"

\section*{- Ezra Pound}

Octave is a fantastic language for solving many problems in science and engineering. However, it is not the only computer language and there are times when you may want to use code written in other languages. Good reasons for doing so include: 1) not re-inventing the wheel; existing function libraries which have been thoroughly tested and debugged or large scale simulation codebases are a good example, 2) accessing unique capabilities of a different language; for example the well-known regular expression functions of Perl (but don't do that because regexp already exists in Octave).

Performance should generally not be a reason for using compiled extensions. Although compiled extensions can run faster, particularly if they replace a loop in Octave code, this is almost never the best path to take. First, there are many techniques to speed up Octave performance while remaining within the language. Second, Octave is a highlevel language that makes it easy to perform common mathematical tasks. Giving that up means shifting the focus from solving the real problem to solving a computer programming problem. It means returning to low-level constructs such as pointers, memory management, mathematical overflow/underflow, etc. Because of the low level nature, and the fact that the compiled code is executed outside of Octave, there is the very real possibility of crashing the interpreter and losing work.

Before going further, you should first determine if you really need to bother writing code outside of Octave.
- Can I get the same functionality using the Octave scripting language alone?

Even when a function already exists outside the language, it may be better to simply reproduce the behavior in an m-file rather than attempt to interface to the outside code.
- Is the code thoroughly optimized for Octave?

If performance is an issue you should always start with the in-language techniques for getting better performance. Chief among these is vectorization (see Chapter 19 [Vectorization and Faster Code Execution], page 625) which not only makes the code concise and more understandable but improves performance (10X-100X). If loops must be used, make sure that the allocation of space for variables takes place outside the loops using an assignment to a matrix of the right size, or zeros.
- Does the code make as much use as possible of existing built-in library routines?

These routines are highly optimized and many do not carry the overhead of being interpreted.
- Does writing a dynamically linked function represent a useful investment of your time, relative to staying in Octave?
It will take time to learn Octave's interface for external code and there will inevitably be issues with tools such as compilers.
With that said, Octave offers a versatile interface for including chunks of compiled code as dynamically linked extensions. These dynamically linked functions can be called from the interpreter in the same manner as any ordinary function. The interface is bi-directional and
external code can call Octave functions (like plot) which otherwise might be very difficult to develop.

The interface is centered around supporting the languages C++, C, and Fortran. Octave itself is written in C++ and can call external C++/C code through its native oct-file interface. The C language is also supported through the mex-file interface for compatibility with matlab. Fortran code is easiest to reach through the oct-file interface.

Because many other languages provide C or \(\mathrm{C}++\) APIs it is relatively simple to build bridges between Octave and other languages. This is also a way to bridge to hardware resources which often have device drivers written in C.

\section*{A. 1 Oct-Files}

\section*{A.1.1 Getting Started with Oct-Files}

Oct-files are pieces of C++ code that have been compiled with the Octave API into a dynamically loadable object. They take their name from the file which contains the object which has the extension .oct.

Finding a C++ compiler, using the correct switches, adding the right include paths for header files, etc. is a difficult task. Octave automates this by providing the mkoctfile command with which to build oct-files. The command is available from within Octave or at the shell command line.
mkoctfile [-options] file...
[output, status] = mkoctfile (...)
The mkoctfile function compiles source code written in C, C++, or Fortran. Depending on the options used with mkoctfile, the compiled code can be called within Octave or can be used as a stand-alone application.
mkoctfile can be called from the shell prompt or from the Octave prompt. Calling it from the Octave prompt simply delegates the call to the shell prompt. Any output is stored in the output variable and the exit status in the status variable. If called with no outputs and the compilation fails then Octave will emit an error. If the programmer requests output or status, however, Octave will merely issue a warning and it is the programmer's responsibility to verify the command was successful.
mkoctfile accepts the following options, all of which are optional except for the filename of the code you wish to compile:
'-I DIR' Add the include directory DIR to compile commands.
'-D DEF' Add the definition DEF to the compiler call.
'-l LIB' Add the library LIB to the link command.
'-L DIR' Add the library directory DIR to the link command.
'-M'
'--depend'
Generate dependency files (.d) for C and C++ source files.
'-R DIR' Add the run-time path to the link command.
'-Wl, ...' Pass options to the linker like "-Wl,-rpath=...". The quotes are needed since commas are interpreted as command separators.
'-W...' Pass options to the assembler like "-Wa,OPTION".
'-c' Compile but do not link.
'-g' Enable debugging options for compilers.
'-o FILE'
'--output FILE'
Output filename. Default extension is .oct (or .mex if '--mex' is specified) unless linking a stand-alone executable.
'-p VAR'
'--print VAR'
Print configuration variable VAR. There are three categories of variables:
Octave configuration variables that users may override with environment variables. These are used in commands that mkoctfile executes.
\begin{tabular}{ll} 
ALL_CFLAGS & INCLUDEDIR \\
ALL_CXXFLAGS & LAPACK_LIBS \\
ALL_FFLAGS & LDFLAGS \\
ALL_LDFLAGS & LD_STATIC_FLAG \\
BLAS_LIBS & LFLAGS \\
CC & LIBDIR \\
CFLAGS & LIBOCTAVE \\
CPICFLAG & LIBOCTINTERP \\
CPPFLAGS & OCTAVE_LINK_OPTS \\
CXX & OCTINCLUDEDIR \\
CXXFLAGS & OCTAVE_LIBS \\
CXXLD & OCTAVE_LINK_DEPS \\
CXXPICFLAG & OCTLIBDIR \\
DL_LDFLAGS & OCT_LINK_DEPS \\
F77 & OCT_LINK_OPTS \\
F77_INTEGER8_FLAG & RDYNAMIC_FLAG \\
FFLAGS & SPECIAL_MATH_LIB \\
FPICFLAG & XTRA_CFLAGS \\
INCFLAGS & XTRA_CXXFLAGS
\end{tabular}

Octave configuration variables as above, but currently unused by mkoctfile.
```

AR
DEPEND_EXTRA_SED_PATTERN
DEPEND_FLAGS
FFTW3F_LDFLAGS
FFTW3F_LIBS
FFTW3_LDFLAGS
FFTW3_LIBS
FFTW_LIBS
FLIBS
LIBS
RANLIB
READLINE_LIBS

```

Octave configuration variables that are provided for informational purposes only. Except for 'OCTAVE_HOME' and 'OCTAVE_EXEC_HOME', users may not override these variables.

If OCTAVE_HOME or OCTAVE_EXEC_HOME are set in the environment, then other variables are adjusted accordingly with OCTAVE_HOME or OCTAVE_EXEC_HOME substituted for the original value of the directory specified by the --prefix or --exec-prefix options that were used when Octave was configured.
\begin{tabular}{ll} 
API_VERSION & LOCALFCNFILEDIR \\
ARCHLIBDIR & LOCALOCTFILEDIR \\
BINDIR & LOCALSTARTUPFILEDIR \\
CANONICAL_HOST_TYPE & LOCALVERARCHLIBDIR \\
DATADIR & LOCALVERFCNFILEDIR \\
DATAROOTDIR & LOCALVEROCTFILEDIR \\
DEFAULT_PAGER & MAN1DIR \\
EXEC_PREFIX & MAN1EXT \\
EXEEXT & MANDIR \\
FCNFILEDIR & OCTAVE_EXEC_HOME \\
IMAGEDIR & OCTAVE_HOME \\
INFODIR & OCTAVE_VERSION \\
INFOFILE & OCTDATADIR \\
LIBEXECDIR & OCTDOCDIR \\
LOCALAPIARCHLIBDIR & OCTFILEDIR \\
LOCALAPIFCNFILEDIR & OCTFONTSDIR \\
LOCALAPIOCTFILEDIR & STARTUPFILEDIR \\
LOCALARCHLIBDIR &
\end{tabular}
‘--link-stand-alone’
Link a stand-alone executable file.
'--mex' Assume creation of a MEX file. Set the default output extension to .mex.
'-s'
'--strip' Strip the output file.
'-v'
'--verbose'
Echo commands as they are executed.
'file' The file to compile or link. Recognized file types are:
\begin{tabular}{ll}
.c & C source \\
.cc & C++ source \\
.cp & C++ source \\
.cpp & C++ source \\
.CPP & C++ source \\
.cxx & C++ source \\
.c++ & C++ source \\
.C & C++ source \\
.f & Fortran source (fixed form) \\
.F & Fortran source (fixed form) \\
.f90 & Fortran source (free form) \\
.F90 & Fortran source (free form) \\
.o & object file \\
.a & library file
\end{tabular}

Consider the following short example which introduces the basics of writing a C++ function that can be linked to Octave.
```

\#include <octave/oct.h>
DEFUN_DLD (helloworld, args, nargout,
"Hello World Help String")
{
octave_stdout << "Hello World has "
<< args.length () << " input arguments and "
<< nargout << " output arguments.\n";
// Return empty matrices for any outputs
octave_value_list retval (nargout);
for (int i = 0; i < nargout; i++)
retval(i) = octave_value (Matrix ());
return retval;
}

```

The first critical line is \#include <octave/oct.h> which makes available most of the definitions necessary for a C++ oct-file. Note that octave/oct.h is a C++ header and cannot be directly \#include'ed in a C source file, nor any other language.

Included by oct.h is a definition for the macro DEFUN_DLD which creates a dynamically loaded function. This macro takes four arguments:
1. The function name as it will be seen in Octave,
2. The list of arguments to the function of type octave_value_list,
3. The number of output arguments, which can be - and often is - omitted if not used, and
4. The string to use for the help text of the function.

The return type of functions defined with DEFUN_DLD is always octave_value_list.
There are a couple of important considerations in the choice of function name. First, it must be a valid Octave function name and so must be a sequence of letters, digits, and underscores not starting with a digit. Second, as Octave uses the function name to define the filename it attempts to find the function in, the function name in the DEFUN_DLD macro must match the filename of the oct-file. Therefore, the above function should be in a file helloworld.cc, and would be compiled to an oct-file using the command
```

mkoctfile helloworld.cc

```

This will create a file called helloworld. oct that is the compiled version of the function. It should be noted that it is perfectly acceptable to have more than one DEFUN_DLD function in a source file. However, there must either be a symbolic link to the oct-file for each of the functions defined in the source code with the DEFUN_DLD macro or the autoload (Section 11.10 [Function Files], page 210) function should be used.

The rest of the function shows how to find the number of input arguments, how to print through the Octave pager, and how to return from the function. After compiling this function as above, an example of its use is
```

helloworld (1, 2, 3)

```
\(\dashv\) Hello World has 3 input arguments and 0 output arguments.
Subsequent sections show how to use specific classes from Octave's core internals. Base classes like dMatrix (a matrix of double values) are found in the directory liboctave/array. The definitive reference for how to use a particular class is the header file itself. However, it is often enough simply to study the examples in the manual in order to be able to use a class.

\section*{A.1.2 Matrices and Arrays in Oct-Files}

Octave supports a number of different array and matrix classes, the majority of which are based on the Array class. The exception are the sparse matrix types discussed separately below. There are three basic matrix types:

Matrix A double precision matrix class defined in dMatrix.h
ComplexMatrix
A complex matrix class defined in CMatrix.h

\section*{BoolMatrix}

A boolean matrix class defined in boolMatrix.h
These are the basic two-dimensional matrix types of Octave. In addition there are a number of multi-dimensional array types including

NDArray A double precision array class defined in dNDArray.h
ComplexNDarray
A complex array class defined in CNDArray.h
boolNDArray
A boolean array class defined in boolNDArray.h
```

int8NDArray
int16NDArray
int32NDArray
int64NDArray
8, 16, 32, and 64-bit signed array classes defined in int8NDArray.h,
int16NDArray.h, etc.
uint8NDArray
uint16NDArray
uint32NDArray
uint64NDArray
8, 16, 32, and 64 -bit unsigned array classes defined in uint8NDArray.h, uint16NDArray.h, etc.

```

There are several basic ways of constructing matrices or multi-dimensional arrays. Using the class Matrix as an example one can
- Create an empty matrix or array with the empty constructor. For example:
```

Matrix a;

```

This can be used for all matrix and array types.
- Define the dimensions of the matrix or array with a dim_vector which has the same characteristics as the vector returned from size. For example:
```

dim_vector dv (2, 3); // 2 rows, 3 columns
Matrix a (dv);

```

This can be used for all matrix and array types.
- Define the number of rows and columns in the matrix. For example:
```

Matrix a (2, 2)

```

This constructor can only be used with matrix types.
These types all share a number of basic methods and operators. Many bear a resemblance to functions that exist in the interpreter. A selection of useful methods include
```

T\& operator () (octave_idx_type)
T\& elem (octave_idx_type)
[Method]
T\& elem (octave_idx_type)
[Method]

```

The () operator or elem method allow the values of the matrix or array to be read or set. These methods take a single argument, which is of type octave_idx_type, that is the index into the matrix or array. Additionally, the matrix type allows two argument versions of the () operator and elem method, giving the row and column index of the value to get or set.

Note that these functions do significant error checking and so in some circumstances the user might prefer to access the data of the array or matrix directly through the fortran_vec method discussed below.
octave_idx_type numel (void) const
[Method] The total number of elements in the matrix or array.
size_t byte_size (void) const
[Method]
The number of bytes used to store the matrix or array.
dim_vector dims (void) const
[Method]
The dimensions of the matrix or array in value of type dim_vector.
int ndims (void) const [Method]
The number of dimensions of the matrix or array. Matrices are always 2-D, but arrays can be N -dimensional.
```

void resize (const dim_vector\&) [Method]

```
void resize (nrows, ncols) [Method]

A method taking either an argument of type dim_vector, or, in the case of a matrix, two arguments of type octave_idx_type defining the number of rows and columns in the matrix.

T* fortran_vec (void)
[Method]
This method returns a pointer to the underlying data of the matrix or array so that it can be manipulated directly, either within Octave or by an external library.

Operators such as,+- , or \(*\) can be used on the majority of the matrix and array types. In addition there are a number of methods that are of interest only for matrices such as transpose, hermitian, solve, etc.

The typical way to extract a matrix or array from the input arguments of DEFUN_DLD function is as follows
```

\#include <octave/oct.h>
DEFUN_DLD (addtwomatrices, args, , "Add A to B")
{
if (args.length () != 2)
print_usage ();
NDArray A = args(0).array_value ();
NDArray B = args(1).array_value ();
return octave_value (A + B);
}

```

To avoid segmentation faults causing Octave to abort, this function explicitly checks that there are sufficient arguments available before accessing these arguments. It then obtains two multi-dimensional arrays of type NDArray and adds these together. Note that the array_value method is called without using the is_matrix_type method. If an error occurs when attempting to extract the value, Octave will print a message and throw an exception. The reason to prefer this coding structure is that the arguments might be a type which is not an NDArray, but for which it would make sense to convert them to one. The array_value method allows this conversion to be performed transparently when possible. If you need to catch errors like this, and perform some kind of cleanup or other operation, you can catch the octave_execution_error exception.
\(A+B\), operating on two NDArray objects returns an NDArray, which is cast to an octave_ value on the return from the function. An example of the use of this demonstration function is
```

addtwomatrices (ones (2, 2), eye (2, 2))
\# 2 1
1 2

```

A list of the basic Matrix and Array types, the methods to extract these from an octave_value, and the associated header file is listed below.
Type
RowVector
ComplexRowVector
ColumnVector
ComplexColumnVector
Matrix
ComplexMatrix
boolMatrix
charMatrix
NDArray
ComplexNDArray
boolNDArray
charNDArray
int8NDArray
int16NDArray
int32NDArray
int64NDArray
uint8NDArray
uint16NDArray
uint32NDArray
uint64NDArray
Function
row_vector_value
complex_row_vector_value
column_vector_value
complex_column_vector_value
matrix_value
complex_matrix_value
bool_matrix_value
char_matrix_value
array_value
complex_array_value
bool_array_value
char_array_value
int8_array_value
int16_array_value
int32_array_value
int64_array_value
uint8_array_value
uint16_array_value
uint32_array_value
uint64_array_value

Source Code
dRowVector.h
CRowVector.h
dColVector.h
CColVector.h
dMatrix.h
CMatrix.h
boolMatrix.h
chMatrix.h
dNDArray.h
CNDArray.h
boolNDArray.h
charNDArray.h
int8NDArray.h
int16NDArray.h
int32NDArray.h
int64NDArray.h
uint8NDArray.h
uint16NDArray.h
uint32NDArray.h
uint64NDArray.h

\section*{A.1.3 Character Strings in Oct-Files}

A character string in Octave is just a special Array class. Consider the example:
```

\#include <octave/oct.h>
DEFUN_DLD (stringdemo, args, , "String Demo")
{
if (args.length () != 1)
print_usage ();
octave_value_list retval;
charMatrix ch = args(0).char_matrix_value ();
retval(1) = octave_value (ch, '\''); // Single Quote String
octave_idx_type nr = ch.rows ();
for (octave_idx_type i = 0; i < nr / 2; i++)

```
```

            {
                        std::string tmp = ch.row_as_string (i);
                        ch.insert (ch.row_as_string (nr-i-1).c_str (), i, 0);
            ch.insert (tmp.c_str (), nr-i-1, 0);
        }
    retval(0) = octave_value (ch, '"'); // Double Quote String
    return retval;
    }

```

An example of the use of this function is
```

s0 = ["First String"; "Second String"];
[s1,s2] = stringdemo (s0)

# s1 = Second String

    First String
    
# s2 = First String

    Second String
    typeinfo (s2)
=> sq_string
typeinfo (s1)

# string

```

One additional complication of strings in Octave is the difference between single quoted and double quoted strings. To find out if an octave_value contains a single or double quoted string use one of the predicate tests shown below.
```

if (args(0).is_sq_string ())
octave_stdout << "First argument is a single quoted string\n";
else if (args(0).is_dq_string ())
octave_stdout << "First argument is a double quoted string\n";

```

Note, however, that both types of strings are represented by the charNDArray type, and so when assigning to an octave_value, the type of string should be specified. For example:
```

octave_value_list retval;
charNDArray ch;
...
// Create single quoted string
retval(1) = octave_value (ch); // default constructor is sq_string
OR
retval(1) = octave_value (ch, '\''); // explicitly create sq_string
// Create a double quoted string
retval(0) = octave_value (ch, '"');

```

\section*{A.1.4 Cell Arrays in Oct-Files}

Octave's cell type is also available from within oct-files. A cell array is just an Array of octave_values, and thus each element of the cell array can be treated like any other octave_value. A simple example is
```

\#include <octave/oct.h>
\#include <octave/Cell.h>
DEFUN_DLD (celldemo, args, , "Cell Demo")
{
if (args.length () != 1)
print_usage ();
Cell c = args(0).cell_value ();
octave_value_list retval;
retval.resize (c.numel ()); // faster code by pre-declaring size
for (octave_idx_type i = 0; i < c.numel (); i++)
{
retval(i) = c(i); // using operator syntax
//retval(i) = c.elem (i); // using method syntax
}
return retval;
}

```

Note that cell arrays are used less often in standard oct-files and so the Cell.h header file must be explicitly included. The rest of the example extracts the octave_values one by one from the cell array and returns them as individual output arguments. For example:
```

[b1, b2, b3] = celldemo ({1, [1, 2], "test"})

# 

b1 = 1
b2 =
1 2
b3 = test

```

\section*{A.1.5 Structures in Oct-Files}

A structure in Octave is a map between a number of fields represented and their values. The Standard Template Library map class is used, with the pair consisting of a std::string and an Octave Cell variable.

A simple example demonstrating the use of structures within oct-files is
```

\#include <octave/oct.h>
\#include <octave/ov-struct.h>

```
```

DEFUN_DLD (structdemo, args, , "Struct Demo")
{
if (args.length () != 2)
print_usage ();
if (! args(0).isstruct ())
error ("structdemo: ARG1 must be a struct");
octave_scalar_map arg0 = args(0).scalar_map_value ();
//octave_map arg0 = args(0).map_value ();
if (! args(1).is_string ())
error ("structdemo: ARG2 must be a character string");
std::string arg1 = args(1).string_value ();
octave_value tmp = arg0.contents (arg1);
//octave_value tmp = arg0.contents (arg1)(0);
if (! tmp.is_defined ())
error ("structdemo: struct does not have a field named '%s'\n",
arg1.c_str ());
octave_scalar_map st;
st.assign ("selected", tmp);
return octave_value (st);
}

```

An example of its use is
```

x.a = 1; x.b = "test"; x.c = [1, 2];
structdemo (x, "b")
=> selected = test

```

The example above specifically uses the octave_scalar_map class which is for representing a single struct. For structure arrays, the octave_map class is used instead. The commented code shows how the demo could be modified to handle a structure array. In that case, the contents method returns a Cell which may have more than one element. Therefore, to obtain the underlying octave_value in the single struct example we would write
```

octave_value tmp = arg0.contents (arg1)(0);

```
where the trailing (0) is the () operator on the Cell object. If this were a true structure array with multiple elements we could iterate over the elements using the () operator.

Structures are a relatively complex data container and there are more functions available in oct-map.h which make coding with them easier than relying on just contents.

\section*{A.1.6 Sparse Matrices in Oct-Files}

There are three classes of sparse objects that are of interest to the user.
```

SparseMatrix
A double precision sparse matrix class
SparseComplexMatrix
A complex sparse matrix class
SparseBoolMatrix
A boolean sparse matrix class

```

All of these classes inherit from the Sparse<T> template class, and so all have similar capabilities and usage. The Sparse<T> class was based on Octave's Array<T> class and users familiar with Octave's Array classes will be comfortable with the use of the sparse classes.

The sparse classes will not be entirely described in this section, due to their similarity with the existing Array classes. However, there are a few differences due the nature of sparse objects, and these will be described. First, although it is fundamentally possible to have N-dimensional sparse objects, the Octave sparse classes do not allow them at this time; All instances of the sparse classes must be 2-dimensional. This means that SparseMatrix is actually more similar to Octave's Matrix class than it is to the NDArray class.

\section*{A.1.6.1 Array and Sparse Class Differences}

The number of elements in a sparse matrix is considered to be the number of nonzero elements, rather than the product of the dimensions. Therefore,
```

SparseMatrix sm;
int nnz = sm.nelem ();

```
returns the number of nonzero elements (like the interpreter function nnz). If the user really requires the number of elements in the matrix, including the nonzero elements, they should use numel rather than nelem. Note that for very large matrices, where the product of the two dimensions is larger than the representation of an unsigned int, numel can overflow. An example is speye (1e6) which will create a matrix with a million rows and columns, but only a million nonzero elements. In this case, the number of rows multiplied by the number of columns is more than two hundred times the maximum value that can be represented by an unsigned 32-bit int. The use of numel should, therefore, be avoided unless it is known that it will not overflow.

Extreme care is also required when using the elem method or the () operator which perform essentially the same function. The reason is that if a sparse object is non-const, then Octave will assume that a request for a zero element in a sparse matrix is in fact a request to create this element so it can be filled. Therefore, a piece of code like
```

SparseMatrix sm;
for (int j = 0; j < nc; j++)
for (int i = 0; i < nr; i++)
std::cerr << " (" << i << "," << j << "): " << sm(i,j) << "\n";

```
is a great way of turning a sparse matrix into a dense one, and a very slow way at that since it reallocates the sparse object for each zero element in the matrix.

A simple way of preventing the above from happening is to create a temporary constant version of the sparse matrix. Note that only the container for the sparse matrix will be copied, while the actual representation of the data will be shared between the two versions of the sparse matrix; This is not a costly operation. The example above, re-written to prevent sparse-to-dense conversion, is
```

SparseMatrix sm;
const SparseMatrix tmp (sm);
for (int j = 0; j < nc; j++)
for (int i = 0; i < nr; i++)
std::cerr << " (" << i << "," << j << "): " << tmp(i,j) << "\n";

```

Finally, because the sparse types aren't represented by a contiguous block of memory, the fortran_vec method of Array<T> is not available. It is, however, replaced by three separate methods ridx, cidx, and data, that access the raw compressed column format that Octave sparse matrices are stored in. These methods can be used in a manner similar to elem to allow the matrix to be accessed or filled. However, it is up to the user to respect the sparse matrix compressed column format or the matrix will become corrupted.

\section*{A.1.6.2 Creating Sparse Matrices in Oct-Files}

There are two useful strategies for creating a sparse matrix. The first is to create three vectors representing the row index, column index, and data values, and from these create the matrix. The second alternative is to create a sparse matrix with the appropriate amount of space, and then fill in the values. Both techniques have their advantages and disadvantages.

Below is an example of creating a small sparse matrix using the first technique
```

int nz, nr, nc;
nz = 4, nr = 3, nc = 4;
ColumnVector ridx (nz);
ColumnVector cidx (nz);
ColumnVector data (nz);
ridx(0) = 1; cidx(0) = 1; data(0) = 1;
ridx(1) = 2; cidx(1) = 2; data(1) = 2;
ridx(2) = 2; cidx(2) = 4; data(2) = 3;
ridx(3) = 3; cidx(3) = 4; data(3) = 4;
SparseMatrix sm (data, ridx, cidx, nr, nc);

```
which creates the matrix given in section Section 22.1.1 [Storage of Sparse Matrices], page 659. Note that the compressed matrix format is not used at the time of the creation of the matrix itself, but is used internally.

As discussed in the chapter on Sparse Matrices, the values of the sparse matrix are stored in increasing column-major ordering. Although the data passed by the user need not respect this requirement, pre-sorting the data will significantly speed up creation of the sparse matrix.

The disadvantage of this technique for creating a sparse matrix is that there is a brief time when two copies of the data exist. For extremely memory constrained problems this may not be the best technique for creating a sparse matrix.

The alternative is to first create a sparse matrix with the desired number of nonzero elements and then later fill those elements in. Sample code:
```

int nz, nr, nc;
nz = 4, nr = 3, nc = 4;
SparseMatrix sm (nr, nc, nz);
sm(0,0) = 1; sm(0,1) = 2; sm(1,3) = 3; sm(2,3) = 4;

```

This creates the same matrix as previously. Again, although not strictly necessary, it is significantly faster if the sparse matrix is created and the elements are added in columnmajor ordering. The reason for this is that when elements are inserted at the end of the current list of known elements then no element in the matrix needs to be moved to allow the new element to be inserted; Only the column indices need to be updated.

There are a few further points to note about this method of creating a sparse matrix. First, it is possible to create a sparse matrix with fewer elements than are actually inserted in the matrix. Therefore,
```

int nr, nc;
nr = 3, nc = 4;
SparseMatrix sm (nr, nc, 0);
sm(0,0) = 1; sm(0,1) = 2; sm(1,3) = 3; sm(2,3) = 4;

```
is perfectly valid. However, it is a very bad idea because as each new element is added to the sparse matrix the matrix needs to request more space and reallocate memory. This is an expensive operation that will significantly slow this means of creating a sparse matrix. It is possible to create a sparse matrix with excess storage, so having \(n z\) greater than 4 in this example is also valid. The disadvantage is that the matrix occupies more memory than strictly needed.

Of course, it is not always possible to know the number of nonzero elements prior to filling a matrix. For this reason the additional unused storage of a sparse matrix can be removed after its creation with the maybe_compress function. In addition to deallocating unused storage, maybe_compress can also remove zero elements from the matrix. The removal of zero elements from the matrix is controlled by setting the argument of the maybe_compress function to be true. However, the cost of removing the zeros is high because it implies re-sorting the elements. If possible, it is better for the user to avoid adding the unnecessary zeros in the first place. An example of the use of maybe_compress is
```

int nz, nr, nc;
nz = 6, nr = 3, nc = 4;
SparseMatrix sm1 (nr, nc, nz);
sm1(0,0) = 1; sm1 (0,1) = 2; sm1 (1,3) = 3; sm1(2,3) = 4;
sm1.maybe_compress (); // No zero elements were added
SparseMatrix sm2 (nr, nc, nz);
sm2(0,0) = 1; sm2 (0,1) = 2; sm(0,2) = 0; sm(1,2) = 0;
sm1(1,3) = 3; sm1 (2,3) = 4;
sm2.maybe_compress (true); // Zero elements were added

```

The use of the maybe_compress function should be avoided if possible as it will slow the creation of the matrix.

A third means of creating a sparse matrix is to work directly with the data in compressed row format. An example of this advanced technique might be
```

octave_value arg;
int nz, nr, nc;
nz = 6, nr = 3, nc = 4; // Assume we know the max \# nz
SparseMatrix sm (nr, nc, nz);
Matrix m = arg.matrix_value ();
int ii = 0;
sm.cidx (0) = 0;
for (int j = 1; j < nc; j++)
{
for (int i = 0; i < nr; i++)
{
double tmp = m(i,j);
if (tmp != 0.)
{
sm.data(ii) = tmp;
sm.ridx(ii) = i;
ii++;
}
}
sm.cidx(j+1) = ii;
}
sm.maybe_compress (); // If don't know a priori the final \# of nz.

```
which is probably the most efficient means of creating a sparse matrix.
Finally, it may sometimes arise that the amount of storage initially created is insufficient to completely store the sparse matrix. Therefore, the method change_capacity exists to reallocate the sparse memory. The above example would then be modified as
```

octave_value arg;

```
```

int nz, nr, nc;
nz = 6, nr = 3, nc = 4; // Guess the number of nz elements
SparseMatrix sm (nr, nc, nz);
Matrix m = arg.matrix_value ();
int ii = 0;
sm.cidx (0) = 0;
for (int j = 1; j < nc; j++)
{
for (int i = 0; i < nr; i++)
{
double tmp = m(i,j);
if (tmp != 0.)
{
if (ii == nz)
{
nz += 2; // Add 2 more elements
sm.change_capacity (nz);
}
sm.data(ii) = tmp;
sm.ridx(ii) = i;
ii++;
}
}
sm.cidx(j+1) = ii;
}
sm.maybe_compress (); // If don't know a priori the final \# of nz.

```

Note that both increasing and decreasing the number of nonzero elements in a sparse matrix is expensive as it involves memory reallocation. Also because parts of the matrix, though not its entirety, exist as old and new copies at the same time, additional memory is needed. Therefore, if possible avoid changing capacity.

\section*{A.1.6.3 Using Sparse Matrices in Oct-Files}

Most of the same operators and functions for sparse matrices that are available from the Octave interpreter are also available within oct-files. The basic means of extracting a sparse matrix from an octave_value, and returning it as an octave_value, can be seen in the following example.
```

octave_value_list retval;
SparseMatrix sm = args(0).sparse_matrix_value ();
SparseComplexMatrix scm = args(1).sparse_complex_matrix_value ();
SparseBoolMatrix sbm = args(2).sparse_bool_matrix_value ();
retval(2) = sbm;
retval(1) = scm;
retval(0) = sm;

```

The conversion to an octave_value is handled by the sparse octave_value constructors, and so no special care is needed.

\section*{A.1.7 Accessing Global Variables in Oct-Files}

Global variables allow variables in the global scope to be accessed. Global variables can be accessed within oct-files by using the support functions global_varval and global_assign from the current interpreter's symbol table. Both functions take as first argument a string representing the variable name to be obtained or assigned. The second argument of global_assign is the value to be assigned. An example of the use of these two functions is
```

\#include <octave/oct.h>
\#include <octave/interpreter.h>
DEFMETHOD_DLD (globaldemo, interp, args, , "Global Demo")
{
if (args.length () != 1)
print_usage ();
octave_value retval;
std::string s = args(0).string_value ();
octave::symbol_table\& symtab = interp.get_symbol_table ();
octave_value tmp = symtab.global_varval (s);
if (tmp.is_defined ())
retval = tmp;
else
retval = "Global variable not found";
symtab.global_assign ("a", 42.0);
return retval;
}

```

An example of its use is
```

global a b
b = 10;
globaldemo ("b")
=> 10
globaldemo ("c")
=> "Global variable not found"
num2str (a)

# 42

```

\section*{A.1.8 Calling Octave Functions from Oct-Files}

There is often a need to be able to call another Octave function from within an oct-file, and there are many examples of such within Octave itself. For example, the quad function is an oct-file that calculates the definite integral by quadrature over a user-supplied function.

There are also many ways in which a function could be given as input. It might be passed as one of
1. Function Handle
2. Anonymous Function Handle
3. String

The code below demonstrates all four methods of passing a function to an oct-file.
```

\#include <octave/oct.h>
\#include <octave/parse.h>
DEFMETHOD_DLD (funcdemo, interp, args, nargout, "Function Demo")
{
int nargin = args.length ();
if (nargin < 2)
print_usage ();
octave_value_list newargs;
for (octave_idx_type i = nargin - 1; i > 0; i--)
newargs(i-1) = args(i);
octave_value_list retval;
if (args(0).is_function_handle () || args(0).is_inline_function ()
|| args(0).is_string ())
retval = interp.feval (args(0), newargs, nargout);
else
error ("funcdemo: INPUT must be string, inline, or function handle");
return retval;
}

```

The first input to the demonstration code is a user-supplied function and the remaining arguments are all passed to the function.
```

funcdemo (@sin, 1)
=> 0.84147
funcdemo (@(x) sin (x), 1)
=> 0.84147
funcdemo ("sin", 1)

# 0.84147

funcdemo (@atan2, 1, 1)

# 0.78540

```

When the user function is passed as a string the treatment of the function is different. In some cases it is necessary to have the user supplied function as an octave_function object. In that case the string argument can be used to create a temporary function as demonstrated below.
```

std::octave fcn_name = unique_symbol_name ("__fcn__");
std::string fcode = "function y = ";
fcode.append (fcn_name);
fcode.append ("(x) y = ");
fcn = extract_function (args(0), "funcdemo", fcn_name,
fcode, "; endfunction");
if (fcn_name.length ())
clear_function (fcn_name);

```

There are two important things to know in this case. First, the number of input arguments to the user function is fixed, and in the above example is a single argument. Second, to avoid leaving the temporary function in the Octave symbol table it should be cleared after use. Also, by convention all internal function names begin and end with the character sequence '__'.

\section*{A.1.9 Calling External Code from Oct-Files}

Linking external C code to Octave is relatively simple, as the C functions can easily be called directly from \(\mathrm{C}++\). One possible issue is that the declarations of the external C functions may need to be explicitly defined as C functions to the compiler. If the declarations of the external C functions are in the header foo.h, then the tactic to ensure that the C++ compiler treats these declarations as C code is
```

\#ifdef __cplusplus
extern "C"
{
\#endif
\#include "foo.h"
\#ifdef __cplusplus
} /* end extern "C" */
\#endif

```

Calling Fortran code, however, can pose more difficulties. This is due to differences in the manner in which compilers treat the linking of Fortran code with C or C++ code. Octave supplies several macros that allow consistent behavior across a number of compilers.

The underlying Fortran code should use the XSTOPX function to replace the Fortran STOP function. XSTOPX uses the Octave exception handler to treat failing cases in the Fortran code explicitly. Note that Octave supplies its own replacement blas Xerbla function, which uses XSTOPX.

If the code calls XSTOPX, then the F77_XFCN macro should be used to call the underlying Fortran function. The Fortran exception state can then be checked with the global variable f77_exception_encountered. If XSTOPX will not be called, then the F77_FCN macro should be used instead to call the Fortran code.

There is no great harm in using F77_XFCN in all cases, except that for Fortran code that is short running and executes a large number of times, there is potentially an overhead in doing so. However, if F77_FCN is used with code that calls XSTOP, Octave can generate a segmentation fault.

An example of the inclusion of a Fortran function in an oct-file is given in the following example, where the C++ wrapper is
```

\#include <octave/oct.h>
\#include <octave/f77-fcn.h>
extern "C"
{
F77_RET_T
F77_FUNC (fortransub, FORTSUB)
(const F77_INT\&, F77_DBLE*, F77_CHAR_ARG_DECL F77_CHAR_ARG_LEN_DECL);
}
DEFUN_DLD (fortrandemo, args, , "Fortran Demo")
{
if (args.length () != 1)
print_usage ();
NDArray a = args(0).array_value ();
double *av = a.fortran_vec ();
octave_idx_type na = a.numel ();
OCTAVE_LOCAL_BUFFER (char, ctmp, 128);
F77_XFCN (fortransub, FORTSUB,
(na, av, ctmp F77_CHAR_ARG_LEN (128)));
return ovl (a, std::string (ctmp));
}
and the Fortran function is

```
```

subroutine fortransub (n, a, s)

```
subroutine fortransub (n, a, s)
implicit none
implicit none
character*(*) s
character*(*) s
real*8 a(*)
real*8 a(*)
integer*4 i, n, ioerr
integer*4 i, n, ioerr
do i = 1, n
do i = 1, n
    if (a(i) .eq. Od0) then
    if (a(i) .eq. Od0) then
        call xstopx ('fortransub: divide by zero')
        call xstopx ('fortransub: divide by zero')
    else
    else
        a(i) = 1d0 / a(i)
        a(i) = 1d0 / a(i)
    endif
    endif
enddo
```

enddo

```
```

write (unit = s, fmt = '(a,i3,a,a)', iostat = ioerr)
\$ 'There are ', n,
\$ ' values in the input vector', char(0)
if (ioerr .ne. 0) then
call xstopx ('fortransub: error writing string')
endif
return
end

```

This example demonstrates most of the features needed to link to an external Fortran function, including passing arrays and strings, as well as exception handling. Both the Fortran and C++ files need to be compiled in order for the example to work.
```

mkoctfile fortrandemo.cc fortransub.f
[b, s] = fortrandemo (1:3)
=>
b = 1.00000 0.50000 0.33333
s = There are 3 values in the input vector
[b, s] = fortrandemo (0:3)
error: fortrandemo: fortransub: divide by zero

```

\section*{A.1.10 Allocating Local Memory in Oct-Files}

Allocating memory within an oct-file might seem easy, as the C++ new/delete operators can be used. However, in that case great care must be taken to avoid memory leaks. The preferred manner in which to allocate memory for use locally is to use the OCTAVE_LOCAL_BUFFER macro. An example of its use is

OCTAVE_LOCAL_BUFFER (double, tmp, len)
that returns a pointer tmp of type double \(*\) of length len.
In this case, Octave itself will worry about reference counting and variable scope and will properly free memory without programmer intervention.

\section*{A.1.11 Input Parameter Checking in Oct-Files}

Because oct-files are compiled functions they open up the possibility of crashing Octave through careless function calls or memory faults. It is quite important that each and every function have a sufficient level of parameter checking to ensure that Octave behaves well.

The minimum requirement, as previously discussed, is to check the number of input arguments before using them to avoid referencing a nonexistent argument. However, in some cases this might not be sufficient as the underlying code imposes further constraints. For example, an external function call might be undefined if the input arguments are not integers, or if one of the arguments is zero, or if the input is complex and a real value was expected. Therefore, oct-files often need additional input parameter checking.

There are several functions within Octave that can be useful for the purposes of parameter checking. These include the methods of the octave_value class like is_real_matrix, is_numeric_type, etc. (see ov.h). Often, with a knowledge of the Octave m-file language, you can guess at what the corresponding C++ routine will. In addition there are some more specialized input validation functions of which a few are demonstrated below.
```

\#include <octave/oct.h>

```
```

DEFUN_DLD (paramdemo, args, nargout, "Parameter Check Demo")
{
if (args.length () != 1)
print_usage ();
NDArray m = args(0).array_value ();
double min_val = -10.0;
double max_val = 10.0;
octave_stdout << "Properties of input array:\n";
if (m.any_element_is_negative ())
octave_stdout << " includes negative values\n";
if (m.any_element_is_inf_or_nan ())
octave_stdout << " includes Inf or NaN values\n";
if (m.any_element_not_one_or_zero ())
octave_stdout << " includes other values than 1 and 0\n";
if (m.all_elements_are_int_or_inf_or_nan ())
octave_stdout << " includes only int, Inf or NaN values\n";
if (m.all_integers (min_val, max_val))
octave_stdout << " includes only integers in [-10,10]\n";
return octave_value_list ();
}

```

An example of its use is:
```

paramdemo ([1, 2, NaN, Inf])
P Properties of input array:
includes Inf or NaN values
includes other values than 1 and 0
includes only int, Inf or NaN values

```

\section*{A.1.12 Exception and Error Handling in Oct-Files}

Another important feature of Octave is its ability to react to the user typing Control-C during extended calculations. This ability is based on the C++ exception handler, where memory allocated by the C++ new/delete methods is automatically released when the exception is treated. When writing an oct-file which may run for a long time the programmer must periodically use the macro OCTAVE_QUIT, in order to allow Octave to check and possibly respond to a user typing Control-C. For example:
```

for (octave_idx_type i = 0; i < a.nelem (); i++)
{
OCTAVE_QUIT;
b.elem (i) = 2. * a.elem (i);
}

```

The presence of the OCTAVE_QUIT macro in the inner loop allows Octave to detect and acknowledge a Control-C key sequence. Without this macro, the user must either wait for the oct-file function to return before the interrupt is processed, or the user must press Control-C three times which will force Octave to exit completely.

The OCTAVE_QUIT macro does impose a very small performance penalty; For loops that are known to be small it may not make sense to include OCTAVE_QUIT.

When creating an oct-file that uses an external library, the function might spend a significant portion of its time in the external library. It is not generally possible to use the OCTAVE_QUIT macro in this case. The alternative code in this case is
```

BEGIN_INTERRUPT_IMMEDIATELY_IN_FOREIGN_CODE;
... some code that calls a "foreign" function ...
END_INTERRUPT_IMMEDIATELY_IN_FOREIGN_CODE;

```

The disadvantage of this is that if the foreign code allocates any memory internally, then this memory might be lost during an interrupt, without being deallocated. Therefore, ideally Octave itself should allocate any memory that is needed by the foreign code, with either the fortran_vec method or the OCTAVE_LOCAL_BUFFER macro.

The Octave unwind_protect mechanism (Section 10.8 [The unwind_protect Statement], page 188) can also be used in oct-files. In conjunction with the exception handling of Octave, it ensures that certain recovery code is always run even if an exception occurs. An example of the use of this mechanism is
```

\#include <octave/oct.h>
\#include <octave/unwind-prot.h>
void
my_err_handler (const char *fmt, ...)
{
// Do nothing!!
}
void
my_err_with_id_handler (const char *id, const char *fmt, ...)
{
// Do nothing!!
}
DEFUN_DLD (unwinddemo, args, nargout, "Unwind Demo")
{
if (args.length () < 2)
print_usage ();

```
```

    NDArray a = args(0).array_value ();
    NDArray b = args(1).array_value ();
    // Create unwind_action objects. At the end of the enclosing scope,
    // destructors for these objects will call the given functions with
    // the specified arguments.
    octave::unwind_action restore_warning_handler
    (set_liboctave_warning_handler, current_liboctave_warning_handler);
    octave::unwind_action restore_warning_with_id_handler
    (set_liboctave_warning_with_id_handler,
        current_liboctave_warning_with_id_handler);
    set_liboctave_warning_handler (my_err_handler);
    set_liboctave_warning_with_id_handler (my_err_with_id_handler);
    return octave_value (quotient (a, b));
    }

```

As can be seen in the example:
```

unwinddemo (1, 0)
| Inf
1 / 0

# warning: division by zero

    Inf
    ```

The warning for division by zero (and in fact all warnings) are disabled in the unwinddemo function.

\section*{A.1.13 Documentation and Testing of Oct-Files}

The documentation for an oct-file is contained in the fourth string parameter of the DEFUN_DLD macro. This string can be formatted in the same manner as the help strings for user functions, however there are some issues that are particular to the formatting of help strings within oct-files.

The major issue is that the help string will typically be longer than a single line of text, and so the formatting of long multi-line help strings needs to be taken into account. There are several possible solutions, but the most common is illustrated in the following example,
```

DEFUN_DLD (do_what_i_want, args, nargout,
"-*- texinfo -*-\n\
@deftypefn {} {} do_what_i_say (@var{n})\n\
A function that does what the user actually wants rather\n\
than what they requested.\n\
@end deftypefn")
{
}

```
where each line of text is terminated by \(\backslash \mathrm{n} \backslash\) which is an embedded newline in the string together with a C++ string continuation character. Note that the final \(\backslash\) must be the last character on the line.

Octave also includes the ability to embed test and demonstration code for a function within the code itself (see Appendix B [Test and Demo Functions], page 1061). This can be used from within oct-files (or in fact any file) with certain provisos. First, the test and demo functions of Octave look for \%! as the first two characters of a line to identify test and demonstration code. This is a requirement for oct-files as well. In addition, the test and demonstration code must be wrapped in a comment block to avoid it being interpreted by the compiler. Finally, the Octave test and demonstration code must have access to the original source code of the oct-file - not just the compiled code - as the tests are stripped from the compiled code. An example in an oct-file might be
```

/*
%!assert (sin ([1,2]), [sin(1),sin(2)])
%!error (sin ())
%!error (sin (1,1))
*/

```

\section*{A. 2 Mex-Files}

Octave includes an interface to allow legacy mex-files to be compiled and used with Octave. This interface can also be used to share compiled code between Octave and matlab users. However, as mex-files expose matlab's internal API, and the internal structure of Octave is different, a mex-file can never have the same performance in Octave as the equivalent octfile. In particular, to support the manner in which variables are passed to mex functions there are a significant number of additional copies of memory blocks when invoking or returning from a mex-file function. For this reason, it is recommended that any new code be written with the oct-file interface previously discussed.

\section*{A.2.1 Getting Started with Mex-Files}

The basic command to build a mex-file is either mkoctfile --mex or mex. The first command can be used either from within Octave or from the command line. To avoid issues with matlab's own mex command, the use of the command mex is limited to within Octave. Compiled mex-files have the extension .mex.
mex [-options] file . . .
status \(=\operatorname{mex}(. .\).
Compile source code written in C, C++, or Fortran, to a MEX file.
status is the return status of the mkoctfile function.
If the compilation fails, and the output argument is not requested, an error is raised. If the programmer requests status, however, Octave will merely issue a warning and it is the programmer's responsibility to verify the command was successful.
This is equivalent to mkoctfile --mex [-options] file.
See also: [mkoctfile], page 1010, [mexext], page 1034.
mexext ()
Return the filename extension used for MEX files.

See also: [mex], page 1034.
Consider the following short example:
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
mexPrintf ("Hello, World!\n");
mexPrintf ("I have %d inputs and %d outputs\n", nrhs, nlhs);
/* Return empty matrices for any outputs */
int i;
for (i = 0; i < nlhs; i++)
plhs[i] = mxCreateDoubleMatrix (0, 0, mxREAL);
}

```

The first line \#include "mex.h" makes available all of the definitions necessary for a mex-file. One important difference between Octave and matlab is that the header file "matrix.h" is implicitly included through the inclusion of "mex.h". This is necessary to avoid a conflict with the Octave file "Matrix.h" for operating systems and compilers that don't distinguish between filenames in upper and lower case.

The entry point into the mex-file is defined by mexFunction. The function takes four arguments:
1. The number of return arguments (\# of left-hand side args).
2. An array of pointers to return arguments.
3. The number of input arguments (\# of right-hand side args).
4. An array of pointers to input arguments.

Note that the function name definition is not explicitly included in mexFunction and so there can only be a single mexFunction entry point per file. Instead, the name of the function as seen in Octave is determined by the name of the mex-file itself minus the extension. If the above function is in the file myhello.c, it can be compiled with
```

mkoctfile --mex myhello.c

```
which creates a file myhello.mex. The function can then be run from Octave as
```

myhello (1,2,3)
=> Hello, World!
=> I have 3 inputs and 0 outputs

```

It should be noted that the mex-file contains no help string. To document mex-files, there should exist an m-file in the same directory as the mex-file itself. Taking the above as an example, there would need to be a file myhello.m which might contain the text
\%MYHELLO Simple test of the functionality of a mex-file.
In this case, the function that will be executed within Octave will be given by the mexfile, while the help string will come from the m-file. This can also be useful to allow a sample implementation of the mex-file within the Octave language itself for testing purposes.

Although there cannot be multiple entry points in a single mex-file, one can use the mexFunctionName function to determine what name the mex-file was called with. This can be used to alter the behavior of the mex-file based on the function name. For example, if
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
const char *nm;
nm = mexFunctionName ();
mexPrintf ("You called function: %s\n", nm);
if (strcmp (nm, "myfunc") == 0)
mexPrintf ("This is the principal function\n", nm);
return;
}

```
is in the file myfunc.c, and is compiled with
```

mkoctfile --mex myfunc.c
ln -s myfunc.mex myfunc2.mex

```
then as can be seen by
```

myfunc ()

# You called function: myfunc

    This is the principal function
    myfunc2 ()

# You called function: myfunc2

```
the behavior of the mex-file can be altered depending on the function's name.
Although the user should only include mex.h in their code, Octave declares additional functions, typedefs, etc., available to the user to write mex-files in the headers mexproto.h and mxarray.h.

\section*{A.2.2 Working with Matrices and Arrays in Mex-Files}

The basic mex type of all variables is mxArray. Any object, such as a matrix, cell array, or structure, is stored in this basic type. mxArray serves essentially the same purpose as the octave_value class in oct-files in that it acts as a container for all the more specialized types.

The mxArray structure contains at a minimum, the name of the variable it represents, its dimensions, its type, and whether the variable is real or complex. It can also contain a number of additional fields depending on the type of the mxArray. There are a number of functions to create mxArray structures, including mxCreateDoubleMatrix, mxCreateCellArray, mxCreateSparse, and the generic mxCreateNumericArray.

The basic function to access the data in an array is mxGetPr. Because the mex interface assumes that real and imaginary parts of a complex array are stored separately, there is an equivalent function mxGetPi that gets the imaginary part. Both of these functions
are only for use with double precision matrices. The generic functions mxGetData and mxGetImagData perform the same operation for all matrix types. For example:
```

mxArray *m;
mwSize *dims;
UINT32_T *pr;
dims = (mwSize *) mxMalloc (2 * sizeof (mwSize));
dims[0] = 2; dims[1] = 2;
m = mxCreateNumericArray (2, dims, mxUINT32_CLASS, mxREAL);
pr = (UINT32_T *) mxGetData (m);

```

There are also the functions mxSetPr, etc., that perform the inverse, and set the data of an array to use the block of memory pointed to by the argument of mxSetPr.

Note the type mwSize used above, and also mwIndex, are defined as the native precision of the indexing in Octave on the platform on which the mex-file is built. This allows both 32 - and 64 -bit platforms to support mex-files. mwSize is used to define array dimensions and the maximum number or elements, while mwIndex is used to define indexing into arrays.

An example that demonstrates how to work with arbitrary real or complex double precision arrays is given by the file mypow2.c shown below.
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
mwSize n;
mwIndex i;
double *vri, *vro;
if (nrhs != 1 || ! mxIsDouble (prhs[0]))
mexErrMsgTxt ("ARG1 must be a double matrix");
n = mxGetNumberOfElements (prhs[0]);
plhs[0] = mxCreateNumericArray (mxGetNumberOfDimensions (prhs[0]),
mxGetDimensions (prhs[0]),
mxGetClassID (prhs[0]),
mxIsComplex (prhs[0]));
vri = mxGetPr (prhs[0]);
vro = mxGetPr (plhs[0]);
if (mxIsComplex (prhs[0]))
{
double *vii, *vio;
vii = mxGetPi (prhs[0]);
vio = mxGetPi (plhs[0]);
for (i = 0; i < n; i++)

```
```

                        {
                        vro[i] = vri[i] * vri[i] - vii[i] * vii[i];
                        vio[i] = 2 * vri[i] * vii[i];
                }
        }
        else
        {
            for (i = 0; i < n; i++)
            vro[i] = vri[i] * vri[i];
        }
    }

```

An example of its use is
```

b = randn (4,1) + 1i * randn (4,1);
all (b. ^2 == mypow2 (b))
=> 1

```

The example above uses the functions mxGetDimensions, mxGetNumberOfElements, and mxGetNumberOfDimensions to work with the dimensions of multi-dimensional arrays. The functions mxGetM, and mxGetN are also available to find the number of rows and columns in a 2-D matrix (MxN matrix).

\section*{A.2.3 Character Strings in Mex-Files}

As mex-files do not make the distinction between single and double quoted strings that Octave does, there is perhaps less complexity in the use of strings and character matrices. An example of their use that parallels the demo in stringdemo.cc is given in the file mystring.c, as shown below.
```

\#include <string.h>
\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
mwSize m, n;
mwIndex i, j;
mxChar *pi, *po;
if (nrhs != 1 || ! mxIsChar (prhs[0])
|| mxGetNumberOfDimensions (prhs[0]) > 2)
mexErrMsgTxt ("ARG1 must be a char matrix");
m = mxGetM (prhs[0]);
n = mxGetN (prhs[0]);
pi = mxGetChars (prhs[0]);
plhs[0] = mxCreateNumericMatrix (m, n, mxCHAR_CLASS, mxREAL);
po = mxGetChars (plhs[0]);

```
```

    for (j = 0; j < n; j++)
    for (i = 0; i < m; i++)
        po[j*m + m - 1 - i] = pi[j*m + i];
    }

```

An example of its expected output is
```

mystring (["First String"; "Second String"])

# Second String

    First String
    ```

Other functions in the mex interface for handling character strings are mxCreateString, mxArrayToString, and mxCreateCharMatrixFromStrings. In a mex-file, a character string is considered to be a vector rather than a matrix. This is perhaps an arbitrary distinction as the data in the mxArray for the matrix is consecutive in any case.

\section*{A.2.4 Cell Arrays with Mex-Files}

One can perform exactly the same operations on Cell arrays in mex-files as in oct-files. An example that duplicates the function of the celldemo.cc oct-file in a mex-file is given by mycell.c as shown below.
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
mwSize n;
mwIndex i;
if (nrhs != 1 || ! mxIsCell (prhs[0]))
mexErrMsgTxt ("ARG1 must be a cell");
n = mxGetNumberOfElements (prhs[0]);
n = (n > nlhs ? nlhs : n);
for (i = 0; i < n; i++)
plhs[i] = mxDuplicateArray (mxGetCell (prhs[0], i));
}

```

The output is identical to the oct-file version as well.
```

[b1, b2, b3] = mycell ({1, [1, 2], "test"})

# 

b1 = 1
b2 =
1 2
b3 = test

```

Note in the example the use of the mxDuplicateArray function. This is needed as the mxArray pointer returned by mxGetCell might be deallocated. The inverse function to \(m x G e t C e l l\), used for setting Cell values, is mxSetCell and is defined as
```

void mxSetCell (mxArray *ptr, int idx, mxArray *val);

```

Finally, to create a cell array or matrix, the appropriate functions are
```

mxArray *mxCreateCellArray (int ndims, const int *dims);
mxArray *mxCreateCellMatrix (int m, int n);

```

\section*{A.2.5 Structures with Mex-Files}

The basic function to create a structure in a mex-file is mxCreateStructMatrix which creates a structure array with a two dimensional matrix, or mxCreateStructArray.
```

mxArray *mxCreateStructArray (int ndims, int *dims,
int num_keys,
const char **keys);
mxArray *mxCreateStructMatrix (int rows, int cols,
int num_keys,
const char **keys);

```

Accessing the fields of the structure can then be performed with mxGetField and mxSetField or alternatively with the mxGetFieldByNumber and mxSetFieldByNumber functions.
```

mxArray *mxGetField (const mxArray *ptr, mwIndex index,
const char *key);
mxArray *mxGetFieldByNumber (const mxArray *ptr,
mwIndex index, int key_num);
void mxSetField (mxArray *ptr, mwIndex index,
const char *key, mxArray *val);
void mxSetFieldByNumber (mxArray *ptr, mwIndex index,
int key_num, mxArray *val);

```

A difference between the oct-file interface to structures and the mex-file version is that the functions to operate on structures in mex-files directly include an index over the elements of the arrays of elements per field; Whereas, the oct-file structure includes a Cell Array per field of the structure.

An example that demonstrates the use of structures in a mex-file can be found in the file mystruct.c shown below.
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
int i;
mwIndex j;
mxArray *v;
const char *keys[] = { "this", "that" };

```
```

    if (nrhs != 1 || ! mxIsStruct (prhs[0]))
        mexErrMsgTxt ("ARG1 must be a struct");
    for (i = 0; i < mxGetNumberOfFields (prhs[0]); i++)
        for (j = 0; j < mxGetNumberOfElements (prhs[0]); j++)
            {
            mexPrintf ("field %s(%d) = ", mxGetFieldNameByNumber (prhs[0], i), j);
            v = mxGetFieldByNumber (prhs[0], j, i);
            mexCallMATLAB (0, NULL, 1, &v, "disp");
        }
    v = mxCreateStructMatrix (2, 2, 2, keys);
    mxSetFieldByNumber (v, 0, 0, mxCreateString ("this1"));
    mxSetFieldByNumber (v, 0, 1, mxCreateString ("that1"));
    mxSetFieldByNumber (v, 1, 0, mxCreateString ("this2"));
    mxSetFieldByNumber (v, 1, 1, mxCreateString ("that2"));
    mxSetFieldByNumber (v, 2, 0, mxCreateString ("this3"));
    mxSetFieldByNumber (v, 2, 1, mxCreateString ("that3"));
    mxSetFieldByNumber (v, 3, 0, mxCreateString ("this4"));
    mxSetFieldByNumber (v, 3, 1, mxCreateString ("that4"));
    if (nlhs)
    plhs[0] = v;
    }

```

An example of the behavior of this function within Octave is then
```

a(1).f1 = "f11"; a(1).f2 = "f12";
a(2).f1 = "f21"; a(2).f2 = "f22";
b = mystruct (a);

# field f1(0) = f11

    field f1(1) = f21
    field f2(0) = f12
    field f2(1) = f22
    b

# x2 struct array containing the fields:

    this
    that
    b(3)
| scalar structure containing the fields:
this = this3
that = that3

```

\section*{A.2.6 Sparse Matrices with Mex-Files}

The Octave format for sparse matrices is identical to the mex format in that it is a compressed column sparse format. Also, in both implementations sparse matrices are required to be two-dimensional. The only difference of importance to the programmer is that the real and imaginary parts of the matrix are stored separately.

The mex-file interface, in addition to using mxGetM, mxGetN, mxSetM, mxSetN, mxGetPr, \(m x G e t P i, m x S e t P r\), and mxSetPi, also supplies the following functions.
```

mwIndex *mxGetIr (const mxArray *ptr);
mwIndex *mxGetJc (const mxArray *ptr);
mwSize mxGetNzmax (const mxArray *ptr);
void mxSetIr (mxArray *ptr, mwIndex *ir);
void mxSetJc (mxArray *ptr, mwIndex *jc);
void mxSetNzmax (mxArray *ptr, mwSize nzmax);

```
mxGetNzmax gets the maximum number of elements that can be stored in the sparse matrix. This is not necessarily the number of nonzero elements in the sparse matrix. mxGetJc returns an array with one additional value than the number of columns in the sparse matrix. The difference between consecutive values of the array returned by mxGetJc define the number of nonzero elements in each column of the sparse matrix. Therefore,
```

mwSize nz, n;
mwIndex *Jc;
mxArray *m;
n = mxGetN (m);
Jc = mxGetJc (m);
nz = Jc[n];

```
returns the actual number of nonzero elements stored in the matrix in nz. As the arrays returned by \(m x G e t P r\) and \(m x G e t P i\) only contain the nonzero values of the matrix, we also need a pointer to the rows of the nonzero elements, and this is given by mxGetIr. A complete example of the use of sparse matrices in mex-files is given by the file mysparse.c shown below.
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
mwSize m, n, nz;
mxArray *v;
mwIndex i;
double *pr, *pi;
double *pr2, *pi2;
mwIndex *ir, *jc;
mwIndex *ir2, *jc2;

```
```

if (nrhs != 1 || ! mxIsSparse (prhs[0]))
mexErrMsgTxt ("ARG1 must be a sparse matrix");
m = mxGetM (prhs[0]);
n = mxGetN (prhs[0]);
nz = mxGetNzmax (prhs[0]);
if (mxIsComplex (prhs[0]))
{
mexPrintf ("Matrix is %d-by-%d complex sparse matrix", m, n);
mexPrintf (" with %d elements\n", nz);
pr = mxGetPr (prhs[0]);
pi = mxGetPi (prhs[0]);
ir = mxGetIr (prhs[0]);
jc = mxGetJc (prhs[0]);
i = n;
while (jc[i] == jc[i-1] \&\& i != 0) i--;
mexPrintf ("last nonzero element (%d, %d) = (%g, %g)\n",
ir[nz-1]+ 1, i, pr[nz-1], pi[nz-1]);
v = mxCreateSparse (m, n, nz, mxCOMPLEX);
pr2 = mxGetPr (v);
pi2 = mxGetPi (v);
ir2 = mxGetIr (v);
jc2 = mxGetJc (v);
for (i = 0; i < nz; i++)
{
pr2[i] = 2 * pr[i];
pi2[i] = 2 * pi[i];
ir2[i] = ir[i];
}
for (i = 0; i < n + 1; i++)
jc2[i] = jc[i];
if (nlhs > 0)
plhs[0] = v;
}
else if (mxIsLogical (prhs[0]))
{
mxLogical *pbr, *pbr2;
mexPrintf ("Matrix is %d-by-%d logical sparse matrix", m, n);
mexPrintf (" with %d elements\n", nz);

```
```

        pbr = mxGetLogicals (prhs[0]);
        ir = mxGetIr (prhs[0]);
        jc = mxGetJc (prhs[0]);
        i = n;
        while (jc[i] == jc[i-1] && i != 0) i--;
        mexPrintf ("last nonzero element (%d, %d) = %d\n",
                        ir[nz-1]+ 1, i, pbr[nz-1]);
        v = mxCreateSparseLogicalMatrix (m, n, nz);
        pbr2 = mxGetLogicals (v);
        ir2 = mxGetIr (v);
        jc2 = mxGetJc (v);
        for (i = 0; i < nz; i++)
            {
                pbr2[i] = pbr[i];
                ir2[i] = ir[i];
            }
        for (i = 0; i < n + 1; i++)
        jc2[i] = jc[i];
        if (nlhs > 0)
        plhs[0] = v;
    }
    else
{
mexPrintf ("Matrix is %d-by-%d real sparse matrix", m, n);
mexPrintf (" with %d elements\n", nz);
pr = mxGetPr (prhs[0]);
ir = mxGetIr (prhs[0]);
jc = mxGetJc (prhs[0]);
i = n;
while (jc[i] == jc[i-1] \&\& i != 0) i--;
mexPrintf ("last nonzero element (%d, %d) = %g\n",
ir[nz-1]+ 1, i, pr[nz-1]);
v = mxCreateSparse (m, n, nz, mxREAL);
pr2 = mxGetPr (v);
ir2 = mxGetIr (v);
jc2 = mxGetJc (v);
for (i = 0; i < nz; i++)
{
pr2[i] = 2 * pr[i];

```
```

                ir2[i] = ir[i];
            }
            for (i = 0; i < n + 1; i++)
            jc2[i] = jc[i];
            if (nlhs > 0)
            plhs[0] = v;
        }
    }

```

A sample usage of mysparse is
```

sm = sparse ([1, 0; 0, pi]);
mysparse (sm)

# 

Matrix is 2-by-2 real sparse matrix with 2 elements
last nonzero element (2, 2) = 3.14159

```

\section*{A.2.7 Calling Other Functions in Mex-Files}

It is possible to call other Octave functions from within a mex-file using mexCallmATLAB. An example of the use of mexCallmATLAB can be see in the example below.
```

\#include "mex.h"
void
mexFunction (int nlhs, mxArray *plhs[],
int nrhs, const mxArray *prhs[])
{
char *str;
mexPrintf ("Starting file myfeval.mex\n");
mexPrintf ("I have %d inputs and %d outputs\n", nrhs, nlhs);
if (nrhs < 1 || ! mxIsChar (prhs[0]))
mexErrMsgTxt ("ARG1 must be a function name");
str = mxArrayToString (prhs[0]);
mexPrintf ("I'm going to call the function %s\n", str);
if (nlhs == 0)
nlhs = 1; // Octave's automatic 'ans' variable
/* Cast prhs just to get rid of 'const' qualifier and stop compile warning */|
mexCallMATLAB (nlhs, plhs, nrhs-1, (mxArray**)prhs+1, str);
mxFree (str);
}

```

If this code is in the file myfeval.c, and is compiled to myfeval.mex, then an example of its use is
```

a = myfeval ("sin", 1)

# Starting file myfeval.mex

    I have 2 inputs and 1 outputs
    I'm going to call the interpreter function sin
    a = 0.84147
    ```

Note that it is not possible to use function handles within a mex-file.

\section*{A. 3 Standalone Programs}

The libraries Octave uses itself can be utilized in standalone applications. These applications then have access, for example, to the array and matrix classes, as well as to all of the Octave algorithms. The following C++ program, uses class Matrix from liboctave.a or liboctave.so.
```

\#include <iostream>
\#include <octave/oct.h>
int
main (void)
{
std::cout << "Hello Octave world!\n";
int n = 2;
Matrix a_matrix = Matrix (n, n);
for (octave_idx_type i = 0; i < n; i++)
for (octave_idx_type j = 0; j < n; j++)
a_matrix(i,j) = (i + 1) * 10 + (j + 1);
std::cout << a_matrix;
return 0;
}

```
mkoctfile can be used to build a standalone application with a command like
```

\$ mkoctfile --link-stand-alone standalone.cc -o standalone
\$ ./standalone
Hello Octave world!
1 1 1 2
2122
\$

```

Note that the application standalone will be dynamically linked against the Octave libraries and any Octave support libraries. The above allows the Octave math libraries to be used by an application. It does not, however, allow the script files, oct-files, or built-in functions of Octave to be used by the application. To do that, the Octave interpreter needs to be initialized first. An example of how to do this can then be seen in the code
```

\#include <iostream>
\#include <octave/oct.h>
\#include <octave/octave.h>
\#include <octave/parse.h>
\#include <octave/interpreter.h>
int
main (void)
{
// Create interpreter.
octave::interpreter interpreter;
try
{
// Inhibit reading history file by calling
//
// interpreter.initialize_history (false);
// Set custom load path here if you wish by calling
//
// interpreter.initialize_load_path (false);
// Perform final initialization of interpreter, including
// executing commands from startup files by calling
//
// interpreter.initialize ();
//
// if (! interpreter.initialized ())
// {
// std::cerr << "Octave interpreter initialization failed!"
// << std::endl;
// exit (status);
// }
//
// You may skip this step if you don't need to do anything
// between reading the startup files and telling the interpreter
// that you are ready to execute commands.
// Tell the interpreter that we're ready to execute commands:
int status = interpreter.execute ();
if (status != 0)
{
std::cerr << "creating embedded Octave interpreter failed!"
<< std::endl;

```
```

                    return status;
            }
                octave_idx_type n = 2;
                octave_value_list in;
                for (octave_idx_type i = 0; i < n; i++)
                    in(i) = octave_value (5 * (i + 2));
                octave_value_list out = octave::feval ("gcd", in, 1);
                if (out.length () > 0)
                        std::cout << "GCD of ["
                        << in(0).int_value ()
                        << ", "
                        << in(1).int_value ()
                        << "] is " << out(0).int_value ()
                        << std::endl;
            else
                        std::cout << "invalid\n";
        }
    catch (const octave::exit_exception& ex)
        {
            std::cerr << "Octave interpreter exited with status = "
                        << ex.exit_status () << std::endl;
        }
    catch (const octave::execution_exception&)
        {
            std::cerr << "error encountered in Octave evaluator!" << std::endl;
        }
    return 0;
    }

```
which, as before, is compiled and run as a standalone application with
```

\$ mkoctfile --link-stand-alone embedded.cc -o embedded
\$ ./embedded
GCD of [10, 15] is 5
\$

```

It is worth re-iterating that, if only built-in functions are to be called from a C++ standalone program then it does not need to initialize the interpreter. The general rule is that for a built-in function named function_name in the interpreter, there will be a C++ function named Ffunction_name (note the prepended capital F) accessible in the C++ API. The declarations for all built-in functions are collected in the header file builtin-defun-decls.h. This feature should be used with care as the list of built-in functions can change. No guarantees can be made that a function that is currently a built-in won't be implemented as a
.m file or as a dynamically linked function in the future. An example of how to call built-in functions from C++ can be seen in the code
```

\#include <iostream>
\#include <octave/oct.h>
\#include <octave/builtin-defun-decls.h>
int
main (void)
{
int n = 2;
Matrix a_matrix = Matrix (n, n);
for (octave_idx_type i = 0; i < n; i++)
for (octave_idx_type j = 0; j < n; j++)
a_matrix(i,j) = (i + 1) * 10 + (j + 1);
std::cout << "This is a matrix:" << std::endl
<< a_matrix << std::endl;
octave_value_list in;
in(0) = a_matrix;
octave_value_list out = Fnorm (in, 1);
double norm_of_the_matrix = out(0).double_value ();
std::cout << "This is the norm of the matrix:" << std::endl
<< norm_of_the_matrix << std::endl;
return 0;
}

```
which is compiled and run as a standalone application with
```

\$ mkoctfile --link-stand-alone standalonebuiltin.cc -o standalonebuiltin
\$ ./standalonebuiltin
This is a matrix:
1112
2122
This is the norm of the matrix:
34.4952
\$

```

\section*{A. 4 Java Interface}

The Java Interface is designed for calling Java functions from within Octave. If you want to do the reverse, and call Octave from within Java, try a library like joPas (http://jopas. sourceforge.net).

\section*{A.4.1 Making Java Classes Available}

Java finds classes by searching a classpath which is a list of Java archive files and/or directories containing class files. In Octave the classpath is composed of two parts:
- the static classpath is initialized once at startup of the JVM, and
- the dynamic classpath which can be modified at runtime.

Octave searches the static classpath first, and then the dynamic classpath. Classes appearing in the static classpath, as well as in the dynamic classpath, will therefore be found in the static classpath and loaded from this location. Classes which will be used frequently, or must be available to all users, should be added to the static classpath. The static classpath is populated once from the contents of a plain text file named javaclasspath.txt (or classpath.txt historically) when the Java Virtual Machine starts. This file contains one line for each individual classpath to be added to the static classpath. These lines can identify directories containing class files, or Java archives with complete class file hierarchies. Comment lines starting with a '\#' or a '\%' character are ignored.

The search rules for the file javaclasspath.txt (or classpath.txt) are:
- First, Octave tries to locate it in the current directory (where Octave was started from). If such a file is found, it is read and defines the initial static classpath. Thus, it is possible to define a static classpath on a 'per Octave invocation' basis.
- Next, Octave searches in the user's home directory. If a file javaclasspath.txt exists here, its contents are appended to the static classpath (if any). Thus, it is possible to build an initial static classpath on a 'per user' basis.
- Finally, Octave looks for a javaclasspath.txt in the m-file directory where Octave Java functions live. This is where the function javaclasspath.m resides, usually something like OCTAVE_HOME/share/octave/OCTAVE_VERSION/m/java/. You can find this directory by executing the command
```

which javaclasspath

```

If this file exists here, its contents are also appended to the static classpath. Note that the archives and class directories defined in this last step will affect all users.

Classes which are used only by a specific script should be placed in the dynamic classpath. This portion of the classpath can be modified at runtime using the javaaddpath and javarmpath functions.

Example:
```

octave> base_path = "C:/Octave/java_files";
octave> \# add two JAR archives to the dynamic classpath
octave> javaaddpath ([base_path, "/someclasses.jar"]);
octave> javaaddpath ([base_path, "/moreclasses.jar"]);
octave> \# check the dynamic classpath
octave> p = javaclasspath;
octave> disp (p{1});
C:/Octave/java_files/someclasses.jar
octave> disp (p{2});

```
```

C:/Octave/java_files/moreclasses.jar
octave> \# remove the first element from the classpath
octave> javarmpath ([base_path, "/someclasses.jar"]);
octave> p = javaclasspath;
octave> disp (p{1});
C:/Octave/java_files/moreclasses.jar
octave> \# provoke an error
octave> disp (p{2});
error: A(I): Index exceeds matrix dimension.

```

Another way to add files to the dynamic classpath exclusively for your user account is to use the file .octaverc which is stored in your home directory. All Octave commands in this file are executed each time you start a new instance of Octave. The following example adds the directory octave to Octave's search path and the archive myclasses.jar in this directory to the Java search path.
```


# contents of .octaverc:

addpath ("~/octave");
javaaddpath ("~/octave/myclasses.jar");

```

\section*{A.4.2 How to use Java from within Octave}

The function [javaObject], page 1054, creates Java objects. In fact it invokes the public constructor of the class with the given name and with the given parameters.

The following example shows how to invoke the constructors BigDecimal (double) and BigDecimal (String) of the builtin Java class java.math.BigDecimal.
```

javaObject ("java.math.BigDecimal", 1.001 );
javaObject ("java.math.BigDecimal", "1.001");

```

Note that parameters of the Octave type double are implicitly converted into the Java type double and the Octave type (array of) char is converted into the java type String. A Java object created by [javaObject], page 1054, is never automatically converted into an Octave type but remains a Java object. It can be assigned to an Octave variable.
```

a = 1.001;
b = javaObject ("java.math.BigDecimal", a);

```

Using [isjava], page 1055 , it is possible to check whether a variable is a Java object and its class can be determined as well. In addition to the previous example:
```

isjava (a)
ans = 0
class (a)
=> ans = double
isjava (b)
a ans = 1
class (b)
=> ans = java.math.BigDecimal

```

The example above can be carried out using only Java objects:
```

a = javaObject ("java.lang.Double", 1.001);
b = javaObject ("java.math.BigDecimal", a);
isjava (a)
a ans = 1
class (a)
=> ans = java.lang.Double
isjava (b)
m ans = 1
class (b)
=> ans = java.math.BigDecimal

```

One can see, that even a java.lang.Double is not converted to an Octave double, when created by [javaObject], page 1054. But ambiguities might arise, if the Java classes java.lang.Double or double are parameters of a method (or a constructor). In this case they can be converted into one another, depending on the context.

Via [javaObject], page 1054, one may create all kinds of Java objects but arrays. The latter are created through [javaArray], page 1054.

It is possible to invoke public member methods on Java objects in Java syntax:
```

a.toString
a ans = 1.001
b.toString
=> ans = 1.000999999999999889865...

```

The second result may be surprising, but simply comes from the fact, that 1.001 cannot exactly be represented as double, due to rounding. Note that unlike in Java, in Octave methods without arguments can be invoked with and without parentheses ().

Currently it is not possible to invoke static methods with a Java like syntax from within Octave. Instead, one has to use the function [javaMethod], page 1056, as in the following example:
```

java.math.BigDecimal.valueOf(1.001); \# does not work
javaMethod ("valueOf", "java.math.BigDecimal", 1.001); \# workaround

```

As mentioned before, method and constructor parameters are converted automatically between Octave and Java types, if appropriate. For functions this is also true with return values, whereas for constructors this is not.

It is also possible to access public fields of Java objects from within Octave using Java syntax, with the limitation of static fields:
```

java.math.BigDecimal.ONE; \# does not work
java_get ("java.math.BigDecimal", "ONE"); \# workaround

```

Accordingly, with [java_set], page 1056, the value of a field can be set. Note that only public Java fields are accessible from within Octave.

The following example indicates that in Octave empty brackets [] represent Java's null value and how Java exceptions are represented.
```

javaObject ("java.math.BigDecimal", []);

# error: [java] java.lang.NullPointerException

```

It is not recommended to represent Java's null value by empty brackets [], because null has no type whereas [] has type double.

In Octave it is possible to provide limited Java reflection by listing the public fields and methods of a Java object, both static or not.
```

fieldnames (<Java object>)
methods (<Java object>)

```

Finally, an examples is shown how to access the stack trace from within Octave, where the function [debug_java], page 1060, is used to set and to get the current debug state. In debug mode, the Java error and the stack trace are displayed.
```

debug_java (true) \# use "false" to omit display of stack trace
debug_java ()
m ans = 1
javaObject ("java.math.BigDecimal", "1") ...
.divide (javaObject ("java.math.BigDecimal", "0"))

```

\section*{A.4.3 Set up the JVM}

In order to execute Java code Octave creates a Java Virtual Machine (JVM). By default the version of the JVM is used that was detected during configuration on Unix-like systems or that is pointed to from the registry keys at HKEY_LOCAL_MACHINE \(\backslash\) SOFTWARE \(\backslash\) JavaSoft \(t \backslash J R E\) or HKEY_LOCAL_MACHINE\SOFTWARE\JavaSoft\Java Runtime Environment on Windows. The default path to the JVM can be overridden by setting the environment variable JAVA_HOME to the path where the JVM is installed. On Windows that might be, for example, C:\Program Files \(\backslash\) Java \(\backslash j r e-10.0 .2\). Make sure that you select a directory that contains the JVM with a bit-ness that matches Octave's.

The JVM is only loaded once per Octave session. Thus, to change the used version of the JVM, you might have to re-start Octave. To check which version of the JVM is currently being used, run version - java.

The JVM allocates a fixed amount of initial memory and may expand this pool up to a fixed maximum memory limit. The default values depend on the Java version (see [javamem], page 1059). The memory pool is shared by all Java objects running in the JVM. This strict memory limit is intended mainly to avoid runaway applications inside web browsers or in enterprise servers which can consume all memory and crash the system. When the maximum memory limit is hit, Java code will throw exceptions so that applications will fail or behave unexpectedly.

You can specify options for the creation of the JVM inside a file named java.opts. This is a text file where enter you enter lines containing -X and -D options that are then passed to the JVM during initialization.

The directory where the Java options file is located is specified by the environment variable OCTAVE_JAVA_DIR. If unset the directory where javaclasspath.m resides is used instead (typically OCTAVE_HOME/share/octave/OCTAVE_VERSION/m/java/). You can find this directory by executing
which javaclasspath
The - X options allow you to increase the maximum amount of memory available to the JVM. The following example allows up to 256 Megabytes to be used by adding the following line to the java.opts file:
-Xmx256m
The maximum possible amount of memory depends on your system. On a Windows system with 2 Gigabytes main memory you should be able to set this maximum to about 1 Gigabyte.

If your application requires a large amount of memory from the beginning, you can also specify the initial amount of memory allocated to the JVM. Adding the following line to the java.opts file starts the JVM with 64 Megabytes of initial memory:

\section*{-Xms64m}

For more details on the available - X options of your Java Virtual Machine issue the command 'java -X' at the operating system command prompt and consult the Java documentation.

The -D options can be used to define system properties which can then be used by Java classes inside Octave. System properties can be retrieved by using the getProperty() methods of the java.lang.System class. The following example line defines the property MyProperty and assigns it the string 12.34.
-DMyProperty=12.34
The value of this property can then be retrieved as a string by a Java object or in Octave:
```

octave> javaMethod ("getProperty", "java.lang.System", "MyProperty");
ans = 12.34

```

See also: javamem.

\section*{A.4.4 Java Interface Functions}

The following functions are the core of the Java Interface. They provide a way to create a Java object, get and set its data fields, and call Java methods which return results to Octave.
jobj = javaObject (classname)
jobj = javaObject (classname, arg1, ...)
Create a Java object of class classsname, by calling the class constructor with the arguments arg1,...
The first example below creates an uninitialized object, while the second example supplies an initial argument to the constructor.
```

x = javaObject ("java.lang.StringBuffer")
x = javaObject ("java.lang.StringBuffer", "Initial string")

```

See also: [javaMethod], page 1056, [javaArray], page 1054.
jary = javaArray (classname, sz)
jary = javaArray (classname, m, n, ...)
Create a Java array of size \(s z\) with elements of class classname.
classname may be a Java object representing a class or a string containing the fully qualified class name. The size of the object may also be specified with individual integer arguments \(m\), \(n\), etc.
The generated array is uninitialized. All elements are set to null if classname is a reference type, or to a default value (usually 0 ) if classname is a primitive type.

Sample code:
```

jary = javaArray ("java.lang.String", 2, 2);
jary(1,1) = "Hello";

```

See also: [javaObject], page 1054.
There are many different variable types in Octave, but only ones created through javaObject can use Java functions. Before using Java with an unknown object the type can be checked with isjava.
isjava (x)
Return true if \(x\) is a Java object.
See also: [class], page 41, [typeinfo], page 41, [isa], page 41, [javaObject], page 1054.
Once an object has been created it is natural to find out what fields the object has, and to read (get) and write (set) them.

In Octave the fieldnames function for structures has been overloaded to return the fields of a Java object. For example:
```

dobj = javaObject ("java.lang.Double", pi);
fieldnames (dobj)

# 

{
[1,1] = public static final double java.lang.Double.POSITIVE_INFINITY
[1,2] = public static final double java.lang.Double.NEGATIVE_INFINITY
[1,3] = public static final double java.lang.Double.NaN
[1,4] = public static final double java.lang.Double.MAX_VALUE
[1,5] = public static final double java.lang.Double.MIN_NORMAL
[1,6] = public static final double java.lang.Double.MIN_VALUE
[1,7] = public static final int java.lang.Double.MAX_EXPONENT
[1,8] = public static final int java.lang.Double.MIN_EXPONENT
[1,9] = public static final int java.lang.Double.SIZE
[1,10] = public static final java.lang.Class java.lang.Double.TYPE
}

```

The analogy of objects with structures is carried over into reading and writing object fields. To read a field the object is indexed with the '.' operator from structures. This is the preferred method for reading fields, but Octave also provides a function interface to read fields with java_get. An example of both styles is shown below.
```

dobj = javaObject ("java.lang.Double", pi);
dobj.MAX_VALUE
=> 1.7977e+308
java_get ("java.lang.Float", "MAX_VALUE")

# 3.4028e+38

val = java_get (obj, name)

```

Get the value of the field name of the Java object obj.
For static fields, obj can be a string representing the fully qualified name of the corresponding class.

When obj is a regular Java object, structure-like indexing can be used as a shortcut syntax. For instance, the following two statements are equivalent
```

java_get (x, "field1")
x.field1

```

See also: [java_set], page 1056, [javaMethod], page 1056, [javaObject], page 1054.
obj = java_set (obj, name, val)
Set the value of the field name of the Java object obj to val.
For static fields, obj can be a string representing the fully qualified named of the corresponding Java class.
When obj is a regular Java object, structure-like indexing can be used as a shortcut syntax. For instance, the following two statements are equivalent
```

java_set (x, "field1", val)
x.field1 = val

```

See also: [java_get], page 1055, [javaMethod], page 1056, [javaObject], page 1054.
To see what functions can be called with an object use methods. For example, using the previously created dobj:
```

methods (dobj)

# 

Methods for class java.lang.Double:
boolean equals(java.lang.Object)
java.lang.String toString(double)
java.lang.String toString()

```
-
To call a method of an object the same structure indexing operator '.' is used. Octave also provides a functional interface to calling the methods of an object through javaMethod. An example showing both styles is shown below.
```

dobj = javaObject ("java.lang.Double", pi);
dobj.equals (3)
O
javaMethod ("equals", dobj, pi)

# 1

ret = javaMethod (methodname, obj)
ret = javaMethod (methodname, obj, arg1, ...)

```

Invoke the method methodname on the Java object obj with the arguments arg1, ....
For static methods, obj can be a string representing the fully qualified name of the corresponding class.
When obj is a regular Java object, structure-like indexing can be used as a shortcut syntax. For instance, the two following statements are equivalent
```

ret = javaMethod ("method1", x, 1.0, "a string")
ret = x.method1 (1.0, "a string")

```
javaMethod returns the result of the method invocation.
See also: [methods], page 890, [javaObject], page 1054.

The following three functions are used to display and modify the class path used by the Java Virtual Machine. This is entirely separate from Octave's PATH variable and is used by the JVM to find the correct code to execute.
```

javaclasspath ()
dpath = javaclasspath ()
[dpath, spath] = javaclasspath ()
clspath = javaclasspath (what)

```

Return the class path of the Java virtual machine in the form of a cell array of strings.
If called with no inputs:
- If no output is requested, the dynamic and static classpaths are printed to the standard output.
- If one output value dpath is requested, the result is the dynamic classpath.
- If two output valuesdpath and spath are requested, the first variable will contain the dynamic classpath and the second will contain the static classpath.
If called with a single input parameter what:
"-dynamic"
Return the dynamic classpath.
"-static"
Return the static classpath.
"-all" Return both the static and dynamic classpath in a single cellstr.
See also: [javaaddpath], page 1057, [javarmpath], page 1057.
javaaddpath (clspath)
javaaddpath (clspath1,...)
javaaddpath (\{clspath1, ...\})
javaaddpath (..., "-end")
Add clspath to the beginning of the dynamic class path of the Java virtual machine.
clspath may either be a directory where .class files are found, or a \(\cdot \mathrm{jar}\) file containing Java classes. Multiple paths may be added at once by specifying additional arguments, or by using a cell array of strings.
If the final argument is "-end", append the new element to the end of the current classpath.
See also: [javarmpath], page 1057, [javaclasspath], page 1057.
javarmpath (clspath)
javarmpath (clspath1,...)
javarmpath (\{clspath1, ...\})
Remove clspath from the dynamic class path of the Java virtual machine.
clspath may either be a directory where .class files are found, or a .jar file containing Java classes. Multiple paths may be removed at once by specifying additional arguments, or by using a cell array of strings.
See also: [javaaddpath], page 1057, [javaclasspath], page 1057.

The following functions provide information and control over the interface between Octave and the Java Virtual Machine.
```

javachk (feature)
javachk (feature, caller)
msg = javachk (...)

```

Check for the presence of the Java feature in the current session. Return an error structure if feature is not available, not enabled, or not recognized.
Possible recognized features are:
"awt" Abstract Window Toolkit for GUIs.
"desktop"
Interactive desktop is running.
"jvm" Java Virtual Machine.
"swing" Swing components for lightweight GUIs.
If feature is not supported, a scalar struct with field "message" and "identifier" is returned. The field "message" contains an error message mentioning feature as well as the optional user-specified caller. This structure is suitable for passing in to the error function.

If feature is supported and available, an empty struct array is returned with fields "message" and "identifier".
javachk determines if specific Java features are available in an Octave session. This function is provided for scripts which may alter their behavior based on the availability of Java or specific Java runtime features.
Compatibility Note: The feature "desktop" is never available since Octave has no Java-based desktop.
See also: [usejava], page 1058, [error], page 239.
usejava (feature)
Return true if the Java element feature is available.
Possible features are:
"awt" Abstract Window Toolkit for GUIs.
"desktop"
Interactive desktop is running.
"jvm" Java Virtual Machine.
"swing" Swing components for lightweight GUIs.
usejava determines if specific Java features are available in an Octave session. This function is provided for scripts which may alter their behavior based on the availability of Java. The feature "desktop" always returns false as Octave has no Java-based desktop. Other features may be available if Octave was compiled with the Java Interface and Java is installed.
See also: [javachk], page 1058.
javamem ()
jmem = javamem ()
Show the current memory usage of the Java virtual machine (JVM) and run the garbage collector.
When no return argument is given the info is printed to the screen. Otherwise, the output cell array jmem contains Maximum, Total, and Free memory (in bytes).
All Java-based routines are run in the JVM's shared memory pool, a dedicated and separate part of memory claimed by the JVM from your computer's total memory (which comprises physical RAM and virtual memory / swap space on hard disk).
The maximum allowable memory usage can be configured using the file java.opts. The directory where this file resides is determined by the environment variable OCTAVE_JAVA_DIR. If unset, the directory where javaaddpath.m resides is used instead (typically OCTAVE_HOME/share/octave/OCTAVE_VERSION/m/java/).
java.opts is a plain text file with one option per line. The default initial memory size and default maximum memory size (which are both system dependent) can be overridden like so:
-Xms64m
-Xmx512m
(in megabytes in this example). You can adapt these values to your own requirements if your system has limited available physical memory or if you get Java memory errors.
"Total memory" is what the operating system has currently assigned to the JVM and depends on actual and active memory usage. "Free memory" is self-explanatory. During operation of Java-based Octave functions the amount of Total and Free memory will vary, due to Java's own cleaning up and your operating system's memory management.
val = java_matrix_autoconversion ()
old_val = java_matrix_autoconversion (new_val)
java_matrix_autoconversion (new_val, "local")
Query or set the internal variable that controls whether Java arrays are automatically converted to Octave matrices.
The default value is false.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [java_unsigned_autoconversion], page 1059, [debug_java], page 1060.
```

val = java_unsigned_autoconversion ()
old_val = java_unsigned_autoconversion (new_val)
java_unsigned_autoconversion (new_val, "local")
Query or set the internal variable that controls how integer classes are converted when java_matrix_autoconversion is enabled.

```

When enabled, Java arrays of class Byte or Integer are converted to matrices of class uint8 or uint32 respectively. The default value is true.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [java_matrix_autoconversion], page 1059, [debug_java], page 1060.
val = debug_java ()
old_val = debug_java (new_val)
debug_java (new_val, "local")
Query or set the internal variable that determines whether extra debugging information regarding the initialization of the JVM and any Java exceptions is printed.
When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.
See also: [java_matrix_autoconversion], page 1059, [java_unsigned_autoconversion], page 1059.

\section*{Appendix B Test and Demo Functions}

Octave includes a number of functions to allow the integration of testing and demonstration code in the source code of the functions themselves.

\section*{B. 1 Test Functions}
```

test name
test name quiet| normal| verbose
test ("name", "quiet| normal| verbose", fid)
test ("name", "quiet| normal| verbose", fname)
success = test (...)
[n, nmax, nxfail, nbug, nskip, nrtskip, nregression] = test (...)
[code, idx] = test ("name", "grabdemo")
test ([], "explain", fid)
test ([], "explain", fname)
Perform built-in self-tests from the first file in the loadpath matching name.

```
test can be called in either command or functional form. The exact operation of test is determined by a combination of mode (interactive or batch), reporting level ("quiet", "normal", "verbose"), and whether a logfile or summary output variable is used.
The default mode when test is called from the command line is interactive. In this mode, tests will be run until the first error is encountered, or all tests complete successfully. In batch mode, all tests are run regardless of any failures, and the results are collected for reporting. Tests which require user interaction, i.e., demo blocks, are never run in batch mode.
Batch mode is enabled by either 1) specifying a logfile using the third argument fname or fid, or 2) requesting an output argument such as success, \(n\), etc.
The optional second argument determines the amount of output to generate and which types of tests to run. The default value is "normal". Requesting an output argument will suppress printing the final summary message and any intermediate warnings, unless verbose reporting is enabled.
"quiet" Print a summary message when all tests pass, or print an error with the results of the first bad test when a failure occurs. Don't run tests which require user interaction.
"normal" Display warning messages about skipped tests or failing xtests during test execution. Print a summary message when all tests pass, or print an error with the results of the first bad test when a failure occurs. Don't run tests which require user interaction.
"verbose"
Display tests before execution. Print all warning messages. In interactive mode, run all tests including those which require user interaction.
The optional third input argument specifies a logfile where results of the tests should be written. The logfile may be a character string (fname) or an open file descriptor

ID (fid). To enable batch processing, but still print the results to the screen, use stdout for fid.
When called with just a single output argument success, test returns true if all of the tests were successful. If called with more than one output argument then the number of successful tests (n), the total number of tests in the file (nmax), the number of xtest failures (nxfail), the number of tests failed due known bugs (nbug), the number of tests skipped due to missing features (nskip), the number of tests skipped due to run-time conditions (nrtskip), and the number of regressions (nregression) are returned.

Example
```

test sind

# 

PASSES 5 out of 5 tests
[n, nmax] = test ("sind")

# 

n = 5
nmax = 5

```

\section*{Additional Calling Syntaxes}

If the second argument is the string "grabdemo", the contents of any built-in demo blocks are extracted but not executed. The text for all code blocks is concatenated and returned as code with idx being a vector of positions of the ends of each demo block. For an easier way to extract demo blocks from files, See [example], page 1070.
If the second argument is "explain" then name is ignored and an explanation of the line markers used in test output reports is written to the file specified by fname or fid.

See also: [assert], page 1067, [fail], page 1068, [demo], page 1069, [example], page 1070, [error], page 239.
test scans the named script file looking for lines which start with the identifier '\%!'. The prefix is stripped off and the rest of the line is processed through the Octave interpreter. If the code generates an error, then the test is said to fail.

Since eval() will stop at the first error it encounters, you must divide your tests up into blocks, with anything in a separate block evaluated separately. Blocks are introduced by valid keywords like test, function, or assert immediately following '\%!'. A block is defined by indentation as in Python. Lines beginning with '\%!<whitespace>' are part of the preceding block.

For example:
```

%!test error ("this test fails!")
%!test "test doesn't fail. it doesn't generate an error"

```

When a test fails, you will see something like:
```

    ***** test error ("this test fails!")
    !!!!! test failed
this test fails!

```

Generally, to test if something works, you want to assert that it produces a correct value. A real test might look something like
```

%!test
%! a = [1, 2, 3; 4, 5, 6]; B = [1; 2];
%! expect = [ a ; 2*a ];
%! get = kron (b, a);
%! if (any (size (expect) != size (get)))
%! error ("wrong size: expected %d,%d but got %d,%d",
%! size (expect), size (get));
%! elseif (any (any (expect != get)))
%! error ("didn't get what was expected.");
%! endif

```

To make the process easier, use the assert function. For example, with assert the previous test is reduced to:
```

%!test
%! a = [1, 2, 3; 4, 5, 6]; b = [1; 2];
%! assert (kron (b, a), [ a; 2*a ]);

```
assert can accept a tolerance so that you can compare results absolutely or relatively. For example, the following all succeed:
```

%!test assert (1+eps, 1, 2*eps) \# absolute error
%!test assert (100+100*eps, 100, -2*eps) \# relative error

```

You can also do the comparison yourself, but still have assert generate the error:
```

%!test assert (isempty ([]))
%!test assert ([1, 2; 3, 4] > 0)

```

Because assert is so frequently used alone in a test block, there is a shorthand form:
```

%!assert (...)

```
which is equivalent to:
```

%!test assert (...)

```

Occasionally a block of tests will depend on having optional functionality in Octave. Before testing such blocks the availability of the required functionality must be checked. A \%! testif HAVE_XXX block will only be run if Octave was compiled with functionality 'HAVE_XXX'. For example, the sparse single value decomposition, svds (), depends on having the arpack library. All of the tests for svds begin with
\%!testif HAVE_ARPACK
Review config.h or __octave_config_info__ ("build_features") to see some of the possible values to check.

Sometimes during development there is a test that should work but is known to fail. You still want to leave the test in because when the final code is ready the test should pass, but you may not be able to fix it immediately. To avoid unnecessary bug reports for these known failures, mark the block with xtest rather than test:
```

%!xtest assert (1==0)
%!xtest fail ("success=1", "error")

```

In this case, the test will run and any failure will be reported. However, testing is not aborted and subsequent test blocks will be processed normally. Another use of xtest is for statistical tests which should pass most of the time but are known to fail occasionally.

Each block is evaluated in its own function environment, which means that variables defined in one block are not automatically shared with other blocks. If you do want to share variables, then you must declare them as shared before you use them. For example, the following declares the variable a, gives it an initial value (default is empty), and then uses it in several subsequent tests.
```

%!shared a
%! a = [1, 2, 3; 4, 5, 6];
%!assert (kron ([1; 2], a), [ a; 2*a ])
%!assert (kron ([1, 2], a), [ a, 2*a ])
%!assert (kron ([1,2; 3,4], a), [ a,2*a; 3*a,4*a ])

```

You can share several variables at the same time:
```

%!shared a, b

```

You can also share test functions:
```

%!function a = fn (b)
%! a = 2*b;
%!endfunction
%!assert (fn(2), 4)

```

Note that all previous variables and values are lost when a new shared block is declared.
Remember that \%!function begins a new block and that \%!endfunction ends this block. Be aware that until a new block is started, lines starting with ‘\%! <space>' will be discarded as comments. The following is nearly identical to the example above, but does nothing.
```

%!function a = fn (b)
%! a = 2*b;
%!endfunction
%! assert (fn(2), 4)

```

Because there is a space after ' \(\%\) !' the assert statement does not begin a new block and this line is treated as a comment.

Error and warning blocks are like test blocks, but they only succeed if the code generates an error. You can check the text of the error is correct using an optional regular expression <pattern>. For example:
```

%!error <passes!> error ("this test passes!")

```

If the code doesn't generate an error, the test fails. For example:
```

%!error "this is an error because it succeeds."

```
produces
```

    ***** error "this is an error because it succeeds."
    !!!!! test failed: no error

```

It is important to automate the tests as much as possible, however some tests require user interaction. These can be isolated into demo blocks, which if you are in batch mode,
are only run when called with demo or the verbose option to test. The code is displayed before it is executed. For example,
```

%!demo
%! t = [0:0.01:2*pi]; x = sin (t);
%! plot (t, x);
%! \# you should now see a sine wave in your figure window

```
produces
```

funcname example 1:
t = [0:0.01:2*pi]; x = sin (t);
plot (t, x);
\# you should now see a sine wave in your figure window
Press <enter> to continue:

```

Note that demo blocks cannot use any shared variables. This is so that they can be executed by themselves, ignoring all other tests.

If you want to temporarily disable a test block, put \# in place of the block type. This creates a comment block which is echoed in the log file but not executed. For example:
```

%!\#demo
%! t = [0:0.01:2*pi]; x = sin (t);
%! plot (t, x);
%! \# you should now see a sine wave in your figure window

```

The following trivial code snippet provides examples for the use of fail, assert, error, and xtest:
```

function output = must_be_zero (input)
if (input != 0)
error ("Nonzero input!")
endif
output = input;
endfunction
%!fail ("must_be_zero (1)")
%!assert (must_be_zero (0), 0)
%!error <Nonzero> must_be_zero (1)
%!xtest error ("This code generates an error")

```

When putting this in a file must_be_zero.m, and running the test, we see
```

test must_be_zero verbose
$\Rightarrow$
>>>>> /path/to/must_be_zero.m
***** fail ("must_be_zero (1)")
***** assert (must_be_zero (0), 0)
***** error <Nonzero> must_be_zero (1)
***** xtest error ("This code generates an error")
!!!!! known failure
This code generates an error
PASSES 3 out of 4 tests (1 expected failure)

```

\section*{Block type summary:}
```

%!test
%!test <MESSAGE> interpreted as for xtest.

```
```

%!testif HAVE_XXX

```
%!testif HAVE_XXX
%!testif HAVE_XXX, HAVE_YYY, ...
%!testif HAVE_XXX, HAVE_YYY, ...
%!testif HAVE_XXX, HAVE_YYY ...; RUNTIME_COND
%!testif HAVE_XXX, HAVE_YYY ...; RUNTIME_COND
%!testif . .. <MESSAGE>
```

%!testif . .. <MESSAGE>

```

Check that entire block is correct. If <MESSAGE> is present, the test block is

Check block only if Octave was compiled with feature HAVE_XXX. RUNTIME_COND is an optional expression to evaluate to check whether some condition is met when the test is executed. If RUNTIME_COND is false, the test is skipped. If <MESSAGE> is present, the test block is interpreted as for xtest.
```

%!xtest
%!xtest <MESSAGE>

```

Check block, report a test failure but do not abort testing. If <MESSAGE> is present, then the text of the message is displayed if the test fails, like this:

\section*{!!!!! known bug: MESSAGE}

If the message is an integer, it is interpreted as a bug ID for the Octave bug tracker and reported as
```

!!!!! known bug: https://octave.org/testfailure/?BUG-ID

```
in which BUG-ID is the integer bug number. The intent is to allow clearer documentation of known problems.

If MESSAGE is an integer preceeded by an asterisk (e.g., *12345), it is interpreted as the id for a bug report that has been closed. That usually means that the issue probed in this test has been resolved. If such tests are failing, they are reported as regressions by the test function:

\footnotetext{
!!!!! regression: https://octave.org/testfailure/?BUG-ID
}
```

%!error
%!error <MESSAGE>
%!warning
%!warning <MESSAGE>

```

Check for correct error or warning message. If <MESSAGE> is supplied it is interpreted as a regular expression pattern that is expected to match the error or warning message.
\%!demo Demo only executes in interactive mode.
\%! \# Comment. Ignore everything within the block
\%! shared \(x, y, z\)
Declare variables for use in multiple tests.
\%!function
Define a function for use in multiple tests.
\%!endfunction
Close a function definition.
\(\%!\) assert ( \(\mathrm{x}, \mathrm{y}, \mathrm{tol}\) )
\%!assert <MESSAGE> ( \(\mathrm{x}, \mathrm{y}, \mathrm{tol}\) )
\(\%\) !fail (CODE, PATTERN)
\%!fail <MESSAGE> (CODE, PATTERN)
Shorthand for \%! test assert ( \(x, y\), tol) or \%!test fail (CODE, PATTERN).
If <MESSAGE> is present, the test block is interpreted as for xtest.
When coding tests the Octave convention is that lines that begin with a block type do not have a semicolon at the end. Any code that is within a block, however, is normal Octave code and usually will have a trailing semicolon. For example,
\#\# bare block instantiation
\(\%\) !assert ( \(\sin (0), 0)\)
but
\#\# test block with normal Octave code
\(\%\) ! test
\(\%\) ! assert ( \(\sin (0), 0)\);
You can also create test scripts for built-in functions and your own C++ functions. To do so, put a file with the bare function name (no .m extension) in a directory in the load path and it will be discovered by the test function. Alternatively, you can embed tests directly in your C++ code:
```

/*
%!test disp ("this is a test")
*/
\#if 0
%!test disp ("this is a test")
\#endif

```
or

However, in this case the raw source code will need to be on the load path and the user will have to remember to type test ("funcname.cc").
```

assert (cond)
assert (cond, errmsg)
assert (cond, errmsg,...)
assert (cond, msg_id, errmsg, ...)
assert (observed, expected)
assert (observed, expected, tol)

```

Produce an error if the specified condition is not met.
assert can be called in three different ways.
assert (cond)
assert (cond, errmsg)
assert (cond, errmsg, ...)
assert (cond, msg_id, errmsg, ...)
Called with a single argument cond, assert produces an error if cond is false (numeric zero).
Any additional arguments are passed to the error function for processing.
assert (observed, expected)
Produce an error if observed is not the same as expected.
Note that observed and expected can be scalars, vectors, matrices, strings, cell arrays, or structures.
assert (observed, expected, tol)
Produce an error if observed is not the same as expected but equality comparison for numeric data uses a tolerance tol.

If tol is positive then it is an absolute tolerance which will produce an error if abs (observed - expected) > abs (tol).
If tol is negative then it is a relative tolerance which will produce an error if abs (observed - expected) > abs (tol * expected).
If expected is zero tol will always be interpreted as an absolute tolerance.
If tol is not scalar its dimensions must agree with those of observed and expected and tests are performed on an element-by-element basis.

See also: [fail], page 1068, [test], page 1061, [error], page 239, [isequal], page 165.
fail (code)
fail (code, pattern)
fail (code, "warning")
fail (code, "warning", pattern)
Return true if code fails with an error message matching pattern, otherwise produce an error.
code must be in the form of a string that is passed to the Octave interpreter via the evalin function, i.e., a (quoted) string constant or a string variable.
Note that if code runs successfully, rather than failing, the error printed is:

If called with two arguments, the return value will be true only if code fails with an error message containing pattern (case sensitive). If the code fails with a different error than the one specified in pattern then the message produced is:
```

expected <pattern>
but got <text of actual error>

```

The angle brackets are not part of the output.
When called with the "warning" option fail will produce an error if executing the code produces no warning.
See also: [assert], page 1067, [error], page 239.

\section*{B. 2 Demonstration Functions}
```

demo name
demo name n
demo ("name")
demo ("name", n)
Run example code block $n$ associated with the function name.
If $n$ is not specified, all examples are run.

```

The preferred location for example code blocks is embedded within the script m-file immediately following the code that it exercises. Alternatively, the examples may be stored in a file with the same name but no extension located on Octave's load path. To separate examples from regular script code all lines are prefixed by \%!. Each example must also be introduced by the keyword "demo" flush left to the prefix with no intervening spaces. The remainder of the example can contain arbitrary Octave code. For example:
```

%!demo
%! t = 0:0.01:2*pi;
%! x = sin (t);
%! plot (t, x);
%! title ("one cycle of a sine wave");
%! \#--------------------------------------------------------
%! \# the figure window shows one cycle of a sine wave

```

Note that the code is displayed before it is executed so that a simple comment at the end suffices for labeling what is being shown. For plots, labeling can also be done with title or text. It is generally not necessary to use disp or printf within the demo.
Demos are run in a stand-alone function environment with no access to external variables. This means that every demo must have separate initialization code. Alternatively, all demos can be combined into a single large demo with the code
\%! input ("Press <enter> to continue: ", "s");
between the sections, but this usage is discouraged. Other techniques to avoid multiple initialization blocks include using multiple plots with a new figure command between each plot, or using subplot to put multiple plots in the same window.

Finally, because demo evaluates within a function context it is not possible to define new functions within the code. Anonymous functions make a good substitute in most instances. If function blocks must be used then the code eval (example ("function", n)) will allow Octave to see them. This has its own problems, however, as eval only evaluates one line or statement at a time. In this case the function declaration must be wrapped with "if 1 <demo stuff> endif" where "if" is on the same line as "demo". For example:
```

%!demo if 1
%! function y = f(x)
%! y = x;
%! endfunction
%! f(3)
%! endif

```

See also: [rundemos], page 1070, [example], page 1070, [test], page 1061.
```

example name

```
example name \(n\)
example ("name")
example ("name", n)
[s, idx] = example (...)

Display the code for example \(n\) associated with the function name, but do not run it.
If \(n\) is not specified, all examples are displayed.
When called with output arguments, the examples are returned in the form of a string \(s\), with idx indicating the ending position of the various examples.
For a complete explanation see [demo], page 1069.
See also: [demo], page 1069, [test], page 1061.
oruntests ()
oruntests (directory)
Execute built-in tests for all m-files in the specified directory.
Test blocks in any C++ source files (*.cc) will also be executed for use with dynamically linked oct-file functions.
If no directory is specified, operate on all directories in Octave's search path for functions.

See also: [rundemos], page 1070, [test], page 1061, [path], page 214.
rundemos ()
rundemos (directory)
Execute built-in demos for all m-files in the specified directory.
Demo blocks in any C++ source files (*.cc) will also be executed for use with dynamically linked oct-file functions.
If no directory is specified, operate on all directories in Octave's search path for functions.
See also: [demo], page 1069, [oruntests], page 1070, [path], page 214.
speed (f, init, max_n, f2, tol)
[order, \(\left.n, T_{-} f, T_{-} f 2\right]=\operatorname{speed}(. .\).
Determine the execution time of an expression ( \(f\) ) for various input values ( \(n\) ).
The \(n\) are log-spaced from 1 to max_n. For each \(n\), an initialization expression (init) is computed to create any data needed for the test. If a second expression (f2) is given then the execution times of the two expressions are compared. When called without output arguments the results are printed to stdout and displayed graphically.
\(f \quad\) The code expression to evaluate.
max_n The maximum test length to run. The default value is 100. Alternatively, use [min_n, max_n] or specify the \(n\) exactly with [ \(n 1, n 2, \ldots, n k\) ].
init Initialization expression for function argument values. Use \(k\) for the test number and \(n\) for the size of the test. This should compute values for all variables used by \(f\). Note that init will be evaluated first for \(k=0\), so things which are constant throughout the test series can be computed once. The default value is \(x=\operatorname{randn}(n, 1)\).
f2 An alternative expression to evaluate, so that the speed of two expressions can be directly compared. The default is [].
tol Tolerance used to compare the results of expression \(f\) and expression \(f 2\). If tol is positive, the tolerance is an absolute one. If tol is negative, the tolerance is a relative one. The default is eps. If tol is Inf, then no comparison will be made.
order The time complexity of the expression \(O\left(a * n^{p}\right)\). This is a structure with fields a and p .
\(n \quad\) The values \(n\) for which the expression was calculated AND the execution time was greater than zero.
T_f The nonzero execution times recorded for the expression \(f\) in seconds.
T_f2 The nonzero execution times recorded for the expression \(f 2\) in seconds. If required, the mean time ratio is simply mean (T_f ./ T_f2).
The slope of the execution time graph shows the approximate power of the asymptotic running time \(O\left(n^{p}\right)\). This power is plotted for the region over which it is approximated (the latter half of the graph). The estimated power is not very accurate, but should be sufficient to determine the general order of an algorithm. It should indicate if, for example, the implementation is unexpectedly \(O\left(n^{2}\right)\) rather than \(O(n)\) because it extends a vector each time through the loop rather than pre-allocating storage. In the current version of Octave, the following is not the expected \(O(n)\).
```

speed ("for i = 1:n, y{i} = x(i); endfor", "", [1000, 10000])

```

But it is if you preallocate the cell array y:
```

speed ("for i = 1:n, y{i} = x(i); endfor", ...
"x = rand (n, 1); y = cell (size (x));", [1000, 10000])

```

An attempt is made to approximate the cost of individual operations, but it is wildly inaccurate. You can improve the stability somewhat by doing more work for each n . For example:
```

speed ("airy(x)", "x = rand (n, 10)", [10000, 100000])

```

When comparing two different expressions ( \(f, f 2\) ), the slope of the line on the speedup ratio graph should be larger than 1 if the new expression is faster. Better algorithms have a shallow slope. Generally, vectorizing an algorithm will not change the slope of the execution time graph, but will shift it relative to the original. For example:
```

speed ("sum (x)", "", [10000, 100000], ...
"v = 0; for i = 1:length (x), v += x(i); endfor")

```

The following is a more complex example. If there was an original version of xcorr using for loops and a second version using an FFT, then one could compare the run speed for various lags as follows, or for a fixed lag with varying vector lengths as follows:
```

speed ("xcorr (x, n)", "x = rand (128, 1);", 100,
"xcorr_orig (x, n)", -100*eps)
speed ("xcorr (x, 15)", "x = rand (20+n, 1);", 100,
"xcorr_orig (x, n)", -100*eps)

```

Assuming one of the two versions is in xcorr_orig, this would compare their speed and their output values. Note that the FFT version is not exact, so one must specify an acceptable tolerance on the comparison \(100 *\) eps. In this case, the comparison should be computed relatively, as abs \(((x-y) . / y)\) rather than absolutely as abs \((x-y)\). Type example ("speed") to see some real examples or demo ("speed") to run them.

\section*{Appendix C Obsolete Functions}

After being marked as deprecated for two major releases, the following functions have been removed from Octave. The third column of the table shows the version of Octave in which the function was removed. Prior to removal, each function in the list was marked as deprecated for at least two major releases. All deprecated functions issue warnings explaining that they will be removed in a future version of Octave, and which function should be used instead.

Replacement functions do not always accept precisely the same arguments as the obsolete function, but should provide equivalent functionality.
\begin{tabular}{|c|c|c|}
\hline Obsolete Function beta_cdf & Replacement betacdf in Octave Forge statistics pkg & Version
3.4.0 \\
\hline beta_inv & betainv in Octave Forge statistics pkg & 3.4.0 \\
\hline beta_pdf & betapdf in Octave Forge statistics pkg & 3.4.0 \\
\hline beta_rnd & betarnd in Octave Forge statistics pkg & 3.4.0 \\
\hline binomial_cdf & binocdf in Octave Forge statistics pkg & 3.4.0 \\
\hline binomial_inv & binoinv in Octave Forge statistics pkg & 3.4.0 \\
\hline binomial_pdf & binopdf in Octave Forge statistics pkg & 3.4.0 \\
\hline binomial_rnd & binornd in Octave Forge statistics pkg & 3.4.0 \\
\hline chisquare_cdf & chi2cdf in Octave Forge statistics pkg & 3.4.0 \\
\hline chisquare_inv & chi2inv in Octave Forge statistics pkg & 3.4.0 \\
\hline chisquare_pdf & chi2pdf in Octave Forge statistics pkg & 3.4.0 \\
\hline chisquare_rnd & chi2rnd in Octave Forge statistics pkg & 3.4.0 \\
\hline clearplot & clf & 3.4.0 \\
\hline com2str & num2str & 3.4 .0 \\
\hline exponential_cdf & expcdf in Octave Forge statistics pkg & 3.4.0 \\
\hline exponential_inv & expinv in Octave Forge statistics pkg & 3.4.0 \\
\hline exponential_pdf & exppdf in Octave Forge statistics pkg & 3.4.0 \\
\hline
\end{tabular}
\begin{tabular}{lll} 
exponential_rnd & \begin{tabular}{l} 
exprnd in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
f_cdf & \begin{tabular}{l} 
fcdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
finv in Octave Forge & 3.4 .0 \\
f_inv & \begin{tabular}{l} 
statistics pkg \\
fpdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
f_pdf & \begin{tabular}{l} 
frnd in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
f_rnd & \begin{tabular}{l} 
gamcdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
gamma_cdf & \begin{tabular}{l} 
gaminv in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
gamma_inv & \begin{tabular}{l} 
gampdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
gamma_pdf & \begin{tabular}{l} 
gamrnd in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
gamma_rnd & \begin{tabular}{l} 
geocdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
geometric_cdf & \begin{tabular}{l} 
geoinv in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
geometric_inv & \begin{tabular}{l} 
geopdf in Octave Forge
\end{tabular} & 3.4 .0 \\
statistics pkg
\end{tabular}\(\quad\)\begin{tabular}{l} 
geornd in Octave Forge
\end{tabular}\(\quad 3.4 .0\)
\begin{tabular}{lll} 
lognormal_cdf & \begin{tabular}{l} 
logncdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
lognormal_inv & \begin{tabular}{l} 
logninv in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
lognormal_pdf & \begin{tabular}{l} 
lognpdf in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
lognormal_rnd & \begin{tabular}{l} 
lognrnd in Octave Forge \\
statistics pkg
\end{tabular} & 3.4 .0 \\
meshgrid
\end{tabular}\(\quad\)\begin{tabular}{l} 
normcdf in Octave Forge
\end{tabular} 3.4.4.0
\begin{tabular}{|c|c|c|}
\hline uniform_cdf & unifcdf in Octave Forge statistics pkg & 3.4.0 \\
\hline uniform_inv & unifinv in Octave Forge statistics pkg & 3.4 .0 \\
\hline uniform_pdf & unifpdf in Octave Forge statistics pkg & 3.4.0 \\
\hline uniform_rnd & unifrnd in Octave Forge statistics pkg & 3.4.0 \\
\hline weibull_cdf & wblcdf in Octave Forge statistics pkg & 3.4.0 \\
\hline weibull_inv & wblinv in Octave Forge statistics pkg & 3.4.0 \\
\hline weibull_pdf & wblpdf in Octave Forge statistics pkg & 3.4.0 \\
\hline weibull_rnd & wblrnd in Octave Forge statistics pkg & 3.4.0 \\
\hline wiener_rnd & wienrnd in Octave Forge statistics pkg & 3.4.0 \\
\hline create_set & unique & 3.6.0 \\
\hline dmult & diag (A) * B & 3.6.0 \\
\hline iscommand & None & 3.6.0 \\
\hline israwcommand & None & 3.6.0 \\
\hline lchol & chol (..., "lower") & 3.6.0 \\
\hline loadimage & load or imread & 3.6.0 \\
\hline mark_as_command & None & 3.6.0 \\
\hline mark_as_rawcommand & None & 3.6.0 \\
\hline spatan2 & atan2 & 3.6.0 \\
\hline spchol & chol & 3.6.0 \\
\hline spchol2inv & chol2inv & 3.6.0 \\
\hline spcholinv & cholinv & 3.6.0 \\
\hline spcumprod & cumprod & 3.6.0 \\
\hline spcumsum & cumsum & 3.6.0 \\
\hline spdet & det & 3.6.0 \\
\hline spdiag & sparse (diag (...)) & 3.6.0 \\
\hline spfind & find & 3.6.0 \\
\hline sphcat & horzcat & 3.6.0 \\
\hline spinv & inv & 3.6.0 \\
\hline spkron & kron & 3.6.0 \\
\hline splchol & chol (..., "lower") & 3.6.0 \\
\hline split & ```
char (strsplit (s,
t))
``` & 3.6.0 \\
\hline splu & lu & 3.6.0 \\
\hline spmax & max & 3.6.0 \\
\hline spmin & min & 3.6.0 \\
\hline spprod & prod & 3.6.0 \\
\hline spqr & qr & 3.6.0 \\
\hline
\end{tabular}
\begin{tabular}{|c|c|c|}
\hline spsum & sum & 3.6.0 \\
\hline spsumsq & sumsq & 3.6 .0 \\
\hline spvcat & vertcat & 3.6 .0 \\
\hline str2mat & char & 3.6 .0 \\
\hline unmark_command & None & 3.6 .0 \\
\hline unmark_rawcommand & None & 3.6 .0 \\
\hline autocor & xcorr in Octave Forge signal pkg & 3.8 .0 \\
\hline autocov & xcov in Octave Forge signal pkg & 3.8 .0 \\
\hline betai & betainc & 3.8.0 \\
\hline cellidx & ismember & 3.8.0 \\
\hline cquad & quadcc & 3.8 .0 \\
\hline dispatch & None & 3.8 .0 \\
\hline fstat & stat & 3.8.0 \\
\hline gammai & gammainc & 3.8.0 \\
\hline glpkmex & glpk & 3.8 .0 \\
\hline is_duplicate_entry & unique & 3.8 .0 \\
\hline is_global & isglobal & 3.8 .0 \\
\hline krylovb & \[
\begin{aligned}
& \text { [Uret, ~, Ucols] = } \\
& \text { krylov (...) }
\end{aligned}
\] & 3.8.0 \\
\hline perror & None & 3.8.0 \\
\hline replot & refresh & 3.8.0 \\
\hline saveimage & imwrite & 3.8 .0 \\
\hline setstr & char & 3.8.0 \\
\hline strerror & None & 3.8.0 \\
\hline values & unique & 3.8 .0 \\
\hline cut & histc & 4.0 .0 \\
\hline cor & corr & 4.0 .0 \\
\hline corrcoef & corr & 4.0.0 \\
\hline __error_text__ & lasterr & 4.0.0 \\
\hline error_text & lasterr & 4.0.0 \\
\hline polyderiv & polyder & 4.0.0 \\
\hline shell_cmd & system & 4.0.0 \\
\hline studentize & zscore & 4.0.0 \\
\hline sylvester_matrix & hadamard (2^k) & 4.0.0 \\
\hline default_save_options & save_default_options & 4.2 .0 \\
\hline gen_doc_cache & doc_cache_create & 4.2 .0 \\
\hline interp1q & interp1 & 4.2 .0 \\
\hline isequalwithequalnans & isequaln & 4.2 .0 \\
\hline java_convert_matrix & \begin{tabular}{l}
java_matrix_ \\
autoconversion
\end{tabular} & 4.2.0 \\
\hline java_debug & debug_java & 4.2.0 \\
\hline java_invoke & javaMethod & 4.2 .0 \\
\hline java_new & javaObject & 4.2 .0 \\
\hline
\end{tabular}
\begin{tabular}{lll} 
java_unsigned_ & java_unsigned_ & 4.2 .0 \\
conversion & autoconversion & \\
javafields & fieldnames & 4.2 .0 \\
javamethods & methods & 4.2 .0 \\
re_read_readline_init_ & readline_re_read_ & 4.2 .0 \\
file & init_file & \\
read_readline_init_ & readline_read_init_ & 4.2 .0 \\
file & file & \\
saving_history & history_save & 4.2 .0 \\
allow_noninteger_ & None & 4.4 .0 \\
range_as_index & & \\
bicubic & interp2 & 4.4 .0 \\
delaunay3 & delaunay & 4.4 .0 \\
do_braindead_ & None & 4.4 .0 \\
shortcircuit_ & & \\
evaluation & & \\
dump_prefs & None 4.4.0 & \\
find_dir_in_path & dir_in_loadpath & 4.4 .0 \\
finite & isfinite & 4.4 .0 \\
fmod & rem & 4.4 .0 \\
fnmatch & glob or regexp & 4.4 .0 \\
gmap40 & None & 4.4 .0 \\
loadaudio & audioread & 4.4 .0 \\
luinc & icholor ilu & 4.4 .0 \\
mouse_wheel_zoom & mousehweelzoom & 4.4 .0 \\
& property & \\
nfields & numfields & 4.4 .0 \\
octave_tmp_file_name & tempname & 4.4 .0 \\
playaudio & audioplayer & 4.4 .0 \\
saveaudio & audiowrite & 4.4 .0 \\
setaudio & None & 4.4 .0 \\
syl & sylvester & 4.4 .0 \\
usage & print_usage & 4.4 .0 \\
bitmax & flintmax & 5.1 .0 \\
mahalanobis & mahal in Octave Forge & 5.1 .0 \\
md5sum & statistics pkg & \\
octave_config_info & hash & 5.1 .0 \\
& -_octave_config_ & 5.1 .0 \\
onenormest & info_- & \\
sleep & normest1 & 5.1 .0 \\
usleep & pause & 5.1 .0 \\
wavread & audioread & 5.1 .0 \\
wavwrite & audiowrite & 5.1 .0 \\
& 5.1.0 \\
& & \\
\hline
\end{tabular}

\section*{Appendix D Known Causes of Trouble}

This section describes known problems that affect users of Octave. Most of these are not Octave bugs per se - if they were, we would fix them. But the result for a user may be like the result of a bug.

Some of these problems are due to bugs in other software, some are missing features that are too much work to add, and some are places where people's opinions differ as to what is best.

\section*{D. 1 Actual Bugs We Haven't Fixed Yet}
- Output that comes directly from Fortran functions is not sent through the pager and may appear out of sequence with other output that is sent through the pager. One way to avoid this is to force pending output to be flushed before calling a function that will produce output from within Fortran functions. To do this, use the command
```

fflush (stdout)

```

Another possible workaround is to use the command
```

    page_screen_output (false);
    to turn the pager off.

```

A list of ideas for future enhancements is distributed with Octave. See the file PROJECTS in the top level directory in the source distribution.

\section*{D. 2 Reporting Bugs}

Your bug reports play an essential role in making Octave reliable.
When you encounter a problem, the first thing to do is to see if it is already known. See Appendix D [Trouble], page 1079. If it isn't known, then you should report the problem.

Reporting a bug may help you by bringing a solution to your problem, or it may not. In any case, the principal function of a bug report is to help the entire community by making the next version of Octave work better. Bug reports are your contribution to the maintenance of Octave.

In order for a bug report to serve its purpose, you must include the information that makes it possible to fix the bug.

\section*{D.2.1 Have You Found a Bug?}

If you are not sure whether you have found a bug, here are some guidelines:
- If Octave gets a fatal signal, for any input whatever, that is a bug. Reliable interpreters never crash.
- If Octave produces incorrect results, for any input whatever, that is a bug.
- Some output may appear to be incorrect when it is in fact due to a program whose behavior is undefined, which happened by chance to give the desired results on another system. For example, the range operator may produce different results because of differences in the way floating point arithmetic is handled on various systems.
- If Octave produces an error message for valid input, that is a bug.
- If Octave does not produce an error message for invalid input, that is a bug. However, you should note that your idea of "invalid input" might be my idea of "an extension" or "support for traditional practice".
- If you are an experienced user of programs like Octave, your suggestions for improvement are welcome in any case.

\section*{D.2.2 Where to Report Bugs}

To report a bug in Octave, submit a bug report to the Octave bug tracker https://bugs . octave.org.

Do not send bug reports to 'help-octave'. Most users of Octave do not want to receive bug reports.

\section*{D.2.3 How to Report Bugs}

Submit bug reports for Octave to the Octave bug tracker https://bugs.octave.org.
The fundamental principle of reporting bugs usefully is this: report all the facts. If you are not sure whether to state a fact or leave it out, state it!

Often people omit facts because they think they know what causes the problem and they conclude that some details don't matter. Thus, you might assume that the name of the variable you use in an example does not matter. Well, probably it doesn't, but one cannot be sure. Perhaps the bug is a stray memory reference which happens to fetch from the location where that name is stored in memory; perhaps, if the name were different, the contents of that location would fool the interpreter into doing the right thing despite the bug. Play it safe and give a specific, complete example.

Keep in mind that the purpose of a bug report is to enable someone to fix the bug if it is not known. Always write your bug reports on the assumption that the bug is not known.

Sometimes people give a few sketchy facts and ask, "Does this ring a bell?" This cannot help us fix a bug. It is better to send a complete bug report to begin with.

Try to make your bug report self-contained. If we have to ask you for more information, it is best if you include all the previous information in your response, as well as the information that was missing.

To enable someone to investigate the bug, you should include all these things:
- The version of Octave. You can get this by noting the version number that is printed when Octave starts, or running it with the ' -v ' option.
- A complete input file that will reproduce the bug.

A single statement may not be enough of an example - the bug might depend on other details that are missing from the single statement where the error finally occurs.
- The command arguments you gave Octave to execute that example and observe the bug. To guarantee you won't omit something important, list all the options.
If we were to try to guess the arguments, we would probably guess wrong and then we would not encounter the bug.
- The type of machine you are using, and the operating system name and version number.
- The command-line arguments you gave to the configure command when you installed the interpreter.
- A complete list of any modifications you have made to the interpreter source.

Be precise about these changes-show a context diff for them.
- Details of any other deviations from the standard procedure for installing Octave.
- A description of what behavior you observe that you believe is incorrect. For example, "The interpreter gets a fatal signal," or, "The output produced at line 208 is incorrect." Of course, if the bug is that the interpreter gets a fatal signal, then one can't miss it. But if the bug is incorrect output, we might not notice unless it is glaringly wrong.
Even if the problem you experience is a fatal signal, you should still say so explicitly. Suppose something strange is going on, such as, your copy of the interpreter is out of sync, or you have encountered a bug in the C library on your system. Your copy might crash and the copy here would not. If you said to expect a crash, then when the interpreter here fails to crash, we would know that the bug was not happening. If you don't say to expect a crash, then we would not know whether the bug was happening. We would not be able to draw any conclusion from our observations.
Often the observed symptom is incorrect output when your program is run. Unfortunately, this is not enough information unless the program is short and simple. It is very helpful if you can include an explanation of the expected output, and why the actual output is incorrect.
- If you wish to suggest changes to the Octave source, send them as context diffs. If you even discuss something in the Octave source, refer to it by context, not by line number, because the line numbers in the development sources probably won't match those in your sources.
Here are some things that are not necessary:
- A description of the envelope of the bug.

Often people who encounter a bug spend a lot of time investigating which changes to the input file will make the bug go away and which changes will not affect it. Such information is usually not necessary to enable us to fix bugs in Octave, but if you can find a simpler example to report instead of the original one, that is a convenience. Errors in the output will be easier to spot, running under the debugger will take less time, etc. Most Octave bugs involve just one function, so the most straightforward way to simplify an example is to delete all the function definitions except the one in which the bug occurs.
However, simplification is not vital; if you don't want to do this, report the bug anyway and send the entire test case you used.
- A patch for the bug. Patches can be helpful, but if you find a bug, you should report it, even if you cannot send a fix for the problem.

\section*{D.2.4 Sending Patches for Octave}

If you would like to write bug fixes or improvements for Octave, that is very helpful. When you send your changes, please follow these guidelines to avoid causing extra work for us in studying the patches.

If you don't follow these guidelines, your information might still be useful, but using it will take extra work. Maintaining Octave is a lot of work in the best of circumstances, and we can't keep up unless you do your best to help.
- Send an explanation with your changes of what problem they fix or what improvement they bring about. For a bug fix, just include a copy of the bug report, and explain why the change fixes the bug.
- Always include a proper bug report for the problem you think you have fixed. We need to convince ourselves that the change is right before installing it. Even if it is right, we might have trouble judging it if we don't have a way to reproduce the problem.
- Include all the comments that are appropriate to help people reading the source in the future understand why this change was needed.
- Don't mix together changes made for different reasons. Send them individually.

If you make two changes for separate reasons, then we might not want to install them both. We might want to install just one.
- Use 'diff -c' to make your diffs. Diffs without context are hard for us to install reliably. More than that, they make it hard for us to study the diffs to decide whether we want to install them. Unified diff format is better than contextless diffs, but not as easy to read as ' \(-c\) ' format.
If you have GNU diff, use 'diff -cp', which shows the name of the function that each change occurs in.
- Write the change log entries for your changes.

Read the ChangeLog file to see what sorts of information to put in, and to learn the style that we use. The purpose of the change log is to show people where to find what was changed. So you need to be specific about what functions you changed; in large functions, it's often helpful to indicate where within the function the change was made.
On the other hand, once you have shown people where to find the change, you need not explain its purpose. Thus, if you add a new function, all you need to say about it is that it is new. If you feel that the purpose needs explaining, it probably does-but the explanation will be much more useful if you put it in comments in the code.
If you would like your name to appear in the header line for who made the change, send us the header line.

\section*{D. 3 How To Get Help with Octave}

The mailing list help@octave.org exists for the discussion of matters related to using and installing Octave. If would like to join the discussion, please send a short note to help-request@octave.org.

Please do not send requests to be added or removed from the mailing list, or other administrative trivia to the list itself.

If you think you have found a bug in Octave or in the installation procedure, however, you should submit a complete bug report to the Octave bug tracker at https://bugs. octave.org. But before you submit a bug report, please read https://www.octave.org/ bugs.html to learn how to submit a useful bug report.

\section*{D. 4 How to Distinguish Between Octave and Matlab}

Octave and matlab are very similar, but handle Java slightly different. Therefore it may be necessary to detect the environment and use the appropriate functions. The following
function can be used to detect the environment. Due to the persistent variable it can be called repeatedly without a heavy performance hit.

Example:
```

%%
%% Return: true if the environment is Octave.
%%
function retval = isOctave
persistent cacheval; % speeds up repeated calls
if isempty (cacheval)
cacheval = (exist ("OCTAVE_VERSION", "builtin") > 0);
end
retval = cacheval;
end

```

\section*{Appendix E Installing Octave}

The procedure for installing Octave from source on a Unix-like system is described next. Building on other platforms will follow similar steps. Note that this description applies to Octave releases. Building the development sources from the Mercurial archive requires additional steps as described in the development source itself.

\section*{E. 1 Build Dependencies}

Octave is a fairly large program with many build dependencies. You may be able to find pre-packaged versions of the dependencies distributed as part of your system, or you may have to build some or all of them yourself.

\section*{E.1.1 Obtaining the Dependencies Automatically}

On some systems you can obtain many of Octave's build dependencies automatically. The commands for doing this vary by system. Similarly, the names of pre-compiled packages vary by system and do not always match exactly the names listed in Section E.1.2 [Build Tools], page 1085, and Section E.1.3 [External Packages], page 1086.

You will usually need the development version of an external dependency so that you get the libraries and header files for building software, not just for running already compiled programs. These packages typically have names that end with the suffix -dev or -devel.

On systems with apt-get (Debian, Ubuntu, etc.), you may be able to install most of the tools and external packages using a command similar to
```

apt-get build-dep octave

```

The specific package name may be octave3.2 or octave3.4. The set of required tools and external dependencies does not change frequently, so it is not important that the version match exactly, but you should use the most recent one available.

On systems with yum (Fedora, Red Hat, etc.), you may be able to install most of the tools and external packages using a command similar to
```

yum-builddep octave

```

The yum-builddep utility is part of the yum-utils package.
For either type of system, the package name may include a version number. The set of required tools and external dependencies does not change frequently, so it is not important that the version exactly match the version you are installing, but you should use the most recent one available.

\section*{E.1.2 Build Tools}

The following tools are required:
C++, C, and Fortran compilers
The Octave sources are primarily written in C++, but some portions are also written in C and Fortran. The Octave sources are intended to be portable. Recent versions of the GNU compiler collection (GCC) should work (https:// gcc.gnu. org). If you use GCC, you should avoid mixing versions. For example, be sure that you are not using the obsolete g77 Fortran compiler with modern versions of gcc and g++.

GNU Make
Tool for building software (https://www.gnu.org/software/make). Octave's build system requires GNU Make. Other versions of Make will not work. Fortunately, GNU Make is highly portable and easy to install.

AWK, sed, and other Unix utilities
Basic Unix system utilities are required for building Octave. All will be available with any modern Unix system and also on Windows with either Cygwin or MinGW and MSYS.

Additionally, the following tools may be needed:
Bison Parser generator (https://www.gnu.org/software/bison). You will need Bison if you modify the oct-parse.yy source file or if you delete the files that are generated from it.

Flex Lexer analyzer (https://www.gnu.org/software/flex). You will need Flex if you modify the lex.ll source file or if you delete the files that are generated from it.

Autoconf Package for software configuration (https://www.gnu.org/software/ autoconf). Autoconf is required if you modify Octave's configure.ac file or other files that it requires.

Automake Package for Makefile generation (https://www.gnu.org/software/automake). Automake is required if you modify Octave's Makefile.am files or other files that they depend on.

Libtool Package for building software libraries (https://www.gnu.org/software/ libtool). Libtool is required by Automake.
gperf Perfect hash function generator (https://www.gnu.org/software/gperf). You will need gperf if you modify the octave.gperf file or if you delete the file that is generated from it.

Texinfo Package for generating online and print documentation (https://www.gnu. org/software/texinfo). You will need Texinfo to build Octave's documentation or if you modify the documentation source files or the docstring of any Octave function.

\section*{E.1.3 External Packages}

The following external packages are required:
BLAS Basic Linear Algebra Subroutine library. Accelerated blas libraries such as OpenBLAS (https://www.openblas.net/) or ATLAS (http://math-atlas. sourceforge.net) are recommended for best performance. The reference implementation (http://www.netlib.org/blas) is slow and suffers from certain bugs in corner case inputs.

LAPACK Linear Algebra Package (http://www.netlib.org/lapack).
PCRE The Perl Compatible Regular Expression library (https://www.pcre.org).

The following external package is optional but strongly recommended:
GNU Readline
Command-line editing library (https://www.gnu.org/s/readline).
If you wish to build Octave without GNU readline installed, you must use the --disable-readline option when running the configure script.

The following external software packages are optional. Octave can be built without them but certain features might be missing:
ARPACK Library for the solution of large-scale eigenvalue problems (https://forge. scilab.org/index.php/p/arpack-ng). ARPACK is required to provide the functions eigs and svds.
cURL Library for transferring data with URL syntax (https://curl.haxx. se). cURL is required to provide the urlread and urlwrite functions and the ftp class.
FFTW3 Library for computing discrete Fourier transforms (http://www.fftw.org). FFTW3 is used to provide better performance for functions that compute discrete Fourier transforms (fft, ifft, fft2, etc.)
FLTK Portable GUI toolkit (http://www.fltk.org). FLTK can be used to provide windows for Octave's OpenGL-based graphics functions.
fontconfig Library for configuring and customizing font access (https://www. freedesktop.org/wiki/Software/fontconfig). Fontconfig is used to manage fonts for Octave's OpenGL-based graphics functions.
FreeType Portable font engine (https://www.freetype.org). FreeType is used to perform font rendering for Octave's OpenGL-based graphics functions.
GLPK GNU Linear Programming Kit (https://www.gnu.org/software/glpk). GPLK is required for the function glpk.
gl2ps OpenGL to PostScript printing library (https://www.geuz.org/gl2ps/ ). gl2ps is required for printing when using OpenGL-based graphics toolkits (currently either FLTK or Qt).
gnuplot Interactive graphics program (http://www.gnuplot.info). gnuplot can be used as a graphics renderer for Octave; prior to Octave 4.0, gnuplot was the default graphics renderer.
GraphicsMagick++
Image processing library (http://www.graphicsmagick.org). GraphicsMagick++ is used to provide the imread and imwrite functions.
HDF5 Library for manipulating portable data files (https://www.hdfgroup.org/ HDF5). HDF5 is required for Octave's load and save commands to read and write HDF data files.

Java Development Kit
Java programming language compiler and libraries. The OpenJDK free software implementation is recommended (http://openjdk.java.net/), although other JDK implementations may work. Java is required to be able to call Java functions from within Octave.

OpenGL API for portable 2-D and 3-D graphics (https://www.opengl.org). An OpenGL implementation can be used to provide a renderer for Octave's graphics functions. Octave's OpenGL-based graphics functions usually outperform the gnuplot-based graphics functions because plot data can be rendered directly instead of sending data and commands to gnuplot for interpretation and rendering. Since Octave 4.0, the default graphics renderer ("qt") has been OpenGL-based.

PortAudio PortAudio (http://www.portaudio.com/) provides a very simple API for recording and/or playing sound using a simple callback function or a blocking read/write interface. It is required for the audio processing functions audioplayer, audiorecorder, and audiodevinfo.

Qhull Computational geometry library (http://www.qhull.org). Qhull is required to provide the functions convhull, convhulln, delaunay, delaunayn, voronoi, and voronoin.

\section*{QRUPDATE}

QR factorization updating library (https://sourceforge.net/projects/ qrupdate). QRUPDATE is used to provide improved performance for the functions qrdelete, qrinsert, qrshift, and qrupdate.

QScintilla Source code highlighter and manipulator; a Qt port of Scintilla (http://www. riverbankcomputing.co.uk/software/qscintilla). QScintilla is used for syntax highlighting and code completion in the GUI.

Qt GUI and utility libraries (https://www.qt.io). Qt is required for building the GUI. It is a large framework, but the only components required are the GUI, core, and network modules. Since Octave 4.0, the default graphics renderer ("qt") has been Qt-based, which has been OpenGL-based.

RapidJSON
A fast JSON parser/generator for C++ with both SAX/DOM style API (https://rapidjson.org/). RapidJSON is required to read or write from or to JSON files with the functions jsondecode and jsonencode.

SuiteSparse
Sparse matrix factorization library (http://faculty.cse.tamu.edu/davis/ suitesparse.html). SuiteSparse is required to provide sparse matrix factorizations and solution of linear equations for sparse systems.

SUNDIALS
The SUite of Nonlinear and DIfferential/ALgebraic Equation Solvers (https:// computation.llnl.gov/projects/sundials) is required for the Ordinary Differential Equations (ODE) solvers ode15i and ode15s.
zlib Data compression library (https://zlib.net). The zlib library is required for Octave's load and save commands to handle compressed data, including matlab v5 MAT files.

\section*{E. 2 Running Configure and Make}
- Run the shell script configure. This will determine the features your system has (or doesn't have) and create a file named Makefile from each of the files named Makefile.in.
For a complete list of configure options, run configure --help. Here is a summary of the configure options that are most frequently used when building Octave:
--help Print a summary of the options recognized by the configure script.
--prefix=prefix
Install Octave in subdirectories below prefix. The default value of prefix is /usr/local.

\section*{--srcdir=dir}

Look for Octave sources in the directory dir.

\section*{--disable-64}

Disable using 64 -bit integers for indexing arrays and use 32 -bit integers instead. On systems with 32 -bit pointers, this option is always disabled. If the configure script determines that your BLAS library uses 32-bit integers, then operations using the following libraries are limited to arrays with dimensions that are smaller than \(2^{31}\) elements:
- blas
- LAPACK
- QRUPDATE
- SuiteSparse
- ARPACK

Additionally, the following libraries use int internally, so maximum problem sizes are always limited:
- GLPK
- Qhull

See Section E. 3 [Compiling Octave with 64-bit Indexing], page 1093, for more details about building Octave with more complete support for large arrays.

Enable compiler options -fsanitize=address and -fomit-framepointer for memory access checking. This option is primarily used for debugging Octave. Building Octave with this option has a negative impact on performance and is not recommended for general use. It may also interfere with proper functioning of the GUI.
--disable-docs
Disable building all forms of the documentation (Info, PDF, HTML). The default is to build documentation, but your system will need functioning Texinfo and \(\mathrm{T}_{\mathrm{E}} \mathrm{X}\) installs for this to succeed.
--enable-float-truncate
This option allows for truncation of intermediate floating point results in calculations. It is only necessary for certain platforms.
--enable-readline
Use the readline library to provide for editing of the command line in terminal environments. This option is on by default.
--enable-shared
Create shared libraries (this is the default). If you are planning to use the dynamic loading features, you will probably want to use this option. It will make your .oct files much smaller and on some systems it may be necessary to build shared libraries in order to use dynamically linked functions.
You may also want to build a shared version of libstdc++, if your system doesn't already have one.
--with-blas=<lib>
By default, configure looks for the best BLAS matrix libraries on your system, including optimized implementations such as the free ATLAS 3.0, as well as vendor-tuned libraries. (The use of an optimized BLAS will generally result in several-times faster matrix operations.) Use this option to specify a particular BLAS library that Octave should use.
--with-lapack=<lib>
By default, configure looks for the best LAPACK matrix libraries on your system, including optimized implementations such as the free ATLAS 3.0, as well as vendor-tuned libraries. (The use of an optimized LAPACK will generally result in several-times faster matrix operations.) Use this option to specify a particular LAPACK library that Octave should use.
--with-magick=<lib>
Select the Magick++ library to use for image I/O. For many distributions, possible values are "GraphicsMagick++" (default) or "ImageMagick++".
--with-sepchar=<char>
Use <char> as the path separation character. This option can help when running Octave on non-Unix systems.
--without-amd
Don't use AMD, disable some sparse matrix functionality.
--without-camd
Don't use CAMD, disable some sparse matrix functionality.
--without-colamd
Don't use COLAMD, disable some sparse matrix functionality.
--without-ccolamd
Don't use CCOLAMD, disable some sparse matrix functionality.
--without-cholmod
Don't use CHOLMOD, disable some sparse matrix functionality.
--without-curl
Don't use the cURL library, disable the ftp objects, urlread and urlwrite functions.
--without-cxsparse
Don't use CXSPARSE, disable some sparse matrix functionality.
--without-fftw3
Use the included fFTPACK library for computing Fast Fourier Transforms instead of the FFTW3 library.
--without-fftw3f
Use the included FFTPACK library for computing Fast Fourier Transforms instead of the FFTW3 library when operating on single precision (float) values.
--without-glpk
Don't use the GLPK library for linear programming.
--without-hdf5
Don't use the HDF5 library, disable reading and writing of HDF5 files.
--without-opengl
Don't use OpenGL, disable native graphics toolkit for plotting. You will need gnuplot installed in order to make plots.
--without-qhull_r
Don't use (re-entrant) Qhull, disable delaunay, convhull, and related functions.
--without-qrupdate
Don't use QRUPDATE, disable QR and Cholesky update functions.
--without-umfpack
Don't use UMFPACK, disable some sparse matrix functionality.
--without-z
Don't use the zlib library, disable data file compression and support for recent MAT file formats.
--without-framework-carbon
Don't use framework Carbon headers, libraries, or specific source code even if the configure test succeeds (the default is to use Carbon framework if available). This is a platform specific configure option for Mac systems.
--without-framework-opengl
Don't use framework OpenGL headers, libraries, or specific source code even if the configure test succeeds. If this option is given then OpenGL headers and libraries in standard system locations are tested (the default value is --with-framework-opengl). This is a platform specific configure option for Mac systems.
See the file INSTALL for more general information about the command line options used by configure. That file also contains instructions for compiling in a directory other than the one where the source is located.
- Run make.

You will need a recent version of GNU Make as Octave relies on certain features not generally available in all versions of make. Modifying Octave's makefiles to work with other make programs is probably not worth your time; instead, we simply recommend installing GNU Make.
There are currently three options for plotting in Octave: the external program gnuplot, the internal graphics engine using OpenGL coupled with either FLTK or Qt widgets. Gnuplot is a command-driven interactive function plotting program.
To compile Octave, you will need a recent version of g++ or other ANSI C++ compiler. In addition, you will need a Fortran 77 compiler or f2c. If you use f2c, you will need a script like fort77 that works like a normal Fortran compiler by combining f2c with your C compiler in a single script.
If you plan to modify the parser you will also need GNU bison and flex. If you modify the documentation, you will need GNU Texinfo.
GNU Make, gcc (and libstdc++), gnuplot, bison, flex, and Texinfo are all available from many anonymous ftp archives. The primary site is ftp://ftp.gnu.org, but it is often very busy. A list of sites that mirror the software on ftp://ftp.gnu.org is available by anonymous ftp from ftp://ftp.gnu.org/pub/gnu/GNUinfo/FTP.
Octave requires approximately 1.4 GB of disk storage to unpack and compile from source (significantly less, 400 MB , if you don't compile with debugging symbols). To compile without debugging symbols try the command
make CFLAGS=-0 CXXFLAGS=-0 LDFLAGS=
instead of just make.
- If you encounter errors while compiling Octave, first see Section E. 4 [Installation Problems], page 1096, for a list of known problems and if there is a workaround or solution for your problem. If not, see Appendix D [Trouble], page 1079, for information about how to report bugs.
- Once you have successfully compiled Octave, run make install.

This will install a copy of Octave, its libraries, and its documentation in the destination directory. As distributed, Octave is installed in the following directories. In the table below, prefix defaults to /usr/local, version stands for the current version number of the interpreter, and arch is the type of computer on which Octave is installed (for example, 'i586-unknown-gnu').
prefix/bin
Octave and other binaries that people will want to run directly.
prefix/lib/octave-version
Libraries like liboctave.a and liboctinterp.a.
prefix/include/octave-version/octave
Include files distributed with Octave.
prefix/share
Architecture-independent data files.
prefix/share/man/man1
Unix-style man pages describing Octave.
```

prefix/share/info
Info files describing Octave.
prefix/share/octave/version/m

```

Function files distributed with Octave. This includes the Octave version, so that multiple versions of Octave may be installed at the same time.
prefix/libexec/octave/version/exec/arch
Executables to be run by Octave rather than the user.
prefix/lib/octave/version/oct/arch
Object files that will be dynamically loaded.
prefix/share/octave/version/imagelib
Image files that are distributed with Octave.

\section*{E. 3 Compiling Octave with 64-bit Indexing}

Note: the following only applies to systems that have 64 -bit pointers. Configuring Octave with --enable-64 cannot magically make a 32 -bit system have a 64 -bit address space.

On 64 -bit systems, Octave uses 64 -bit integers for indexing arrays by default. If the configure script determines that your BLAs library uses 32 -bit integers, then operations using the following libraries are limited to arrays with dimensions that are smaller than \(2^{31}\) elements:
- BLAS
- Lapack
- QRUPDATE
- SuiteSparse
- SUNDIALS IDA
- ARPACK

Additionally, the following libraries use int internally, so maximum problem sizes are always limited:
- GLPK
- Qhull

Except for GLPK and Qhull, these libraries may also be configured to use 64-bit integers, but most systems do not provide packages built this way. If you wish to experiment with large arrays, the following information may be helpful.

To determine the integer size of the BLAS library used by Octave, the following code can be executed:
```

clear all;
N = 2^31;

## The following line requires about 8 GB of RAM!

a = b = ones (N, 1, "single");
c = a' * b

```

If the blas library uses 32 -bit integers, an error will be thrown:
```

error: integer dimension or index out of range for Fortran
INTEGER type

```

Otherwise, if the BLAS library uses 64-bit integers, the result is:
```

c = 2^31 = 2147483648

```

Note that the test case above usually requires twice the memory, if a and \(b\) are not assigned by \(\mathrm{a}=\mathrm{b}=\ldots\). Note further, that the data type single has a precision of about 23 binary bits. In this particular example no rounding errors occur.

Generally, it is best to have all of these libraries in versions that support 32-bit indexing, or all of these libraries must support 64 -bit indexing. Mixing libraries with 64 -bit indexing with libraries with 32-bit indexing can cause unpredictable behavior including program crashes with possible loss of data.

The following instructions were tested with the development version of Octave and GCC 4.3.4 on an x86_64 Debian system and may be out of date now. Please report any problems or corrections on the Octave bug tracker.

The versions listed below are the versions used for testing. If newer versions of these packages are available, you should try to use them, although there may be some differences.

All libraries and header files will be installed in subdirectories of \$prefix64 (you must choose the location of this directory).
- BLAS and LAPACK (http://www.netlib.org/lapack)

Reference versions for both libraries are included in the reference LAPACK 3.2.1 distribution from http://www.netlib.org/.
- Copy the file make.inc. example and name it make.inc. The options -fdefault-integer-8 and -fPIC (on 64-bit CPU) have to be added to the variable OPTS and NOOPT.
- Once you have compiled this library make sure that you use it for compiling Suite Sparse and Octave. In the following we assume that you installed the LAPACK library as \$prefix64/lib/liblapack.a.
- QRUPDATE (https://sourceforge.net/projects/qrupdate)

In the Makeconf file:
- Add -fdefault-integer-8 to FFLAGS.
- Adjust the BLAS and LAPACK variables as needed if your 64-bit aware BLAS and LAPACK libraries are in a non-standard location.
- Set PREFIX to the top-level directory of your install tree.
- Run make solib to make a shared library.
- Run make install to install the library.
- SuiteSparse (http://faculty.cse.tamu.edu/davis/suitesparse.html)

Pass the following options to make to enable 64-bit integers for BLAS library calls. On 64-bit Windows systems, use -DLONGBLAS="long long" instead.
```

CFLAGS=' -DLONGBLAS=long'
CXXFLAGS=' -DLONGBLAS=long '

```

The SuiteSparse makefiles don't generate shared libraries. On some systems, you can generate them by doing something as simple as
```

top=$(pwd)
for f in *.a; do
    mkdir tmp
    cd tmp
    ar vx ../$f
gcc -shared -o ../\${f%%.a}.so *.o
cd \$top
rm -rf tmp
done

```

Other systems may require a different solution.
- SUNDIALS IDA (https://computing.llnl.gov/projects/sundials/ida)

When configuring with cmake add the flag -DSUNDIALS_INDEX_SIZE=64.
- ARPACK (https://forge.scilab.org/index.php/p/arpack-ng/)
- Add -fdefault-integer-8 to FFLAGS when running configure.
- Run make to build the library.
- Run make install to install the library.
- ATLAS instead of reference blas and lapack

Suggestions on how to compile ATLAS would be most welcome.
- GLPK
- Qhull (http://www.qhull.org)

Both GLPK and Qhull use int internally so maximum problem sizes may be limited.
- Octave

Octave's 64-bit index support is activated with the configure option --enable-64.
```

./configure \
LD_LIBRARY_PATH="$prefix64/lib" \
    CPPFLAGS="-I$prefix64/include" LDFLAGS="-L\$prefix64/lib" \
--enable-64

```

You must ensure that all Fortran sources except those in the liboctave/external/ranlib】 directory are compiled such that INTEGERS are 8 -bytes wide. If you are using gfortran, the configure script should automatically set the Makefile variable F77_INTEGER_8_FLAG to -fdefault-integer-8. If you are using another compiler, you must set this variable yourself. You should NOT set this flag in FFLAGS, otherwise the files in liboctave/external/ranlib will be miscompiled.
- Other dependencies

Probably nothing special needs to be done for the following dependencies. If you discover that something does need to be done, please submit a bug report.
- pcre
- zlib
- hdf5
- fftw3
- cURL
- GraphicsMagick++
- OpenGL
- freetype
- fontconfig
- fltk

\section*{E. 4 Installation Problems}

This section contains a list of problems (and some apparent problems that don't really mean anything is wrong) that may show up during installation of Octave.
- On some SCO systems, info fails to compile if HAVE_TERMIOS_H is defined in config.h. Simply removing the definition from info/config.h should allow it to compile.
- If configure finds dlopen, dlsym, dlclose, and dlerror, but not the header file dlfcn.h, you need to find the source for the header file and install it in the directory usr/include. This is reportedly a problem with Slackware 3.1. For Linux/GNU systems, the source for dlfcn.h is in the ldso package.
- Building . oct files doesn't work.

You should probably have a shared version of libstdc++. A patch is needed to build shared versions of version 2.7.2 of libstdc++ on the HP-PA architecture. You can find the patch at ftp://ftp.cygnus.com/pub/g++/libg++-2.7.2-hppa-gcc-fix.
- On some DEC alpha systems there may be a problem with the libdxml library, resulting in floating point errors and/or segmentation faults in the linear algebra routines called by Octave. If you encounter such problems, then you should modify the configure script so that SPECIAL_MATH_LIB is not set to -ldxml.
- On FreeBSD systems Octave may hang while initializing some internal constants. The fix appears to be to use
options GPL_MATH_EMULATE
rather than
options MATH_EMULATE
in the kernel configuration files (typically found in the directory /sys/i386/conf). After making this change, you'll need to rebuild the kernel, install it, and reboot.
- If you encounter errors like
```

passing 'void (*)()' as argument 2 of
'octave_set_signal_handler(int, void (*)(int))'
warning: ANSI C++ prohibits conversion from '(int)'
to '(...)'

```
or
while compiling sighandlers.cc, you may need to edit some files in the gcc include subdirectory to add proper prototypes for functions there. For example, Ultrix 4.2 needs proper declarations for the signal function and the SIG_IGN macro in the file signal.h.
On some systems the SIG_IGN macro is defined to be something like this:
```

\#define SIG_IGN (void (*)())1

```
when it should really be something like:
```

\#define SIG_IGN (void (*)(int))1

```
to match the prototype declaration for the signal function. This change should also be made for the SIG_DFL and SIG_ERR symbols. It may be necessary to change the definitions in sys/signal.h as well.

The gcc fixincludes and fixproto scripts should probably fix these problems when gcc installs its modified set of header files, but I don't think that's been done yet.

You should not change the files in /usr/include. You can find the gcc include directory tree by running the command
```

gcc -print-libgcc-file-name

```

The directory of gcc include files normally begins in the same directory that contains the file libgcc.a.
- Some of the Fortran subroutines may fail to compile with older versions of the Sun Fortran compiler. If you get errors like
```

zgemm.f:
zgemm:
warning: unexpected parent of complex expression subtree
zgemm.f, line 245: warning: unexpected parent of complex
expression subtree
warning: unexpected parent of complex expression subtree
zgemm.f, line 304: warning: unexpected parent of complex
expression subtree
warning: unexpected parent of complex expression subtree
zgemm.f, line 327: warning: unexpected parent of complex
expression subtree
pcc_binval: missing IR_CONV in complex op
make[2]: *** [zgemm.o] Error 1

```
when compiling the Fortran subroutines in the liboctave/external subdirectory, you should either upgrade your compiler or try compiling with optimization turned off.
- On NeXT systems, if you get errors like this:
```

/usr/tmp/cc007458.s:unknown:Undefined local
symbol LBB7656
/usr/tmp/cc007458.s:unknown:Undefined local
symbol LBE7656

```
when compiling Array.cc and Matrix.cc, try recompiling these files without -g.
- Some people have reported that calls to system() and the pager do not work on SunOS systems. This is apparently due to having G_HAVE_SYS_WAIT defined to be 0 instead of 1 when compiling libg++.
- On systems where the reference Blas library is used the following matrix-by-vector multiplication incorrectly handles NaN values of the form \(\mathrm{NaN} * 0\).
```

[NaN, 1; 0, 0] * [0; 1]

# 

[ 1
0 ]
correct result }
[ NaN
0 ]

```

Install a different BLAS library such as OpenBLAS or ATLAS to correct this issue.
- On NeXT systems, linking to libsys_s.a may fail to resolve the following functions
```

_tcgetattr
_tcsetattr
_tcflow

```
which are part of libposix.a. Unfortunately, linking Octave with -posix results in the following undefined symbols.
```

.destructors_used
.constructors_used
_objc_msgSend
_NXGetDefaultValue
_NXRegisterDefaults
.objc_class_name_NXStringTable
.objc_class_name_NXBundle

```

One kluge around this problem is to extract termios.o from libposix.a, put it in Octave's src directory, and add it to the list of files to link together in the makefile. Suggestions for better ways to solve this problem are welcome!
- If Octave crashes immediately with a floating point exception, it is likely that it is failing to initialize the IEEE floating point values for infinity and NaN .
If your system actually does support IEEE arithmetic, you should be able to fix this problem by modifying the function octave_ieee_init in the file lo-ieee.cc to correctly initialize Octave's internal infinity and NaN variables.
If your system does not support IEEE arithmetic but Octave's configure script incorrectly determined that it does, you can work around the problem by editing the file config.h to not define HAVE_ISINF, HAVE_FINITE, and HAVE_ISNAN.
In any case, please report this as a bug since it might be possible to modify Octave's configuration script to automatically determine the proper thing to do.
- If Octave is unable to find a header file because it is installed in a location that is not normally searched by the compiler, you can add the directory to the include search path by specifying (for example) CPPFLAGS=-I/some/nonstandard/directory as an argument to configure. Other variables that can be specified this way are CFLAGS, CXXFLAGS, FFLAGS, and LDFLAGS. Passing them as options to the configure script also records them in the config.status file. By default, CPPFLAGS and LDFLAGS are empty, CFLAGS and CXXFLAGS are set to "-g -02" and FFLAGS is set to "-0".

\section*{Appendix F Grammar and Parser}

This appendix will eventually contain a semi-formal description of Octave's language.

\section*{F. 1 Keywords}

The following identifiers are keywords, and may not be used as variable or function names:
\begin{tabular}{lll} 
__FILE__ & __LINE__ & break \\
case & catch & classdef \\
continue & do & else \\
elseif & end & end_try_catch \\
end_unwind_protect & endclassdef & endenumeration \\
endevents & endfor & endfunction \\
endif & endmethods & endparfor \\
endproperties & endswitch & endwhile \\
enumeration & events & for \\
function & global & if \\
methods & otherwise & parfor \\
persistent & properties & return \\
switch & try & until \\
unwind_protect & unwind_protect_cleanup & while
\end{tabular}

The function iskeyword can be used to quickly check whether an identifier is reserved by Octave.
iskeyword ()
iskeyword (name)
Return true if name is an Octave keyword.
If name is omitted, return a list of keywords.
See also: [isvarname], page 137, [exist], page 144.

\section*{F. 2 Parser}

The parser has a number of variables that affect its internal operation. These variables are generally documented in the manual alongside the code that they affect.

In addition, there are three non-specific parser customization functions. add_input_ event_hook can be used to schedule a user function for periodic evaluation. remove_ input_event_hook will stop a user function from being evaluated periodically.
```

id = add_input_event_hook (fcn)
id = add_input_event_hook (fcn, data)

```

Add the named function or function handle fcn to the list of functions to call periodically when Octave is waiting for input.
The function should have the form
```

fcn (data)

```

If data is omitted, Octave calls the function without any arguments.

The returned identifier may be used to remove the function handle from the list of input hook functions.

See also: [remove_input_event_hook], page 1100.
remove_input_event_hook (name)
remove_input_event_hook ( \(f\) cn_id)
Remove the named function or function handle with the given identifier from the list of functions to call periodically when Octave is waiting for input.
See also: [add_input_event_hook], page 1099.
Finally, when the parser cannot identify an input token it calls a particular function to handle this. By default, this is the internal function "__unimplemented__" which makes suggestions about possible Octave substitutes for matlab functions.
```

val = missing_function_hook ()
old_val = missing_function_hook (new_val)
missing_function_hook (new_val, "local")

```

Query or set the internal variable that specifies the function to call to provide extra information when an unknown identifier is referenced.

When called from inside a function with the "local" option, the variable is changed locally for the function and any subroutines it calls. The original variable value is restored when exiting the function.

See also: [missing_component_hook], page 1008.

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\section*{Concept Index}
A
acknowledgements .....  1
addition ..... 161, 900
and operator ..... 166, 900
anonymous functions ..... 233
ans ..... 137
answers, incorrect ..... 1079, 1081
application-defined data ..... 494
apply ..... 631
area series ..... 500
arguments in function call ..... 157
arithmetic operators ..... 161, 900
array, creating a Java array ..... 1054
assignment expressions ..... 169
assignment operators ..... 169
axes graphics object ..... 423

\section*{B}
bar series ..... 501
batch processing ..... 37
block comments ..... 39
body of a loop ..... 183
boolean expressions ..... 166, 900
boolean operators ..... 166, 900
break statement ..... 186
broadcast ..... 627
broadcasting ..... 627
BSX . ..... 627
bug criteria ..... 1079
bug tracker ..... 1080
bugs. ..... 1079
bugs, investigating ..... 1081
bugs, known ..... 1079
bugs, reporting ..... 1080
built-in data types ..... 41
C
callbacks ..... 492
calling Java from Octave ..... 1049
calling Octave from Java ..... 1049
case statement ..... 181
catch. ..... 188
cell arrays ..... 46, 124
character strings ..... 45, 71
Cholesky factorization ..... 595
Citations .....  5
Citing Octave .....  5
classes, making available to Octave ..... 1050
classpath, adding new path ..... 1057
classpath, difference between
static and dynamic ..... 1050
classpath, displaying ..... 1057
classpath, dynamic ..... 1057
classpath, removing path ..... 1057
classpath, setting ..... 1050
classpath, static ..... 1057
classpath.txt ..... 1050
clearing the screen ..... 26
code profiling ..... 264
colors, graphics ..... 491
comma separated lists ..... 134
command and output logs ..... 34
command completion ..... 28
command descriptions ..... 13
command echoing ..... 34
command history ..... 28
command options ..... 15
command-line editing ..... 25
commands functions ..... 236
comments ..... 38
comparison expressions ..... 164, 900
complex-conjugate transpose ..... 161, 900
containers ..... 111
continuation lines ..... 189
continue statement ..... 187
contour series ..... 502
contributing to Octave ..... 6
contributors ..... 1
conversion specifications (printf) ..... 299
conversion specifications (scanf) ..... 305
copy-on-write ..... 638
copyright ..... 1101
core dump ..... 1079
COW ..... 638
creating graphics objects ..... 423
cs-lists ..... 134
customizing readline ..... 32
customizing the prompt ..... 33
D
DAE ..... 719
data sources in object groups ..... 500
data structures ..... 46, 111
data types ..... 41
data types, built-in ..... 41
data types, user-defined ..... 46
decrement operator ..... 171
default arguments ..... 202
default graphics properties ..... 489
defining functions ..... 191
Degree Symbol ..... 408
deprecated functions ..... 1073
description format ..... 12
diagonal and permutation matrices ..... 651
diagonal matrix expressions ..... 654
dialog, displaying a dialog for selecting directories................................ . 915
dialog, displaying a dialog for selecting files.... 915 ..... 915
dialog, displaying a dialog for storing files
dialog, displaying a font selection dialog ..... 921
dialog, displaying a help dialog ..... 917
dialog, displaying a list dialog ..... 918
dialog, displaying a message dialog ..... 919
dialog, displaying a modal dialog ..... 922
dialog, displaying a question dialog ..... 920
dialog, displaying a warning dialog ..... 921
dialog, displaying an error dialog ..... 917
dialog, displaying an input dialog ..... 918
diary of commands and output ..... 34
differential equations ..... 719
diffs, submitting ..... 1081
distribution of Octave ..... 6
division ..... 161, 900
do-until statement ..... 184
documentation fonts ..... 11
documentation notation ..... 11
documenting functions ..... 39
documenting Octave programs ..... 38
documenting user scripts ..... 39
Dulmage-Mendelsohn decomposition ..... 678
dynamic classpath ..... 1050, 1057
dynamic classpath, adding new path ..... 1057
dynamic naming ..... 116
dynamic-linking ..... 1009
Dynamically Linked Functions ..... 1009
E
echoing executing commands ..... 34
editing the command line ..... 25
element-by-element evaluation ..... 166
else statement ..... 179
elseif statement ..... 179
end statement ..... 179
end, indexing ..... 153
end: and :end ..... 153
end_try_catch ..... 188
end_unwind_protect ..... 188
endfor statement ..... 184
endfunction statement ..... 191
endif statement ..... 179
endswitch statement ..... 181
endwhile statement ..... 183
equality operator ..... 164, 900
equality, tests for ..... 164, 900
equations, nonlinear ..... 641
erroneous messages ..... 1079
erroneous results ..... 1079, 1081
error bar series ..... 503
error ids ..... 244
error message notation ..... 12
error messages ..... 36
error messages, incorrect ..... 1079
escape sequence notation ..... 71
evaluation notation ..... 11
executable scripts ..... 37
execution speed ..... 638
exiting octave ..... 7, 19
exponentiation ..... 161, 900
expression, range ..... 56
expressions ..... 151
expressions, assignment ..... 169
expressions, boolean ..... 166, 900
expressions, comparison ..... 164, 900
expressions, logical ..... 166, 900
F
factorial function ..... 159
fatal signal ..... 1079
field, returning value of Java object field ..... 1055
field, setting value of Java object field ..... 1056
fields, displaying available fields
of a Java object ..... 1055
figure deletefcn ..... 160
figure graphics object ..... 423
finding minimums ..... 645
finish.m ..... 19
flag character (printf) ..... 301
flag character (scanf) ..... 306
for statement ..... 184
Frobenius norm ..... 592
function application. ..... 631
function descriptions ..... 12
function file. ..... 210
function handle ..... 160
function statement ..... 191
functions, deprecated ..... 1073
functions, obsolete ..... 1073
functions, user-defined ..... 191
funding Octave development ..... 6
G
general p-norm ..... 592
global statement ..... 139
global variables ..... 139
GNUTERM ..... 508
grammar rules ..... 1099
graphics ..... 315
graphics colors ..... 491
graphics data structures ..... 421
graphics line styles ..... 492
graphics marker styles ..... 492
graphics object properties ..... 433
graphics object, axes. ..... 423
graphics object, figure ..... 423
graphics object, image ..... 423
graphics object, light ..... 423
graphics object, line ..... 423
graphics object, patch ..... 423
graphics object, root ..... 423
graphics object, surface ..... 423
graphics object, text ..... 423
graphics objects ..... 423
graphics objects, saving ..... 432
graphics properties, default ..... 489
graphics toolkits ..... 507
greater than operator ..... 164, 900
group objects ..... 504, 505, 506
H
handle functions ..... 427
handle, function handles ..... 233
hash table ..... 124
help, online ..... 21
help, user-defined functions ..... 39
help, where to find ..... 1082
Hermitian operator ..... 161, 900
Hessenberg decomposition ..... 597
history ..... 1
history of commands ..... 28
I
if statement............................................. . 179
image graphics object ..... 423
improving Octave ..... 1080, 1081
incorrect error messages ..... 1079
incorrect output ..... 1079, 1081
incorrect results ..... 1079, 1081
increment operator ..... 171
indirect function call ..... 160
infinity norm ..... 592
initialization ..... 18
input conversions, for scanf ..... 306
input history ..... 28
~/.inputrc ..... 32
installation trouble ..... 1079
installing Octave ..... 1085
instance, creating a Java instance ..... 1054
introduction ..... 7
introduction to graphics structures ..... 421
invalid input ..... 1079
J
Java, calling from Octave ..... 1049
Java, using with Octave. ..... 1049
javaclasspath.txt ..... 1050
Jupyter Notebooks ..... 232
K
Kendall's Tau ..... 788
key/value store ..... 124
keywords ..... 1099
known causes of trouble ..... 1079
L
language definition ..... 1099
less than operator ..... 164, 900
light graphics object ..... 423
line graphics object ..... 423
line series ..... 504
line styles, graphics ..... 492
linear algebra ..... 585
linear algebra, techniques ..... 585
loading data ..... 278
local minimum ..... 645
logging commands and output .....  34
logical expressions ..... 166, 900
logical operators ..... 166, 900
loop ..... 183
looping over structure elements ..... 185
LP ..... 741
LU decomposition ..... 597
lvalue ..... 169
M
Map ..... 124
map ..... 631
marker styles, graphics ..... 492
matching failure, in scanf ..... 305
matrices ..... 52
matrices, diagonal and permutation ..... 651
matrix factorizations ..... 595
matrix functions, basic ..... 585
matrix multiplication ..... 161, 900
matrix, functions of ..... 609
matrix, permutation functions ..... 656
matrix, specialized solvers ..... 611
matrix, zero elements ..... 657
maximum field width (scanf) ..... 306
memory management ..... 638
memory, displaying Java memory status ..... 1059
memory, limitations on JVM ..... 1053
messages, error ..... 36
method, invoking a method of
a Java object ..... 1056
methods, displaying available methods of a Java object ..... 1056
mex ..... 1034
mex-files ..... 1034
minimum field width (printf) ..... 301
missing data ..... 45
mkoctfile ..... 1010
multi-line comments ..... 39
multiplication ..... 161, 900
N
negation ..... 161, 900
NLP ..... 741
nonlinear equations ..... 641
nonlinear programming ..... 741
not operator ..... 166, 900
numeric constant ..... 44, 51
numeric value ..... 44, 51
O
object groups ..... 495
object, creating a Java object ..... 1054
obsolete functions ..... 1073
oct ..... 1010
oct-files. ..... 1010
Octave and matlab, how to distinguish between ..... 1082
Octave API ..... 1009
Octave command options ..... 15
Octave, calling from Java ..... 1049
. octaverc ..... 16, 19
~/.octaverc ..... 16, 19
ODE ..... 719
online help ..... 21
opengl rendering slow windows. ..... 508
opengl single precision date time ..... 509
operator precedence ..... 172
operators, arithmetic ..... 161, 900
operators, assignment ..... 169
operators, boolean ..... 166, 900
operators, decrement ..... 171
operators, increment ..... 171
operators, logical ..... 166, 900
operators, relational ..... 164, 900
optimization ..... 638, 741
options, Octave command ..... 15
--braindead ..... 17
--built-in-docstrings-file filename ..... 15
--debug ..... 15
--doc-cache-file filename ..... 15
--echo-commands ..... 15
--eval code. ..... 15
--exec-path path ..... 15
--gui ..... 15
--help ..... 16
--image-path path ..... 16
--info-file filename ..... 16
--info-program program ..... 16
--interactive ..... 16
--line-editing ..... 16
--no-gui ..... 16
--no-history ..... 16
--no-init-file ..... 16
--no-init-path ..... 16
--no-line-editing ..... 16
--no-site-file ..... 16
--no-window-system ..... 16
--norc ..... 16
--path path ..... 17
--persist ..... 17
--quiet ..... 17
--silent ..... 17
--texi-macros-file filename ..... 17
--traditional ..... 17
--verbose ..... 17
--version ..... 18
-d. ..... 15
-f. ..... 16
-h. ..... 16
-H. ..... 16
-i. ..... 16
-p path ..... 17
-q. ..... 17
-v. ..... 18
-V. ..... 17
-W. ..... 16
-x ..... 15
or operator ..... 166, 900
oregonator ..... 721
otherwise statement ..... 181
output conversions, for printf ..... 301
P
parser ..... 1099
patch graphics object ..... 423
patches, submitting ..... 1081
path, adding to classpath ..... 1057
path, removing from classpath ..... 1057
permutation matrix functions ..... 656
persistent statement ..... 140
persistent variables ..... 140
personal startup file ..... 19
PKG_ADD ..... 1004
PKG_DEL ..... 1004
plotting. ..... 315
plotting, high-level ..... 315
plotting, multiple plot windows ..... 399
plotting, multiple plots per figure ..... 398
plotting, object manipulation ..... 400
plotting, saving and printing plots ..... 409
plotting, three-dimensional ..... 357
plotting, two-dimensional functions ..... 353
plotting, window manipulation ..... 401
precision (printf) ..... 301
printing notation. ..... 11
printing plots ..... 409
profiler ..... 264
program, self contained ..... 37
Progress Bar ..... 923
project startup file ..... 19
prompt customization ..... 33
pseudoinverse. ..... 593, 654
Q
QP ..... 741
QR factorization ..... 599
quadratic programming ..... 741
quitting octave ..... 7, 19
quiver group ..... 504
quotient ..... 161, 900
R
range expressions ..... 56
readline customization ..... 32
recycling. ..... 627
relational operators ..... 164, 900
reporting bugs ..... 1079, 1080
results, incorrect ..... 1079, 1081
root graphics object ..... 423
S
saving data ..... 278
saving graphics objects ..... 432
saving plots ..... 409
Schur decomposition ..... 604
script files ..... 191
scripts ..... 37
select JVM version ..... 1053
self contained programs ..... 37
series objects ..... 506
short-circuit evaluation ..... 167
side effect ..... 169
SIMD ..... 627
singular value decomposition ..... 607
site exiting file ..... 19
site startup file ..... 16, 18
Spearman's Rho ..... 788
speedups ..... 638
stair group ..... 505
startup ..... 18
startup files ..... 18
startup.m ..... 19
statements ..... 179
static classpath ..... 1050, 1057
stem series ..... 506
strings ..... 45, 71
structural rank ..... 685
structure elements, looping over ..... 185
structures. ..... 46, 111
submitting diffs ..... 1081
submitting patches ..... 1081
subtraction ..... 161, 900
suggestions ..... 1080
surface graphics object ..... 423
surface group ..... 506
switch statement ..... 181
T
test functions ..... 1061
tests for equality ..... 164, 900
text graphics object ..... 423
toolkit customization ..... 508
toolkits, graphics ..... 507
transform groups ..... 507
transpose ..... 161, 900
transpose, complex-conjugate. ..... 161, 900
troubleshooting ..... 1079
try statement ..... 188
U
uitable properties ..... 481
unary minus ..... 161, 900
undefined behavior ..... 1079
undefined function value ..... 1079
unwind_protect statement ..... 188
unwind_protect_cleanup ..... 188
use of comments ..... 38
user-defined data types ..... 46
user-defined functions ..... 191
user-defined variables ..... 137
using Octave with Java ..... 1049
V
validating arguments ..... 202
varargin ..... 199
varargout ..... 198
variable-length argument lists ..... 199
variable-length return lists ..... 198
variables, global ..... 139
variables, persistent ..... 140
variables, user-defined ..... 137
vectorization ..... 625
vectorize ..... 625
version startup file ..... 19
W
warning ids ..... 248
warranty ..... 1101
while statement ..... 183
wrong answers ..... 1079, 1081

\section*{Function Index}

\section*{A}
audioinfo ..... 881
abs............................................................ . 549

accumdim................................................... 637
acos............................................................ 551
acosd......................................................... 554
acosh........................................................... 552
acot......................................................... 552
acotd......................................................... 554
acoth..................................................... 553
acsc......................................................... 552
acscd............................................................. 554
acsch......................................................... 553
add_input_event_hook............................. . . . 1099
addlistener............................................... 497
addpath................................................... 213
addpref.................................................... . . . 937
addproperty............................................... . . 496
addtodate................................................. . . . . . 951
airy.......................................................... . . . . . . . . . . . . . 565
all.......................................................... . 511
allchild.................................................... 431
amd.......................................................... . . . . 675
ancestor................................................... . 431
and............................................................... . . 166
angle........................................................ . . . . 549
annotation.................................................... . 396
any............................................................. . . 511
arch_fit................................................... . 850
arch_rnd..................................................... . . 851
arch_test.................................................. . 851
area................................................................ 346
arg............................................................ . . 549
argv.......................................................... . . 18
arma_rnd................................................... . . . 851
arrayfun.................................................. . . 631
ascii........................................................... 967
asctime...................................................... 943
asec............................................................ 551
asecd.......................................................... . . 554
asech........................................................ 552
asin........................................................... . 551
asind.......................................................... 554
asinh......................................................... 552
assert..................................................... . 1068
assignin..................................................... 178
atan......................................................... . 551
atan2......................................................... . . 553
atan2d........................................................ . 554
atand......................................................... 554
atanh......................................................... . . 552
atexit........................................................ 20
audiodevinfo........................................... 882
audioformats.......................................... 882
audioplayer ..... 883
audioread ..... 881
audiorecorder ..... 884
audiowrite ..... 882
auto_repeat_debug_command ..... 263
autoload ..... 221
autoreg_matrix ..... 852
autumn ..... 872
available_graphics_toolkits ..... 508
axes ..... 424
axis ..... 349

\section*{B}
balance..................................................... . . 585
bandwidth................................................ . . . 586
bar............................................................. . 321
barh........................................................... 322
bartlett................................................... . . 852
base2dec.................................................. . 100
base64_decode.......................................... . . 971
base64_encode.......................................... . . . 971
beep........................................................ . 241
beep_on_error .......................................... . . . . 241
besselh................................................... . . . 568
besseli....................................................... . 567
besselj...................................................... 566
besselk..................................................... . . 568
bessely..................................................... . . . . . . . . . 566
beta.......................................................... . 569
betainc..................................................... . . 569
betaincinv................................................... . . 569
betaln........................................................ . . 570
bicg......................................................... 611
bicgstab................................................. 613
bin2dec.......................................................... 98
binary...................................................... . . 967
bincoeff..................................................... . . 570
bitand.......................................................... 63
bitcmp........................................................... 63
bitget........................................................ 62
bitor.......................................................... 63
bitpack...................................................... 43
bitset......................................................... 62
bitshift..................................................... 64
bitunpack..................................................... . 44
bitxor........................................................ 63
blackman.................................................... . . 852
blanks.......................................................... . 71
blkdiag...................................................... . . 525
blkmm......................................................... . 610
bone........................................................ . . . 872
bounds ....................................................... . . . 759
box............................................................ 394
brighten................................................. . 876 ..... 876

clock....................................................... 945
bsxfun ..... 628
built_in_docstrings_file ..... 24
builtin ..... 220
bunzip2 ..... 963
byte_size ..... 1015
bzip2. ..... 965
Ccalendar.................................................. 951
camlight ..... 372
camlookat ..... 376
camorbit ..... 377
campos ..... 376
camroll ..... 377
camtarget ..... 378
camup ..... 379
camva. ..... 379
camzoom ..... 380
canonicalize_file_name ..... 962
cart2pol ..... 578
cart2sph ..... 579
cast. ..... 42
cat ..... 517
caxis ..... 350
cbrt ..... 548
ccolamd ..... 675
cd. ..... 967, 982
ceil. ..... 557
cell. ..... 127
cell2mat ..... 134
cell2struct ..... 134
celldisp ..... 125
cellfun ..... 633
cellindexmat ..... 132
cellslices ..... 130
cellstr ..... 133
center ..... 784
cgs ..... 615
char. ..... 78
chdir ..... 982
chol. ..... 595
chol2inv ..... 596
choldelete ..... 596
cholinsert ..... 596
cholinv ..... 595
cholshift ..... 597
cholupdate ..... 596
circshift ..... 519
citation .....  5
cla. ..... 404
clabel ..... 393
class. ..... 41
clc ..... 27
clear ..... 146
clearvars ..... 147
clf ..... 404
close............................................... . 405, 966
closereq.................................................. . . 405
cmpermute.................................................. . . . . . . . . . . . 878
cmunique ................................................. . . 877
colamd........................................................ . 676
colloc.................................................... 712
colon....................................................... . . 898
colorbar.......................................... . 394, 395
colorcube.................................................. . . 872

colperm................................................... . . . 677
colstyle................................................... . 492
columns.................................................... 46
comet......................................................... . 348
comet3.................................................. . . 348
command_line_path..................................... . . 216
commandhistory ...................................... 939
commandwindow.......................................... . 939
common_size............................................... . . 513
commutation_matrix................................... . 570
compan................................................. 798
compare_versions...................................... . . 989
compass................................................... 344
completion_append_char............................. . 28
completion_matches..................................... 28
complex...................................................... 52
computer.................................................. . . 985
cond.......................................................... 586

condest.................................................. . . 683
confirm_recursive_rmdir.......................... . . . 955
conj.......................................................... 549
containers.Map......................................... . . 124
contour..................................................... . . . . 330
contour3.................................................... . 332
contourc.................................................. . . 331
contourf.................................................. . . . 331
contrast................................................... . 876
conv.......................................................... . . 799
conv2..................................................... . . . 800
convhull................................................... . . 836
convhulln................................................ . 837
convn....................................................... . 799
cool......................................................... 873
copper.................................................... . . 873
copyfile................................................... . . . 953
copyobj................................................. 432
corr......................................................... . 787
corrcoef................................................. . . 787
cos........................................................... 551
cosd........................................................... . 553
cosh............................................................ 552
cosint.................................................... . 571
cospi............................................................ 555
cot............................................................ 551
cotd......................................................... . 554
coth........................................................ 552
cov........................................................ 786 dec2base ..... 100
cplxpair ....................................................... 549
cputime946
crash_dumps_octave_core ..... 292
cross ..... 561
csc ..... 551
cscd. ..... 554
csch. ..... 552
cstrcat ..... 80
csvread ..... 285
csvwrite ..... 285
csymamd ..... 677
ctime ..... 942
ctranspose ..... 162
cubehelix ..... 873
cummax ..... 559
cummin ..... 560
cumprod ..... 557
cumsum ..... 556
cumtrapz ..... 711
curl ..... 562
cylinder ..... 388
D
daspect ..... 384
daspk ..... 722
daspk_options ..... 723
dasrt ..... 727
dasrt_options ..... 729
dassl ..... 725
dassl_options ..... 726
date. ..... 946
datenum ..... 948
datestr ..... 949
datetick ..... 952
datevec ..... 951
dawson ..... 571
dbclear ..... 260, 261
dbcont ..... 258
dbdown ..... 264
dblist ..... 262
dblquad ..... 713
dbnext ..... 263
dbquit ..... 258
dbstack ..... 263
dbstatus ..... 260
dbstep ..... 263
dbstop ..... 258
dbtype ..... 262
dec2bin ..... 99
dec2hex ..... 99
decic. ..... 736
deconv ..... 800
deg2rad ..... 550
del2 ..... 562
delaunay ..... 825
delaunayn ..... 826
delete ..... 405, 967
dellistener ..... 497
demo ..... 1069
det ..... 587
detrend ..... 852
diag ..... 525
dialog ..... 922
diary ..... 14, 35
diff ..... 512
diffpara ..... 852
diffuse ..... 370
dims ..... 1016
dir ..... 967, 983
dir_encoding ..... 216
dir_in_loadpath ..... 216
discrete_cdf ..... 789
discrete_inv ..... 789
discrete_pdf ..... 789
discrete_rnd ..... 789
disp ..... 271
display ..... 903
divergence ..... 561
dlmread ..... 284
dlmwrite ..... 284
dmperm ..... 678
do_string_escapes ..... 76
doc ..... 21
doc_cache_create ..... 24
doc_cache_file ..... 24
dos ..... 973
dot ..... 561
double ..... 51
drawnow ..... 401
dsearch ..... 831
dsearchn ..... 832
dup2 ..... 977
duplication_matrix ..... 571
durbinlevinson ..... 853
dbup. ..... 264
dbwhere ..... 262
deal. ..... 198
deblank ..... 75
debug_java ..... 1060
debug_on_error ..... 257
debug_on_interrupt ..... 257
debug_on_warning ..... 257

e........................................................... 580

factor.................................................... 563
echo.......................................................... . 35
edit ..... 211
edit_history ..... 29
EDITOR ..... 32
eig ..... 587
elem. ..... 1015
ellipj. ..... 571
ellipsoid ..... 389
empirical_cdf ..... 789
empirical_inv ..... 789
empirical_pdf ..... 789
empirical_rnd ..... 790
endgrent ..... 985 ..... 985
endpwent ..... 984
endsWith ..... 86
eomday ..... 952
eps ..... 582
eq. ..... 165
erase ..... 90
erf ..... 572
erfc ..... 572
erfcinv ..... 573
erfcx ..... 572
erfi. ..... 573
erfinv ..... 573
errno ..... 245
errno_list ..... 245
error ..... 239
errorbar ..... 333
errordlg ..... 917
etime. ..... 946
etree ..... 669
etreeplot ..... 669
eval. ..... 175
evalc. ..... 175
evalin ..... 178
example ..... 1070
exec ..... 976
EXEC_PATH ..... 976
exist ..... 145
exp ..... 547
expint ..... 573
expm. ..... 609
expm1 ..... 547
eye ..... 525
ezcontour ..... 355
ezcontourf ..... 355
ezmesh ..... 385
ezmeshc ..... 386
ezplot ..... 354
ezplot3 ..... 384
ezpolar ..... 356
ezsurf ..... 386
ezsurfc ..... 387
factorial.................................................. . . 563
fail.......................................................... 1068
false............................................................. 65
fclear.................................................... . . 313
fclose...................................................... 296
fcntl........................................................ . . 979
fdisp........................................................... 283
feather..................................................... . . 345
feof............................................................. 312
ferror...................................................... . . . 313
feval............................................................. 176
fflush..................................................... . . 276
fft............................................................... 845
fft2......................................................... . . . 845
fftconv..................................................... . . 848
fftfilt...................................................... 848
fftn........................................................... 846
fftshift........................................................ . 853
fftw............................................................. . . 847
fgetl........................................................ . . 297
fgets........................................................... . 298
fieldnames................................................. . . . 119
figure............................................... 399, 400
file_in_loadpath...................................... . . 215
file_in_path ............................................ . 960
fileattrib.................................................. 958
filebrowser.................................................... 939
fileparts.................................................. . 961
fileread................................................... . . 283
filesep................................................... . . . . . . . . . 960
fill........................................................... 346
fill3.......................................................... 347
filter...................................................... . . 848
filter2......................................................... 849
find........................................................ . . . 513
findall.................................................... . . 489
findfigs................................................... . . 431
findobj.................................................... 488
findstr..................................................... . . 87
fix........................................................... 557
fixed_point_format...................................... . . . 55
flag........................................................... . . 873
flintmax.................................................... 60
flip............................................................. 515
fliplr...................................................... . . 515
flipud....................................................... . . . . . . . . . . . . . 515
floor......................................................... . 557
fminbnd...................................................... . . . . . 645
fminsearch.............................................. . . 647
fminunc....................................................... 646
foo................................................................ . 13
fopen........................................................ . . . 295
fork............................................................ . . . 976
format....................................................... 272
fortran_vec............................................. . . . 1016
fplot........................................................ 353
fprintf................................................... 299 getppid ..... 980
fputs. ..... 297
fractdiff ..... 853
frame2im ..... 870
fread ..... 308
freport ..... 313
freqz. ..... 849
freqz_plot ..... 850
frewind ..... 314
fscanf ..... 304
fseek. ..... 313
fskipl ..... 298
fsolve ..... 641
ftell. ..... 313
ftp ..... 966
full. ..... 664
fullfile ..... 961
func2str ..... 234
functions ..... 234
fwrite ..... 310
fzero ..... 644
G
gallery ............... 537, 538, 539, 540, 541, 542
gamma ..... 574
gammainc ..... 574
gammaincinv ..... 574
gammaln ..... 576
gca ..... 429
gcbf. ..... 494
gcbo ..... 494
gcd. ..... 563
gcf. ..... 428
gco. ..... 429
ge. ..... 165
genpath ..... 214
get ..... 430, 884, 886
get_first_help_sentence ..... 25
get_help_text ..... 25
get_help_text_from_file ..... 25
get_home_directory ..... 981
getappdata ..... 495
getaudiodata ..... 885
getegid ..... 981
getenv ..... 981
geteuid ..... 980
getfield ..... 121
getframe ..... 869
getgid ..... 981
getgrent ..... 985
getgrgid ..... 985
getgrnam ..... 985
gethostname ..... 966
getpgrp ..... 980
getpid ..... 980
getpixelposition ..... 934
getplayer ..... 885
getpref ..... 937
getpwent ..... 984
getpwnam ..... 984
getpwuid ..... 984
getrusage ..... 992
getuid ..... 981
ginput ..... 419
givens ..... 588
glob. ..... 959
glpk. ..... 741
gls. ..... 752
gmres ..... 617
gmtime ..... 942
gnuplot_binary ..... 508
gplot ..... 669
grabcode ..... 227
gradient ..... 560
graphics_toolkit ..... 507
gray ..... 873
gray2ind ..... 868
grid. ..... 394
griddata ..... 838
griddata3 ..... 839
griddatan ..... 839
groot. ..... 428
gsvd. ..... 589
gt ..... 165
gtext. ..... 420
guidata ..... 933
guihandles ..... 934
gunzip ..... 963
gzip. ..... 963
H
hadamard ..... 542
hamming ..... 853
hankel ..... 542
hanning ..... 853
hash. ..... 994
have_window_system ..... 934
hdl2struct ..... 432
help. ..... 21
helpdlg ..... 917
hess. ..... 597
hex2dec ..... 100
hex2num ..... 101
hggroup ..... 495
hgload ..... 418
hgsave ..... 417
hgtransform ..... 507
hidden ..... 360
hilb. ..... 543
hist. ..... 323
histc. ..... 784
history ..... 29
history_control ..... 31
history_file ..... 31
history_save ..... 30
history_size. ..... 32
history_timestamp_format_string ..... 32
hold. ..... 403
home ..... 27
horzcat ..... 517
hot ..... 874
housh. ..... 608
hsv ..... 874
hsv2rgb ..... 879
humps ..... 649
hurst. ..... 854
hypot. ..... 560
I
581i....
695
idivide ..... 61
ifelse ..... 168
ifft. ..... 845
ifft2. ..... 846
ifftn. ..... 846
ifftshift ..... 853
ignore_function_time_stamp ..... 213
ilu ..... 696, 697
im2double ..... 867
im2frame ..... 870
imag.. ..... 550
image. ..... 866
IMAGE_PATH. ..... 861
imagesc. ..... 866
imfinfo. ..... 862
imformats ..... 864
importdata. ..... 291
imread ..... 859
imshow ..... 865
imwrite. ..... 860
ind2gray ..... 868
ind2rgb ..... 868
ind2sub ..... 156
index. ..... 87
Inf ..... 581
inf ..... 581
inferiorto. ..... 901
info. ..... 22
info_file ..... 22
info_program. ..... 23
inpolygon. ..... 835
input. ..... 277
inputdlg ..... 918
inputname ..... 194
inputParser ..... 208
inputParser.CaseSensitive. ..... 209
inputParser.FunctionName ..... 209
inputParser.KeepUnmatched. ..... 209
inputParser.Parameters ..... 208
inputParser.Results ..... 209
inputParser.StructExpand ..... 209
inputParser.Unmatched ..... 209
inputParser.UsingDefaults ..... 209
int16 ..... 58
int2str ..... 97
int32 ..... 59
int64 ..... 59
int8. ..... 58
integral ..... 709
integral2 ..... 715
integral3 ..... 717
interp1 ..... 815
interp2 ..... 819
interp3 ..... 820
interpft ..... 817
interpn ..... 821
intersect ..... 793
intmax. ..... 59
intmin ..... 60
inv ..... 590
inverse ..... 590
invhilb ..... 543
ipermute ..... 518
iqr ..... 759
is_absolute_filename ..... 962
is_dq_string. ..... 72
is_function_handle. ..... 234
is_leap_year ..... 946
is_rooted_relative_filename ..... 962
is_same_file ..... 962
is_sq_string ..... 72
is_valid_file_id ..... 296
isa ..... 41
isalnum ..... 108
isalpha ..... 108
isappdata ..... 495
isargout ..... 201
isascii ..... 110
isaxes ..... 427
isbanded ..... 69
isbool. ..... 67
iscell ..... 126
iscellstr ..... 133
ischar. ..... 73
iscntrl ..... 109
iscolormap ..... 871
iscolumn ..... 68
iscomplex ..... 67
isdebugmode ..... 263
isdefinite ..... 69
isdeployed ..... 987
isdiag ..... 70
isdigit ..... 109
isdir. ..... 959
isempty ..... 48
isequal ..... 165
isequaln ..... 165
isfield................................................... 120
isfigure ..... 427
isfile ..... 959
isfinite ..... 513
isfloat ..... 67
isfolder ..... 959
isglobal ..... 140
isgraph ..... 109
isgraphics ..... 427
isguirunning ..... 934
ishandle ..... 427
ishermitian ..... 69
ishghandle. ..... 427
ishold ..... 403
isieee ..... 987
isindex ..... 157
isinf. ..... 512
isinteger ..... 58
isjava ..... 1055
iskeyword ..... 1099
isletter ..... 108
islogical ..... 67
islower ..... 109
ismac. ..... 986
ismatrix ..... 68
ismember ..... 795
ismethod ..... 891
isna. ..... 45
isnan. ..... 512
isnull. ..... 48
isnumeric ..... 67
isobject ..... 890
isocaps ..... 366
isocolors ..... 367
isonormals ..... 365
isosurface. ..... 363
ispc. ..... 986
isplaying ..... 884
ispref ..... 938
isprime ..... 70
isprint ..... 110
isprop ..... 423
ispunct. ..... 109
isreal. ..... 67
isrecording ..... 885
isrow. ..... 68
isscalar ..... 68
issorted ..... 522
isspace ..... 109
issparse ..... 666
issquare ..... 68
isstring ..... 73
isstrprop ..... 110
isstruct ..... 119
isstudent ..... 987
issymmetric ..... 68
istril. ..... 70
istriu ..... 70
isunix
isunix ..... 986 ..... 986
isupper ..... 109
isvarname ..... 137
isvector ..... 68
isxdigit ..... 109
I. ..... 581
J
j. ..... 581
java_get ..... 1055
java_matrix_autoconversion ..... 1059
java_set ..... 1056
java_unsigned_autoconversion ..... 1059
javaaddpath ..... 1057
javaArray ..... 1054
javachk ..... 1058
javaclasspath ..... 1057
javamem ..... 1059
javaMethod ..... 1056
javaObject ..... 1054
javarmpath ..... 1057
jet ..... 874
jsondecode ..... 106
jsonencode ..... 104
jupyter_notebook ..... 232
J. ..... 581
K
kbhit. ..... 278
kendall ..... 788
keyboard ..... 261
kill. ..... 980
kron. ..... 610
krylov ..... 609
kurtosis ..... 763
L
lasterr ..... 243
lasterror ..... 242
lastwarn ..... 248
1 cm ..... 563
ldivide ..... 162
le. ..... 165
legend ..... 390
legendre ..... 575
length ..... 47
lgamma ..... 576
license ..... 989, 990
light. ..... 426
lightangle ..... 373
lighting. ..... 371
lin2mu ..... 886
line. ..... 424
lines. ..... 874
link ..... 954
linkaxes...................................................... 499 menu ..... 277
linkprop..................................................... . 498
linsolve ..... 590
linspace ..... 530
list_in_columns ..... 271
list_primes ..... 565
listdlg ..... 918
listfonts ..... 934
load ..... 281, 282
loaded_graphics_toolkits ..... 508
loadobj ..... 894
localfunctions ..... 217
localtime ..... 942
log ..... 547
log10 ..... 547
\(\log 1 p\) ..... 547
log2 ..... 547
logical ..... 65
loglog ..... 320
loglogerr ..... 336
logm ..... 610
logspace ..... 530
lookfor ..... 21, 22
lookup ..... 514
lower ..... 75
ls ..... 982
ls_command ..... 982
lscov ..... 753
lsode ..... 719
lsode_options ..... 720
lsqnonneg ..... 752
lstat ..... 956
lt ..... 165
lu ..... 597
luupdate ..... 598
M
mad ..... 760
magic ..... 543
make_absolute_filename ..... 962
makeinfo_program ..... 23
mat2cell ..... 128
mat2str ..... 95
material ..... 371
matlab.lang.makeUniqueStrings ..... 138
matlab.lang.makeValidName ..... 137
matlab.net.base64decode ..... 972
matlab.net.base64encode ..... 971
matlabroot ..... 987
matrix_type ..... 591
\(\max\) ..... 558
max_recursion_depth ..... 159
max_stack_depth ..... 160
mean ..... 757
meansq ..... 760
median ..... 758
memory ..... 990
mesh. ..... 358
meshc ..... 359
meshgrid ..... 373
meshz ..... 359
methods ..... 890, 891
mex ..... 1034
mexext ..... 1034
mfilename ..... 212
mget ..... 966
mgorth ..... 593
min ..... 559
minus ..... 162
mislocked ..... 222
missing_component_hook ..... 1008
missing_function_hook ..... 1100
mkdir ..... 954, 967
mkfifo ..... 955
mkoctfile ..... 1010
mkpp ..... 811
mkstemp ..... 311
mktime ..... 943
mldivide ..... 162
mlock ..... 222
mod ..... 564
mode ..... 758
moment ..... 763
more ..... 275
movefile ..... 953
movegui ..... 935
movfun ..... 766
movie ..... 869
movmad ..... 768
movmax ..... 770
movmean ..... 772
movmedian ..... 773
movmin ..... 775
movprod ..... 776
movslice ..... 768
movstd ..... 778
movsum ..... 780
movvar ..... 782
mpoles ..... 799
mpower ..... 163
mput ..... 967
mrdivide ..... 163
msgbox ..... 919
mtimes ..... 163
mu2lin ..... 887
munlock ..... 222
mustBeFinite ..... 206
mustBeGreaterThan ..... 206
mustBeGreaterThanOrEqual ..... 206
mustBeInteger ..... 206
mustBeLessThan ..... 206
mustBeLessThanOrEqual ..... 207
mustBeMember ..... 207
mustBeNegative......................................... . 207
mustBeNonempty ....................................... . . . 207
mustBeNonNan ........................................... 207
O
ocean. ..... 874
mustBeNonnegative ..... 207
mustBeNonpositive ..... 207
mustBeNonsparse ..... 207
mustBeNonzero ..... 208
mustBeNumeric ..... 208
mustBeNumericOrLogical ..... 208
mustBePositive ..... 208
mustBeReal ..... 208
N
namedargs2cell ..... 124
namelengthmax ..... 138
nan ..... 582
NaN ..... 582
nargin ..... 193
narginchk ..... 203
nargout ..... 197
nargoutchk ..... 203
native_float_format ..... 283
native2unicode ..... 95
NA ..... 45
nchoosek ..... 785
ndgrid ..... 374
ndims ..... 46, 1016
ne. ..... 165
newline ..... 77
newplot ..... 402
news ..... 22
nextpow2 ..... 548
nnz ..... 666
nonzeros ..... 666
norm. ..... 592
normest ..... 682
normest1 ..... 682
not ..... 167
now ..... 941
nproc ..... 986
nth_element ..... 522
nthargout ..... 196
nthroot ..... 548
null. ..... 593
num2cell ..... 127
num2hex ..... 101
num2str ..... 96
numel 46, 1015
numfields ..... 119
nzmax ..... 666
octave_core_file_limit ..... 293
octave_core_file_name ..... 294
octave_core_file_options ..... 293
OCTAVE_EXEC_HOME ..... 987
OCTAVE_HOME ..... 987
OCTAVE_VERSION ..... 987
ode15i ..... 735
ode15s ..... 733, 734
ode23 ..... 731, 732
ode23s ..... 732, 733
ode45 ..... 731
odeget ..... 739
odeplot ..... 739
odeset ..... 737
ols ..... 751
onCleanup ..... 245
ones ..... 526
open ..... 973
openfig ..... 418
openvar ..... 935
operator ..... 1015
optimget ..... 755
optimize_diagonal_matrix ..... 651
optimize_permutation_matrix ..... 651
optimize_range ..... 56
optimize_subsasgn_calls ..... 896
optimset ..... 753
or ..... 167
ordeig ..... 606
orderfields ..... 121
ordqz ..... 605
ordschur ..... 605
orient ..... 417
orth. ..... 593
oruntests ..... 1070
ostreamtube ..... 342
ostrsplit ..... 83
output_precision ..... 54
P
P_tmpdir ..... 312
pack ..... 148
padecoef ..... 810
page_output_immediately ..... 276
page_screen_output ..... 276
PAGER ..... 275
PAGER_FLAGS ..... 275
pan ..... 400
pareto ..... 329
parseparams ..... 200
pascal ..... 544
patch. ..... 425
path. ..... 214
pathdef ..... 215
pathsep ..... 215
pause............................... . 884, 885, 947, 948 print_empty_dimensions ..... 55
pbaspect................................................. 384 ..... 113 ..... 113
pcg ..... 690
pchip. ..... 854
pclose ..... 975
pcolor ..... 345
pcr. ..... 693
peaks ..... 421
periodogram ..... 854
perl. ..... 974
perms ..... 785
permute ..... 518
pi. ..... 580
pie ..... 338
pie3 ..... 338
pink ..... 875
pinv. ..... 593
pipe. ..... 976
pkg ..... 997, 998
planerot ..... 589
play ..... 883, 885
playblocking ..... 883
plot. ..... 316
plot3. ..... 374
plotmatrix ..... 328
plotyy ..... 318
plus. ..... 163
pol2cart ..... 578
polar ..... 337
poly ..... 813
polyaffine ..... 803
polyarea ..... 834
polyder ..... 802
polyeig ..... 798
polyfit ..... 803
polygcd ..... 800
polyint ..... 803
polyout ..... 813
polyreduce ..... 813
polyval. ..... 797
polyvalm. ..... 797
popen ..... 974
popen2 ..... 975
postpad ..... 525
pow2 ..... 548
power ..... 163
powerset ..... 795
ppder ..... 812
ppint ..... 812
ppjumps ..... 812
ppval. ..... 812
print_usage ..... 241
printd ..... 325
printf ..... 298
prism. ..... 875
prod ..... 556
profexplore ..... 266
profexport ..... 266
profile ..... 265
profshow ..... 266
program_invocation_name ..... 18
program_name. ..... 18
properties ..... 909
PS1 ..... 33
PS2 ..... 34
PS4 ..... 34
psi ..... 576
publish ..... 225
putenv ..... 981
puts ..... 297
pwd ..... 983
python ..... 974
Q
qmr ..... 620
qp ..... 747
qr ..... 599
qrdelete ..... 602
qrinsert ..... 601
qrshift ..... 602
qrupdate ..... 601
quad. ..... 704
quad_options ..... 705
quad2d ..... 714
quadcc ..... 708
quadgk ..... 707
quadl ..... 706
quadv ..... 705
quantile ..... 764
questdlg ..... 920
quit ..... 20
quiver ..... 339
quiver3 ..... 339
qz. ..... 602
qzhess ..... 603
pqpnonneg ..... 748
prctile ..... 765
prefdir ..... 938
preferences ..... 938
prepad ..... 524
primes ..... 564
print ..... 410

SEEK_END .................................................. . . 314 specular ..... 370
speed
speed ..... 1071 ..... 1071SEEK_SET ............................................................................................... 319
semilogxerr ..... 335319
semilogy
semilogyerr ..... 336
set........................................ \(430,884,886\)
494
setappdata
794
setenv ..... 981
setfield ..... 120
setgrent ..... 985
setpref ..... 937
setpwent ..... 984
setxor ..... 794
shading ..... 382
shg. ..... 404
shift. ..... 520
shiftdim ..... 520
shrinkfaces ..... 369
sighup_dumps_octave_core. ..... 292
sign ..... 565
signbit ..... 565
sigquit_dumps_octave_core ..... 292
sigterm_dumps_octave_core ..... 293
SIG ..... 980
silent_functions ..... 194
\(\sin\) ..... 551
sinc. ..... 850
sind ..... 553
sinetone ..... 855
sinewave ..... 855
single. ..... 57
sinh. ..... 552
sinint ..... 576
sinpi. ..... 554
size ..... 47
size_equal ..... 49
sizemax ..... 53
sizeof. .....  48
skewness ..... 762
slice. ..... 381
smooth3 ..... 367
sombrero ..... 420
sort ..... 521
sortrows ..... 522
sound. ..... 887
soundsc ..... 887
source ..... 225
spalloc ..... 664
sparse ..... 664
sparse_auto_mutate ..... 671
spaugment ..... 686
spconvert ..... 665
spdiags ..... 661
spearman ..... 788
spectral_adf ..... 855
spectral_xdf ..... 856
spencer ..... 856
speye. ..... 662
spfun. ..... 632
sph2cart ..... 579
sphere ..... 388
spinmap ..... 877
spline ..... 819
splinefit ..... 804
split_long_rows ..... 54
spones ..... 662
spparms ..... 684
sprand ..... 662
sprandn ..... 662
sprandsym ..... 663
sprank ..... 685
spring ..... 875
sprintf ..... 299
spstats ..... 666
spy ..... 668
sqp ..... 749
sqrt ..... 548
sqrtm. ..... 610
squeeze ..... 49
sscanf ..... 305
stairs ..... 325
startsWith ..... 85
stat. ..... 956
statistics ..... 765
std. ..... 760
stderr ..... 294
stdin. ..... 294
stdout ..... 294
stem ..... 325, 326
stem3. ..... 327
stemleaf ..... 324
stft ..... 856
stop ..... 884, 885
str2double ..... 97
str2func ..... 234
str2num ..... 98
strcat. ..... 79
strchr ..... 87
strcmp. ..... 84
strcmpi ..... 85
stream2 ..... 343
stream3 ..... 344
streamline ..... 343
streamribbon ..... 340
streamtube ..... 341
strfind ..... 88
strftime ..... 943
string_fill_char ..... 73
strjoin ..... 84
strjust ..... 74
strmatch ..... 89
strncmp ..... 84
strncmpi..................................................... 85 ..... 945strread
strrep ..... 102
strsplit ..... 81
strtok. ..... 81
strtrim ..... 75
strtrunc ..... 76
struct ..... 118
struct_levels_to_print ..... 113
struct2cell ..... 123
struct2hdl ..... 432
structfun ..... 635
strvcat ..... 79
sub2ind ..... 155
subplot ..... 398
subsasgn ..... 896
subsindex ..... 897
subspace ..... 606
subsref ..... 894
substr ..... 80
substruct ..... 122
sum ..... 555
summer ..... 875
sumsq ..... 557
superiorto ..... 901
suppress_verbose_help_message ..... 24
surf ..... 360
surface ..... 426
surfc ..... 361
surfl ..... 362
surfnorm ..... 363
svd ..... 607
svd_driver ..... 608
svds. ..... 689
swapbytes ..... 43
sylvester ..... 611
symamd ..... 679
symbfact ..... 685
symlink ..... 954
symrcm ..... 680
symvar ..... 235
synthesis ..... 856
system ..... 972
T
tan ..... 551
tand. ..... 553
tanh. ..... 552
tar ..... 964
tempdir ..... 312
tempname ..... 311
terminal_size ..... 272
test. ..... 1061
tetramesh ..... 829
texi_macros_file ..... 23
text ..... 392
unicode_idx............................................. 88 webread ..... 969
unicode2native ..... 94
union ..... 793
unique ..... 791
uniquetol ..... 792
unix. ..... 973
unlink ..... 954
unmkpp ..... 811
unpack ..... 965
unsetenv ..... 981
untabify ..... 76
untar. ..... 964
unwrap ..... 850
unzip ..... 964
uplus ..... 164
upper ..... 75
urlread ..... 968
urlwrite ..... 969
usejava ..... 1058
user_config_dir ..... 987
user_data_dir ..... 987
V
validateattributes. ..... 204
validatestring ..... 203
vander ..... 544
var ..... 761
vec ..... 524
vech ..... 524
vecnorm ..... 594
ver ..... 988
verLessThan ..... 989
version ..... 988
vertcat ..... 517
view. ..... 375
viridis ..... 876
voronoi ..... 832
voronoin ..... 833
W
waitbar ..... 923
waitfor ..... 936
waitforbuttonpress ..... 419
waitpid ..... 977
warndlg ..... 921
warning ..... 246
warranty ..... 22
waterfall ..... 383
WCONTINUE ..... 977
WCOREDUMP ..... 978
web ..... 968
weboptions ..... 970
webwrite ..... 970
weekday ..... 952
WEXITSTATUS ..... 978
what ..... 148
which ..... 148
white. ..... 876
whitebg ..... 877
who ..... 142
whos. ..... 143
whos_line_format ..... 144
WIFCONTINUED ..... 978
WIFEXITED ..... 978
WIFSIGNALED ..... 978
WIFSTOPPED ..... 978
wilkinson ..... 545
winqueryreg ..... 993
winter ..... 876
WNOHANG ..... 979
workspace ..... 939
WSTOPSIG ..... 979
WTERMSIG ..... 979
WUNTRACED ..... 979
X
xlabel ..... 393
xlim. ..... 351
xor ..... 512
xtickangle ..... 352
xticklabels ..... 352
xticks ..... 351
Y
yes_or_no................................................ . . 277
ylabel ..... 393
ylim. ..... 351
ytickangle ..... 352
yticklabels ..... 352
yticks ..... 351
yulewalker ..... 857
Z
zeros ..... 526, 527
zip ..... 964
zlabel ..... 393
zlim ..... 351
zoom. ..... 401
zscore ..... 784
ztickangle ..... 352
zticklabels ..... 352
zticks ..... 351

\section*{Operator Index}

\begin{tabular}{|c|c|c|c|}
\hline \multirow[t]{2}{*}{^} & & \multicolumn{2}{|l|}{\}} \\
\hline & 161, 163, 900 & \} & 125 \\
\hline \multicolumn{4}{|l|}{@} \\
\hline @ class methods.. & ........ 889 & 1. & .166, 167, 900 \\
\hline @ function handle & ....... 160 & 11. & .......... 167 \\
\hline \multicolumn{4}{|l|}{\(\backslash \sim\)} \\
\hline \multirow[t]{2}{*}{\begin{tabular}{l}
\....................... \\
\ continuation marker
\end{tabular}} & 161, 162, 900 & \(\sim\) & 166, 167 \\
\hline & .......... 189 & \(\sim\) & 164, 165, 900 \\
\hline \(\{\) & & & \\
\hline , & .......... . 125 & & \\
\hline
\end{tabular}

\section*{Graphics Properties Index}
Axes Properties ..... 441
Figure Properties ..... 434
Image Properties ..... 455
Legend Properties ..... 449
Light Properties ..... 468
Line Properties ..... 450
Patch Properties ..... 457
Root Properties ..... 433
Scatter Properties ..... 461
Surface Properties ..... 465
Text Properties ..... 453
Uibuttongroup Properties ..... 472
Uicontextmenu Properties ..... 474
Uicontrol Properties ..... 478
Uimenu Properties ..... 470
Uipanel Properties ..... 476
Uipushtool Properties ..... 485
Uitoggletool Properties ..... 486
Uitoolbar Properties ..... 483
A
axes alim ..... 441
axes alimmode ..... 441
axes alphamap ..... 441
axes alphascale ..... 441
axes ambientlightcolor ..... 441
axes beingdeleted ..... 441
axes box ..... 441
axes boxstyle ..... 441
axes busyaction ..... 441
axes buttondownfcn ..... 441
axes cameraposition ..... 441
axes camerapositionmode ..... 442
axes cameratarget ..... 442
axes cameratargetmode ..... 442
axes cameraupvector ..... 442
axes cameraupvectormode ..... 442
axes cameraviewangle ..... 442
axes cameraviewanglemode ..... 442
axes children ..... 442
axes clim ..... 442
axes climmode ..... 442
axes clipping ..... 442
axes clippingstyle ..... 442
axes color ..... 442
axes colormap ..... 442
axes colororder ..... 442
axes colororderindex ..... 442
axes colorscale ..... 442
axes contextmenu. ..... 442
axes createfcn. ..... 442
axes currentpoint ..... 442
axes dataaspectratio ..... 442
axes dataaspectratiomode ..... 443
axes deletefcn ..... 443
axes fontangle ..... 443
axes fontname ..... 443
axes fontsize ..... 443
axes fontsizemode ..... 443
axes fontsmoothing ..... 443
axes fontunits. ..... 443
axes fontweight ..... 443
axes gridalpha ..... 443
axes gridalphamode ..... 443
axes gridcolor ..... 443
axes gridcolormode ..... 443
axes gridlinestyle ..... 443
axes handlevisibility ..... 443
axes hittest ..... 443
axes innerposition ..... 444
axes interactions ..... 444
axes interruptible ..... 444
axes labelfontsizemultiplier ..... 444
axes layer ..... 444
axes layout ..... 444
axes legend ..... 444
axes linestyleorder ..... 444
axes linestyleorderindex ..... 444
axes linewidth ..... 444
axes minorgridalpha ..... 444
axes minorgridalphamode ..... 444
axes minorgridcolor ..... 444
axes minorgridcolormode ..... 444
axes minorgridlinestyle ..... 444
axes mousewheelzoom ..... 444
axes nextplot ..... 444
axes nextseriesindex ..... 445
axes outerposition ..... 445
axes parent ..... 445
axes pickableparts ..... 445
axes plotboxaspectratio ..... 445
axes plotboxaspectratiomode ..... 445
axes position ..... 445
axes positionconstraint ..... 445
axes projection ..... 445
axes selected ..... 445
axes selectionhighlight ..... 445
axes sortmethod ..... 445
axes tag ..... 445
axes tickdir ..... 446
axes tickdirmode ..... 446
axes ticklabelinterpreter ..... 446
axes ticklength ..... 446
axes tightinset ..... 446
axes title ..... 446
axes titlefontsizemultiplier ..... 446
axes titlefontweight................................... . . . 446
axes toolbar ............................................... . . 446 figure alphamap ..... 434
axes type ..... 446
axes units ..... 446
axes userdata ..... 446
axes view ..... 446
axes visible ..... 446
axes xaxis ..... 446
axes xaxislocation ..... 446
axes xcolor ..... 446
axes xcolormode ..... 446
axes xdir ..... 447
axes xgrid ..... 447
axes xlabel ..... 447
axes xlim ..... 447
axes xlimmode ..... 447
axes xminorgrid ..... 447
axes xminortick ..... 447
axes xscale ..... 447
axes xtick ..... 447
axes xticklabel ..... 447
axes xticklabelmode. ..... 447
axes xticklabelrotation ..... 447
axes xtickmode ..... 447
axes yaxis ..... 447
axes yaxislocation ..... 447
axes ycolor ..... 447
axes ycolormode ..... 447
axes ydir ..... 447
axes ygrid ..... 447
axes ylabel ..... 447
axes ylim ..... 447
axes ylimmode ..... 447
axes yminorgrid ..... 448
axes yminortick ..... 448
axes yscale ..... 448
axes ytick ..... 448
axes yticklabel ..... 448
axes yticklabelmode ..... 448
axes yticklabelrotation ..... 448
axes ytickmode ..... 448
axes zaxis ..... 448
axes zcolor ..... 448
axes zcolormode ..... 448
axes zdir ..... 448
axes zgrid ..... 448
axes zlabel ..... 448
axes zlim ..... 448
axes zlimmode ..... 448
axes zminorgrid ..... 448
axes zminortick ..... 448
axes zscale ..... 448
axes ztick ..... 448
axes zticklabel ..... 448
axes zticklabelmode ..... 448
axes zticklabelrotation ..... 448
axes ztickmode ..... 448
figure beingdeleted ..... 434
figure busyaction. ..... 434
figure buttondownfcn ..... 435
figure children ..... 435
figure clipping ..... 435
figure closerequestfcn ..... 435
figure color ..... 435
figure colormap ..... 435
figure contextmenu ..... 435
figure createfcn ..... 435
figure currentaxes ..... 435
figure currentcharacter ..... 435
figure currentobject ..... 435
figure currentpoint ..... 435
figure deletefcn ..... 436
figure dockcontrols ..... 436
figure filename ..... 436
figure graphicssmoothing ..... 436
figure handlevisibility ..... 436
figure hittest ..... 436
figure integerhandle ..... 436
figure interruptible ..... 436
figure inverthardcopy ..... 436
figure keypressfcn ..... 436
figure keyreleasefcn ..... 436
figure menubar ..... 437
figure name ..... 437
figure nextplot ..... 437
figure number ..... 437
figure numbertitle ..... 437
figure outerposition ..... 437
figure paperorientation ..... 437
figure paperposition. ..... 437
figure paperpositionmode ..... 437
figure papersize ..... 438
figure papertype ..... 438
figure paperunits. ..... 438
figure parent ..... 438
figure pickableparts ..... 438
figure pointer ..... 438
figure pointershapecdata ..... 438
figure pointershapehotspot ..... 438
figure position ..... 438
figure renderer ..... 439
figure renderermode ..... 439
figure resize ..... 439
figure resizefcn ..... 439
figure selected ..... 439
figure selectionhighlight ..... 439
figure selectiontype ..... 439
figure sizechangedfcn ..... 439
figure tag ..... 439
figure toolbar ..... 439
figure type. ..... 439
figure units ..... 439
figure userdata ..... 440
figure visible ..... 440
light children ..... 468
figure windowbuttondownfcn ..... 440
figure windowbuttonmotionfcn ..... 440
figure windowbuttonupfen ..... 440
figure windowkeypressfcn ..... 440
figure windowkeyreleasefcn ..... 440
figure windowscrollwheelfcn ..... 440
figure windowstyle ..... 440
I
image alphadata ..... 455
image alphadatamapping ..... 455
image beingdeleted ..... 455
image busyaction ..... 456
image buttondownfcn ..... 456
image cdata ..... 456
image cdatamapping ..... 456
image children ..... 456
image clipping ..... 456
image contextmenu ..... 456
image createfcn ..... 456
image deletefcn ..... 456
image handlevisibility ..... 456
image hittest ..... 456
image interruptible ..... 456
image parent ..... 457
image pickableparts ..... 457
image selected ..... 457
image selectionhighlight ..... 457
image tag ..... 457
image type ..... 457
image userdata ..... 457
image visible ..... 457
image xdata ..... 457
image ydata ..... 457
L
legend autoupdate ..... 449
legend box ..... 449
legend color ..... 449
legend edgecolor ..... 449
legend fontangle ..... 449
legend fontname ..... 449
legend fontsize ..... 449
legend fontunits ..... 449
legend fontweight ..... 449
legend location ..... 449
legend orientation ..... 449
legend position ..... 450
legend string ..... 450
legend textcolor ..... 450
legend title ..... 450
light clipping ..... 469
light color ..... 469
light contextmenu ..... 469
light createfcn ..... 469
light deletefcn ..... 469
light handlevisibility ..... 469
light hittest ..... 469
light interruptible ..... 469
light parent ..... 469
light pickableparts ..... 469
light position ..... 470
light selected ..... 470
light selectionhighlight ..... 470
light style ..... 470
light tag ..... 470
light type ..... 470
light userdata ..... 470
light visible ..... 470
line beingdeleted ..... 450
line busyaction ..... 450
line buttondownfcn ..... 450
line children ..... 450
line clipping ..... 450
line color ..... 450
line contextmenu ..... 451
line createfcn ..... 451
line deletefcn ..... 451
line displayname ..... 451
line handlevisibility ..... 451
line hittest ..... 451
line interruptible ..... 451
line linejoin ..... 451
line linestyle ..... 451
line linewidth ..... 451
line marker ..... 451
line markeredgecolor ..... 451
line markerfacecolor ..... 452
line markersize ..... 452
line parent ..... 452
line pickableparts ..... 452
line selected ..... 452
line selectionhighlight ..... 452
line tag ..... 452
line type ..... 452
line userdata ..... 452
line visible ..... 452
line xdata ..... 452
line xdatasource ..... 452
line ydata ..... 452
line ydatasource ..... 452
line zdata ..... 452
line zdatasource ..... 452
legend units ..... 450
light beingdeleted ..... 468
light busyaction. ..... 468
light buttondownfen ..... 468

\section*{P}
patch alphadatamapping ..... 457
patch ambientstrength ..... 457
patch backfacelighting ..... 457
patch beingdeleted ..... 458
patch busyaction. ..... 458
patch buttondownfcn ..... 458
patch cdata ..... 458
patch cdatamapping ..... 458
patch children ..... 458
patch clipping ..... 458
patch contextmenu ..... 458
patch createfcn ..... 458
patch deletefcn ..... 458
patch diffusestrength ..... 458
patch displayname ..... 459
patch edgealpha ..... 459
patch edgecolor ..... 459
patch edgelighting ..... 459
patch facealpha ..... 459
patch facecolor ..... 459
patch facelighting ..... 459
patch facenormals ..... 459
patch facenormalsmode ..... 459
patch faces ..... 459
patch facevertexalphadata ..... 459
patch facevertexcdata ..... 459
patch handlevisibility ..... 459
patch hittest ..... 459
patch interruptible ..... 460
patch linestyle ..... 460
patch linewidth ..... 460
patch marker ..... 460
patch markeredgecolor ..... 460
patch markerfacecolor ..... 460
patch markersize ..... 460
patch parent ..... 460
patch pickableparts ..... 460
patch selected ..... 460
patch selectionhighlight ..... 460
patch specularcolorreflectance ..... 460
patch specularexponent ..... 460
patch specularstrength ..... 460
patch tag ..... 461
patch type ..... 461
patch userdata ..... 461
patch vertexnormals ..... 461
patch vertexnormalsmode ..... 461
patch vertices ..... 461
patch visible ..... 461
patch xdata ..... 461
patch ydata ..... 461
patch zdata ..... 461

\section*{R}
root beingdeleted...................................... . . 433
root busyaction....................................... . 433
root buttondownfen.................................... . . . 433
root callbackobject...................................... . . 433
root children.............................................. . . . . 433
root clipping........................................... . . 433
root commandwindowsize . . . . . . . . . . . . . . . . . . . . . 433
root contextmenu ...................................... . . . . 433
root createfcn.............................................. . . . 433
root currentfigure....................................... . . . 433
root deletefcn............................................ . . . . 433
root fixedwidthfontname . . . . . . . . . . . . . . . . . . . . . . . 433
root handlevisibility................................... . . . . 433
root hittest.............................................. . . . 433
root interruptible . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 433
root monitorpositions . . . . . . . . . . . . . . . . . . . . . . . . . 433
root parent.............................................. . . . 434
root pickableparts...................................... . . . 434
root pointerlocation.................................. . . . 434
root pointerwindow . . . . . . . . . . . . . . . . . . . . . . . . . . . 434
root screendepth........................................ . . . . . 434
root screenpixelsperinch............................. . . . 434
root screensize. . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 434
root selected. . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 434
root selectionhighlight . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 434
root showhiddenhandles............................. . . . 434
root tag.................................................. . . 434
root type.................................................. . . . . . . 434
root units..................................................... 434
root userdata.............................................. . . 434
root visible................................................ . . . 434

\section*{S}
scatter annotation.................................... . . . 461
scatter beingdeleted.................................... . . 461
scatter busyaction . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 461
scatter buttondownfcn .............................. . . . . 461
scatter cdata.............................................. . . 461
scatter cdatamode . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter cdatasource . . . . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter children......................................... . . . 462
scatter clipping . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter contextmenu . . . . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter createfcn......................................... . 462
scatter datatiptemplate ............................ . . 462
scatter deletefcn . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter displayname................................. . . . . . . . . . . . . 462
scatter handlevisibility . . . . . . . . . . . . . . . . . . . . . . . . . 462
scatter hittest .......................................... . . 462
scatter interruptible.................................. . 462
scatter latitudedata.................................... . . . . 462
scatter latitudedatasource.......................... . . . 463
scatter linewidth......................................... . . . 463
scatter longitudedata ............................... . . . 463
scatter longitudedatasource . . . . . . . . . . . . . . . . . . . . 463
scatter marker............................................ . . . 463
scatter markeredgealpha ..... 463
surface markersize ..... 467
scatter markeredgecolor ..... 463
scatter markerfacealpha ..... 463
scatter markerfacecolor ..... 463
scatter parent ..... 463
scatter pickableparts ..... 463
scatter rdata ..... 463
scatter rdatasource ..... 464
scatter selected ..... 464
scatter selectionhighlight ..... 464
scatter seriesindex ..... 464
scatter sizedata ..... 464
scatter sizedatasource ..... 464
scatter tag ..... 464
scatter thetadata ..... 464
scatter thetadatasource ..... 464
scatter type. ..... 464
scatter userdata ..... 464
scatter visible ..... 464
scatter xdata ..... 464
scatter xdatasource ..... 464
scatter ydata ..... 464
scatter ydatasource ..... 464
scatter zdata ..... 464
scatter zdatasource ..... 464
surface alphadata ..... 465
surface alphadatamapping ..... 465
surface ambientstrength ..... 465
surface backfacelighting ..... 465
surface beingdeleted ..... 465
surface busyaction ..... 465
surface buttondownfcn ..... 465
surface cdata ..... 465
surface cdatamapping ..... 465
surface cdatasource ..... 465
surface children ..... 465
surface clipping ..... 465
surface contextmenu ..... 465
surface createfcn ..... 465
surface deletefcn ..... 465
surface diffusestrength ..... 466
surface displayname ..... 466
surface edgealpha ..... 466
surface edgecolor ..... 466
surface edgelighting ..... 466
surface facealpha ..... 466
surface facecolor ..... 466
surface facelighting ..... 466
surface facenormals ..... 466
surface facenormalsmode ..... 466
surface handlevisibility ..... 466
surface hittest ..... 466
surface interruptible ..... 467
surface linestyle ..... 467
surface linewidth ..... 467
surface marker ..... 467
surface markeredgecolor ..... 467
surface markerfacecolor ..... 467
surface meshstyle ..... 467
surface parent ..... 467
surface pickableparts ..... 467
surface selected ..... 467
surface selectionhighlight ..... 467
surface specularcolorreflectance ..... 467
surface specularexponent ..... 467
surface specularstrength ..... 468
surface tag. ..... 468
surface type ..... 468
surface userdata ..... 468
surface vertexnormals ..... 468
surface vertexnormalsmode ..... 468
surface visible ..... 468
surface xdata ..... 468
surface xdatasource ..... 468
surface ydata ..... 468
surface ydatasource ..... 468
surface zdata ..... 468
surface zdatasource ..... 468
text backgroundcolor ..... 453
text beingdeleted ..... 453
text busyaction ..... 453
text buttondownfen ..... 453
text children ..... 453
text clipping ..... 453
text color ..... 453
text contextmenu ..... 453
text createfcn ..... 453
text deletefcn ..... 453
text edgecolor ..... 453
text editing ..... 453
text extent ..... 453
text fontangle. ..... 453
text fontname ..... 454
text fontsize ..... 454
text fontsmoothing ..... 454
text fontunits ..... 454
text fontweight ..... 454
text handlevisibility ..... 454
text hittest ..... 454
text horizontalalignment ..... 454
text interpreter ..... 454
text interruptible ..... 454
text linestyle ..... 454
text linewidth ..... 454
text margin ..... 454
text parent ..... 455
text pickableparts ..... 455
text position ..... 455
text rotation ..... 455
text selected ..... 455
text selectionhighlight ..... 455
text string ..... 455
text tag ..... 455
text type ..... 455
text units ..... 455
text userdata ..... 455
text verticalalignment ..... 455
text visible ..... 455
U
uibuttongroup backgroundcolor ..... 472
uibuttongroup beingdeleted ..... 472
uibuttongroup bordertype. ..... 472
uibuttongroup borderwidth ..... 472
uibuttongroup busyaction ..... 472
uibuttongroup buttondownfen ..... 472
uibuttongroup children ..... 472
uibuttongroup clipping ..... 472
uibuttongroup contextmenu .. ..... 472
uibuttongroup createfcn ..... 472
uibuttongroup deletefcn ..... 472
uibuttongroup fontangle ..... 472
uibuttongroup fontname ..... 473
uibuttongroup fontsize ..... 473
uibuttongroup fontunits. ..... 473
uibuttongroup fontweight ..... 473
uibuttongroup foregroundcolor ..... 473
uibuttongroup handlevisibility ..... 473
uibuttongroup highlightcolor . ..... 473
uibuttongroup hittest ..... 473
uibuttongroup interruptible ..... 473
uibuttongroup parent ..... 473
uibuttongroup pickableparts ..... 473
uibuttongroup position ..... 474
uibuttongroup resizefcn ..... 474
uibuttongroup selected ..... 474
uibuttongroup selectedobject ..... 474
uibuttongroup selectionchangedfcn ..... 474
uibuttongroup selectionhighlight ..... 474
uibuttongroup shadowcolor ..... 474
uibuttongroup sizechangedfcn ..... 474
uibuttongroup tag ..... 474
uibuttongroup title ..... 474
uibuttongroup titleposition ..... 474
uibuttongroup type ..... 474
uibuttongroup units ..... 474
uibuttongroup userdata ..... 474
uibuttongroup visible ..... 474
uicontextmenu beingdeleted ..... 474
uicontextmenu busyaction ..... 474
uicontextmenu buttondownfcn ..... 474
uicontextmenu callback ..... 474
uicontextmenu children ..... 474
uicontextmenu clipping ..... 474
uicontextmenu contextmenu ..... 475
uicontextmenu createfcn ..... 475
uicontextmenu deletefcn ..... 475
uicontextmenu handlevisibility ..... 475
uicontextmenu hittest ..... 475
uicontextmenu interruptible ..... 475
uicontextmenu parent ..... 475
uicontextmenu pickableparts ..... 475
uicontextmenu position ..... 475
uicontextmenu selected ..... 476
uicontextmenu selectionhighlight ..... 476
uicontextmenu tag ..... 476
uicontextmenu type ..... 476
uicontextmenu userdata ..... 476
uicontextmenu visible ..... 476
uicontrol backgroundcolor ..... 478
uicontrol beingdeleted ..... 478
uicontrol busyaction ..... 478
uicontrol buttondownfen ..... 478
uicontrol callback ..... 478
uicontrol cdata ..... 479
uicontrol children ..... 479
uicontrol clipping ..... 479
uicontrol contextmenu ..... 479
uicontrol createfcn ..... 479
uicontrol deletefcn ..... 479
uicontrol enable ..... 479
uicontrol extent ..... 479
uicontrol fontangle ..... 479
uicontrol fontname ..... 479
uicontrol fontsize ..... 479
uicontrol fontunits ..... 479
uicontrol fontweight ..... 479
uicontrol foregroundcolor ..... 479
uicontrol handlevisibility ..... 480
uicontrol hittest ..... 480
uicontrol horizontalalignment ..... 480
uicontrol interruptible ..... 480
uicontrol keypressfcn. ..... 480
uicontrol listboxtop ..... 480
uicontrol max ..... 480
uicontrol min ..... 480
uicontrol parent ..... 480
uicontrol pickableparts ..... 480
uicontrol position ..... 480
uicontrol selected ..... 480
uicontrol selectionhighlight ..... 480
uicontrol sliderstep ..... 480
uicontrol string ..... 480
uicontrol style ..... 480
uicontrol tag ..... 480
uicontrol tooltipstring ..... 480
uicontrol type ..... 480
uicontrol units.. ..... 480
uicontrol userdata ..... 481
uicontrol value ..... 481
uicontrol verticalalignment ..... 481
uicontrol visible ..... 481
uimenu accelerator ..... 470
uimenu beingdeleted ..... 470
uimenu busyaction ..... 470
uimenu buttondownfen. ..... 470
uimenu callback ..... 470
uimenu checked ..... 470
uimenu children ..... 470
uimenu clipping ..... 470
uimenu contextmenu ..... 470
uimenu createfcn ..... 470
uimenu deletefcn ..... 471
uimenu enable ..... 471
uimenu foregroundcolor ..... 471
uimenu handlevisibility ..... 471
uimenu hittest ..... 471
uimenu interruptible ..... 471
uimenu label ..... 471
uimenu parent ..... 471
uimenu pickableparts ..... 471
uimenu position ..... 471
uimenu selected ..... 471
uimenu selectionhighlight ..... 471
uimenu separator ..... 471
uimenu tag ..... 471
uimenu text ..... 471
uimenu type ..... 472
uimenu userdata ..... 472
uimenu visible ..... 472
uipanel backgroundcolor ..... 476
uipanel beingdeleted ..... 476
uipanel bordertype ..... 476
uipanel borderwidth ..... 476
uipanel busyaction ..... 476
uipanel buttondownfcn ..... 476
uipanel children ..... 476
uipanel clipping ..... 476
uipanel contextmenu ..... 476
uipanel createfcn. ..... 476
uipanel deletefcn ..... 476
uipanel fontangle ..... 477
uipanel fontname ..... 477
uipanel fontsize ..... 477
uipanel fontunits ..... 477
uipanel fontweight ..... 477
uipanel foregroundcolor ..... 477
uipanel handlevisibility ..... 477
uipanel highlightcolor ..... 477
uipanel hittest ..... 477
uipanel interruptible ..... 477
uipanel parent ..... 477
uipanel pickableparts ..... 477
uipanel position ..... 478
uipanel resizefcn ..... 478
uipanel selected ..... 478
uipanel selectionhighlight ..... 478
uipanel shadowcolor ..... 478
uipanel sizechangedfcn ..... 478
uipanel tag ..... 478
uipanel title ..... 478
uipanel titleposition ..... 478
uipanel type ..... 478
uipanel units ..... 478
uipanel userdata ..... 478
uipanel visible ..... 478
uipushtool __named_icon ..... 485
uipushtool beingdeleted ..... 485
uipushtool busyaction ..... 485
uipushtool buttondownfcn ..... 485
uipushtool cdata ..... 485
uipushtool children ..... 485
uipushtool clickedcallback ..... 485
uipushtool clipping ..... 485
uipushtool contextmenu ..... 485
uipushtool createfcn ..... 485
uipushtool deletefcn ..... 485
uipushtool enable ..... 485
uipushtool handlevisibility ..... 485
uipushtool hittest ..... 485
uipushtool interruptible ..... 486
uipushtool parent ..... 486
uipushtool pickableparts ..... 486
uipushtool selected ..... 486
uipushtool selectionhighlight ..... 486
uipushtool separator ..... 486
uipushtool tag ..... 486
uipushtool tooltipstring ..... 486
uipushtool type ..... 486
uipushtool userdata ..... 486
uipushtool visible ..... 486
uitable backgroundcolor ..... 481
uitable beingdeleted ..... 481
uitable busyaction ..... 481
uitable buttondownfcn ..... 481
uitable celleditcallback ..... 481
uitable cellselectioncallback ..... 481
uitable children ..... 481
uitable clipping ..... 481
uitable columneditable ..... 481
uitable columnformat ..... 481
uitable columnname ..... 481
uitable columnwidth ..... 481
uitable contextmenu ..... 481
uitable createfcn ..... 481
uitable data ..... 481
uitable deletefcn ..... 482
uitable enable. ..... 482
uitable extent ..... 482
uitable fontangle ..... 482
uitable fontname ..... 482
uitable fontsize ..... 482
uitable fontunits ..... 482
uitable fontweight ..... 482
uitable foregroundcolor ..... 482
uitable handlevisibility ..... 482
uitable hittest ..... 482
uitable interruptible ..... 482
uitable keypressfcn. ..... 482
uitable keyreleasefcn ..... 482
uitable parent ..... 482
uitable pickableparts ..... 482
uitable position ..... 483
uitable rearrangeablecolumns ..... 483
uitable rowname ..... 483
uitable rowstriping ..... 483
uitable selected ..... 483
uitable selectionhighlight ..... 483
uitable tag ..... 483
uitable tooltipstring ..... 483
uitable type ..... 483
uitable units ..... 483
uitable userdata ..... 483
uitable visible ..... 483
uitoggletool __named_icon ..... 486
uitoggletool beingdeleted ..... 486
uitoggletool busyaction ..... 486
uitoggletool buttondownfcn ..... 487
uitoggletool cdata. ..... 487
uitoggletool children ..... 487
uitoggletool clickedcallback. ..... 487
uitoggletool clipping ..... 487
uitoggletool contextmenu ..... 487
uitoggletool createfcn ..... 487
uitoggletool deletefcn ..... 487
uitoggletool enable ..... 487
uitoggletool handlevisibility ..... 487
uitoggletool hittest ..... 487
uitoggletool interruptible ..... 487
uitoggletool offcallback ..... 487
uitoggletool oncallback ..... 488
uitoggletool parent ..... 488
uitoggletool pickableparts ..... 488
uitoggletool selected ..... 488
uitoggletool selectionhighlight ..... 488
uitoggletool separator ..... 488
uitoggletool state ..... 488
uitoggletool tag ..... 488
uitoggletool tooltipstring ..... 488
uitoggletool type ..... 488
uitoggletool userdata. ..... 488
uitoggletool visible ..... 488
uitoolbar beingdeleted ..... 483
uitoolbar busyaction ..... 483
uitoolbar buttondownfen ..... 483
uitoolbar children ..... 483
uitoolbar clipping ..... 483
uitoolbar contextmenu ..... 483
uitoolbar createfcn ..... 484
uitoolbar deletefcn ..... 484
uitoolbar handlevisibility ..... 484
uitoolbar hittest ..... 484
uitoolbar interruptible ..... 484
uitoolbar parent ..... 484
uitoolbar pickableparts ..... 484
uitoolbar selected ..... 484
uitoolbar selectionhighlight ..... 484
uitoolbar tag ..... 484
uitoolbar type ..... 484
uitoolbar userdata ..... 485
uitoolbar visible ..... 485```


[^0]:    ${ }^{1}$ The '\#!' mechanism works on Unix systems derived from Berkeley Unix, System V Release 4, and some System V Release 3 systems.

[^1]:    ${ }^{1}$ Comma-separated lists are also sometimes informally referred to as cs-lists.

[^2]:    ${ }^{1}$ Some of Octave's functions are implemented in terms of functions that cannot be called recursively. For example, the ODE solver lsode is ultimately implemented in a Fortran subroutine that cannot be called recursively, so lsode should not be called either directly or indirectly from within the user-supplied function that lsode requires. Doing so will result in an error.
    ${ }^{2}$ It would be much better to use prod (1:n), or gamma ( $n+1$ ) instead, after first checking to ensure that the value n is actually a positive integer.

[^3]:    ${ }^{1}$ We only know it is the binary multiplication operator, but fortunately this operator appears only at one place in the code and thus we know which occurrence takes so much time. If there were multiple places, we would have to use the hierarchical profile to find out the exact place which uses up the time which is not covered in this example.

