

# Maxima Manual

Maxima is a computer algebra system, implemented in Lisp.

Maxima is derived from the Macsyma system, developed at MIT in the years 1968 through 1982 as part of Project MAC. MIT turned over a copy of the Macsyma source code to the Department of Energy in 1982; that version is now known as DOE Macsyma. A copy of DOE Macsyma was maintained by Professor William F. Schelter of the University of Texas from 1982 until his death in 2001. In 1998, Schelter obtained permission from the Department of Energy to release the DOE Macsyma source code under the GNU Public License, and in 2000 he initiated the Maxima project at SourceForge to maintain and develop DOE Macsyma, now called Maxima.



# 1 Introduction to Maxima

Start Maxima with the command "maxima". Maxima will display version information and a prompt. End each Maxima command with a semicolon. End the session with the command "quit()". Here's a sample session:

```
[wfs@chromium]$ maxima
Maxima 5.9.1 http://maxima.sourceforge.net
Using Lisp CMU Common Lisp 19a
Distributed under the GNU Public License. See the file COPYING.
Dedicated to the memory of William Schelter.
This is a development version of Maxima. The function bug_report()
provides bug reporting information.
(%i1) factor(10!);

(%o1)
      8 4 2
      2 3 5 7
(%i2) expand ((x + y)^6);
      6 5 2 4 3 3 4 2 5 6
(%o2) y + 6 x y + 15 x y + 20 x y + 15 x y + 6 x y + x
(%i3) factor (x^6 - 1);

(%o3)
      2 2
      (x - 1) (x + 1) (x - x + 1) (x + x + 1)
(%i4) quit();
[wfs@chromium]$
```

Maxima can search the info pages. Use the *describe* command to show all the commands and variables containing a string, and optionally their documentation. The question mark ? is an abbreviation for describe:

```
(%i1) ? integ

0: (maxima.info)Introduction to Elliptic Functions and Integrals.
1: Definitions for Elliptic Integrals.
2: Integration.
3: Introduction to Integration.
4: Definitions for Integration.
5: askinteger :Definitions for Simplification.
6: integerp :Definitions for Miscellaneous Options.
7: integrate :Definitions for Integration.
8: integrate_use_rootsof :Definitions for Integration.
9: integration_constant_counter :Definitions for Integration.
Enter space-separated numbers, 'all' or 'none': 6 5
```

Info from file /usr/local/info/maxima.info:

- Function: integerp (<expr>)
  - Returns 'true' if <expr> is an integer, otherwise 'false'.
- Function: askinteger (expr, integer)
- Function: askinteger (expr)
- Function: askinteger (expr, even)

- Function: askinteger (expr, odd)  
 'askinteger (expr, integer)' attempts to determine from the 'assume' database whether 'expr' is an integer. 'askinteger' will ask the user if it cannot tell otherwise, and attempt to install the information in the database if possible. 'askinteger (expr)' is equivalent to 'askinteger (expr, integer)'.  
 'askinteger (expr, even)' and 'askinteger (expr, odd)' likewise attempt to determine if 'expr' is an even integer or odd integer, respectively.

(%o1) false

To use a result in later calculations, you can assign it to a variable or refer to it by its automatically supplied label. In addition, % refers to the most recent calculated result:

(%i1) u: expand ((x + y)^6);  
 (%o1)  $y^6 + 6xy^5 + 15x^2y^4 + 20x^3y^3 + 15x^4y^2 + 6x^5y + x^6$   
 (%i2) diff (u, x);  
 (%o2)  $6y^5 + 30x^4y^4 + 60x^2y^3 + 60x^3y^2 + 30x^4y + 6x^5$   
 (%i3) factor (%o2);  
 (%o3)  $6(y + x)^5$

Maxima knows about complex numbers and numerical constants:

(%i1) cos(%pi);  
 (%o1) - 1  
 (%i2) exp(%i\*%pi);  
 (%o2) - 1

Maxima can do differential and integral calculus:

(%i1) u: expand ((x + y)^6);  
 (%o1)  $y^6 + 6xy^5 + 15x^2y^4 + 20x^3y^3 + 15x^4y^2 + 6x^5y + x^6$   
 (%i2) diff (%o1, x);  
 (%o2)  $6y^5 + 30x^4y^4 + 60x^2y^3 + 60x^3y^2 + 30x^4y + 6x^5$   
 (%i3) integrate (1/(1 + x^3), x);  
 (%o3)  $-\frac{\log(x^2 - x + 1)}{6} + \frac{\operatorname{atan}\left(\frac{2x - 1}{\sqrt{3}}\right)}{\sqrt{3}} + \frac{\log(x + 1)}{3}$

Maxima can solve linear systems and cubic equations:

(%i1) linsolve ([3\*x + 4\*y = 7, 2\*x + a\*y = 13], [x, y]);  
 (%o1)  $\left[x = \frac{7a - 52}{3a - 8}, y = \frac{25}{3a - 8}\right]$   
 (%i2) solve (x^3 - 3\*x^2 + 5\*x = 15, x);

```
(%o2) [x = - sqrt(5) %i, x = sqrt(5) %i, x = 3]
```

Maxima can solve nonlinear sets of equations. Note that if you don't want a result printed, you can finish your command with \$ instead of ;.

```
(%i1) eq_1: x^2 + 3*x*y + y^2 = 0$
```

```
(%i2) eq_2: 3*x + y = 1$
```

```
(%i3) solve ([eq_1, eq_2]);
```

```
(%o3) [[y = -  $\frac{3 \sqrt{5} + 7}{2}$ , x =  $\frac{\sqrt{5} + 3}{2}$ ],
```

```
      [y =  $\frac{3 \sqrt{5} - 7}{2}$ , x = -  $\frac{\sqrt{5} - 3}{2}$ ]]
```

Maxima can generate plots of one or more functions:

```
(%i1) eq_1: x^2 + 3*x*y + y^2 = 0$
```

```
(%i2) eq_2: 3*x + y = 1$
```

```
(%i3) solve ([eq_1, eq_2]);
```

```
(%o3) [[y = -  $\frac{3 \sqrt{5} + 7}{2}$ , x =  $\frac{\sqrt{5} + 3}{2}$ ],
```

```
      [y =  $\frac{3 \sqrt{5} - 7}{2}$ , x = -  $\frac{\sqrt{5} - 3}{2}$ ]]
```

```
(%i4) kill(labels);
```

```
(%o0) done
```

```
(%i1) plot2d (sin(x)/x, [x, -20, 20]);
```

```
(%o1)
```

```
(%i2) plot2d ([atan(x), erf(x), tanh(x)], [x, -5, 5]);
```

```
(%o2)
```

```
(%i3) plot3d (sin(sqrt(x^2 + y^2))/sqrt(x^2 + y^2), [x, -12, 12], [y, -12, 12]);
```

```
(%o3)
```



## 2 Bug Detection and Reporting

### 2.1 Introduction to Bug Detection and Reporting

Like all large programs, Maxima contains both known and unknown bugs. This chapter describes the built-in facilities for running the Maxima test suite as well as reporting new bugs.

### 2.2 Definitions for Bug Detection and Reporting

**run\_testsuite** () Function  
**run\_testsuite** (*boolean*) Function  
**run\_testsuite** (*boolean, boolean*) Function

Run the Maxima test suite. Tests producing the desired answer are considered “passes,” as are tests that do not produce the desired answer, but are marked as known bugs.

`run_testsuite` () displays only tests that do not pass.

`run_testsuite` (`true`) displays tests that are marked as known bugs, as well as failures.

`run_testsuite` (`true, true`) displays all tests.

`run_testsuite` changes the Maxima environment. Typically a test script executes `kill` to establish a known environment (namely one without user-defined functions and variables) and then defines functions and variables appropriate to the test.

`run_testsuite` returns `done`.

**bug\_report** () Function

Prints out Maxima and Lisp version numbers, and gives a link to the Maxima project bug report web page. The version information is the same as reported by `build_info`.

When a bug is reported, it is helpful to copy the Maxima and Lisp version information into the bug report.

`bug_report` returns an empty string "".

**build\_info** () Function

Prints out a summary of the parameters of the Maxima build.

`build_info` returns an empty string "".





## 3 Help

### 3.1 Introduction to Help

The primary on-line help function is `describe`, which is typically invoked by the question mark `?` at the interactive prompt. `? foo` (with a space between `?` and `foo`) is equivalent to `describe ("foo")`, where `foo` is the name or part of the name of a function or topic; `describe` then finds all documented items which contain the string `foo` in their titles. If there is more than one such item, Maxima asks the user to select an item or items to display.

```
(%i1) ? integ
0: (maxima.info)Introduction to Elliptic Functions and Integrals.
1: Definitions for Elliptic Integrals.
2: Integration.
3: Introduction to Integration.
4: Definitions for Integration.
5: askinteger :Definitions for Simplification.
6: integerp :Definitions for Miscellaneous Options.
7: integrate :Definitions for Integration.
8: integrate_use_rootsof :Definitions for Integration.
9: integration_constant_counter :Definitions for Integration.
Enter space-separated numbers, 'all' or 'none': 7 8
```

```
Info from file /use/local/maxima/doc/info/maxima.info:
- Function: integrate (expr, var)
- Function: integrate (expr, var, a, b)
  Attempts to symbolically compute the integral of 'expr' with
  respect to 'var'. 'integrate (expr, var)' is an indefinite
  integral, while 'integrate (expr, var, a, b)' is a definite
  integral, [...]
```

In this example, items 7 and 8 were selected. All or none of the items could have been selected by entering `all` or `none`, which can be abbreviated `a` or `n`, respectively.

### 3.2 Lisp and Maxima

Maxima is written in Lisp, and it is easy to access Lisp functions and variables from Maxima and vice versa. Lisp and Maxima symbols are distinguished by a naming convention. A Lisp symbol which begins with a dollar sign `$` corresponds to a Maxima symbol without the dollar sign. A Maxima symbol which begins with a question mark `?` corresponds to a Lisp symbol without the question mark. For example, the Maxima symbol `foo` corresponds to the Lisp symbol `$foo`. while the Maxima symbol `?foo` corresponds to the Lisp symbol `foo`, Note that `?foo` is written without a space between `?` and `foo`; otherwise it might be mistaken for `describe ("foo")`.

Lisp symbols appearing in Maxima code must follow the rules for Maxima identifiers (apart from beginning with a question mark). In particular, symbols containing hyphen - or other special characters cannot appear in Maxima code.

Lisp code may be executed from within a Maxima session. A single line of Lisp (containing one or more forms) may be executed by the special command `:lisp`. For example,

```
(%i1) :lisp (foo $x $y)
```

calls the Lisp function `foo` with Maxima variables `x` and `y` as arguments. The `:lisp` construct can appear at the interactive prompt or in a file processed by `batch` or `demo`, but not in a file processed by `load`, `batchload`, `translate_file`, or `compile_file`.

The function `to_lisp()` opens an interactive Lisp session. Entering `(to-maxima)` closes the Lisp session and returns to Maxima.

Lisp functions and variables which are to be visible in Maxima as functions and variables with ordinary names (no special punctuation) must have Lisp names beginning with the dollar sign `$`.

Maxima is case-sensitive, distinguishing between lowercase and uppercase letters in identifiers, while Lisp is not. There are some rules governing the translation of names between Lisp and Maxima.

1. A Lisp identifier not enclosed in vertical bars corresponds to a Maxima identifier in lowercase. Whether the Lisp identifier is uppercase, lowercase, or mixed case, is ignored. E.g., Lisp `$foo`, `$FOO`, and `$Foo` all correspond to Maxima `foo`.
2. A Lisp identifier which is all uppercase or all lowercase and enclosed in vertical bars corresponds to a Maxima identifier with case reversed. That is, uppercase is changed to lowercase and lowercase to uppercase. E.g., Lisp `|$FOO|` and `|$foo|` correspond to Maxima `foo` and `FOO`, respectively.
3. A Lisp identifier which is mixed uppercase and lowercase and enclosed in vertical bars corresponds to a Maxima identifier with the same case. E.g., Lisp `|$Foo|` corresponds to Maxima `Foo`.

The `#$` Lisp macro allows the use of Maxima expressions in Lisp code. `#$expr$` expands to a Lisp expression equivalent to the Maxima expression `expr`.

```
(msetq $foo #$(x, y)$)
```

This has the same effect as entering

```
(%i1) foo: [x, y];
```

The Lisp function `displa` prints an expression in Maxima format.

```
(%i1) :lisp #$(x, y, z)$
((MLIST SIMP) $X $Y $Z)
(%i1) :lisp (displa '((MLIST SIMP) $X $Y $Z))
[x, y, z]
NIL
```

Functions defined in Maxima are not ordinary Lisp functions. The Lisp function `mfuncall` calls a Maxima function. For example:

```
(%i1) foo(x,y) := x*y$
(%i2) :lisp (mfuncall '$foo 'a 'b)
((MTIMES SIMP) A B)
```

Some Lisp functions are shadowed in the Maxima package, namely the following.

`complement`, `continue`, `//`, `float`, `functionp`, `array`, `exp`, `listen`, `signum`, `atan`, `asin`, `acos`, `asinh`, `acosh`, `atanh`, `tanh`, `cosh`, `sinh`, `tan`, `break`, and `gcd`.

### 3.3 Garbage Collection

Symbolic computation tends to create a good deal of garbage, and effective handling of this can be crucial to successful completion of some programs.

Under GCL, on UNIX systems where the `mprotect` system call is available (including SUN OS 4.0 and some variants of BSD) a stratified garbage collection is available. This limits the collection to pages which have been recently written to. See the GCL documentation under `ALLOCATE` and `GBC`. At the Lisp level doing `(setq si::*notify-gbc* t)` will help you determine which areas might need more space.

### 3.4 Documentation

The Maxima on-line user's manual can be viewed in different forms. From the Maxima interactive prompt, the user's manual is viewed as plain text by the `? command` (i.e., the `describe` function). The user's manual is viewed as `info` hypertext by the `info viewer` program and as a web page by any ordinary web browser.

`example` displays examples for many Maxima functions. For example,

```
(%i1) example (integrate);
yields
(%i2) test(f):=block([u],u:integrate(f,x),ratsimp(f-diff(u,x)))
(%o2) test(f) := block([u], u : integrate(f, x),
                                ratsimp(f - diff(u, x)))

(%i3) test(sin(x))
(%o3)                                0
(%i4) test(1/(x+1))
(%o4)                                0
(%i5) test(1/(x^2+1))
(%o5)                                0
```

and additional output.

### 3.5 Definitions for Help

**demo** (*filename*)

Function

Evaluates Maxima expressions in *filename* and displays the results. `demo` pauses after evaluating each expression and continues after the user enters a carriage return. (If running in Xmaxima, `demo` may need to see a semicolon ; followed by a carriage return.)

`demo` searches the list of directories `file_search_demo` to find *filename*. If the file has the suffix `dem`, the suffix may be omitted. See also `file_search`.

`demo` evaluates its argument. `demo` returns the name of the demonstration file.

Example:

```
(%i1) demo ("disol");

batching /home/wfs/maxima/share/simplification/disol.dem
```

At the \_ prompt, type ';' followed by enter to get next demo  
 (%i2) load(disol)

exp1 : a (e (g + f) + b (d + c))  
 (%o3) a (e (g + f) + b (d + c))

disolate(exp1, a, b, e)  
 (%t4) d + c

(%t5) g + f

(%o5) a (%t5 e + %t4 b)

(%i5) demo ("rncomb");

batching /home/wfs/maxima/share/simplification/rncomb.dem  
 At the \_ prompt, type ';' followed by enter to get next demo  
 (%i6) load(rncomb)

exp1 :  $\frac{z}{y + x} + \frac{x}{2 (y + x)}$   
 (%o7)  $\frac{z}{y + x} + \frac{x}{2 (y + x)}$

combine(exp1)  
 (%o8)  $\frac{z}{y + x} + \frac{x}{2 (y + x)}$

rncombine(%)  
 (%o9)  $\frac{2 z + x}{2 (y + x)}$

exp2 :  $\frac{d}{3} + \frac{c}{3} + \frac{b}{2} + \frac{a}{2}$   
 (%o10)  $\frac{d}{3} + \frac{c}{3} + \frac{b}{2} + \frac{a}{2}$

```

-
(%i11)          combine(exp2)
              2 d + 2 c + 3 (b + a)
(%o11)  -----
              6

```

```

-
(%i12)          rncombine(exp2)
              2 d + 2 c + 3 b + 3 a
(%o12)  -----
              6

```

```

-
(%i13)

```

**describe** (*string*)

Function

Finds all documented items which contain *string* in their titles. If there is more than one such item, Maxima asks the user to select an item or items to display. At the interactive prompt, ? foo (with a space between ? and foo) is equivalent to describe ("foo").

describe ("") yields a list of all topics documented in the on-line manual.

describe quotes its argument. describe always returns false.

Example:

```

(%i1) ? integ
0: (maxima.info)Introduction to Elliptic Functions and Integrals.
1: Definitions for Elliptic Integrals.
2: Integration.
3: Introduction to Integration.
4: Definitions for Integration.
5: askinteger :Definitions for Simplification.
6: integerp :Definitions for Miscellaneous Options.
7: integrate :Definitions for Integration.
8: integrate_use_rootsof :Definitions for Integration.
9: integration_constant_counter :Definitions for Integration.
Enter space-separated numbers, 'all' or 'none': 7 8

```

```

Info from file /use/local/maxima/doc/info/maxima.info:

```

- Function: integrate (expr, var)
- Function: integrate (expr, var, a, b)
  - Attempts to symbolically compute the integral of 'expr' with respect to 'var'. 'integrate (expr, var)' is an indefinite integral, while 'integrate (expr, var, a, b)' is a definite integral, [...]

In this example, items 7 and 8 were selected. All or none of the items could have been selected by entering all or none, which can be abbreviated a or n, respectively.

see [Section 3.1 \[Introduction to Help\]](#), page 9

**example** (*topic*) Function  
**example** () Function

**example** (*topic*) displays some examples of *topic*, which is a symbol (not a string). Most topics are function names. **example** () returns the list of all recognized topics.

The name of the file containing the examples is given by the global variable `manual_demo`, which defaults to "manual\_demo".

**example** quotes its argument. **example** returns `done` unless there is an error or there is no argument, in which case **example** returns the list of all recognized topics.

Examples:

```
(%i1) example (append);
(%i2) append([x+y,0,-3.2],[2.5E+20,x])
(%o2)      [y + x, 0, - 3.2, 2.5E+20, x]
(%o2)      done
(%i3) example (coeff);
(%i4) coeff(b+tan(x)+2*a*tan(x) = 3+5*tan(x),tan(x))
(%o4)      2 a + 1 = 5
(%i5) coeff(1+x*e^x+y,x,0)
(%o5)      y + 1
(%o5)      done
```

## 4 Command Line

### 4.1 Introduction to Command Line

**''** operator

The single quote operator `'` prevents evaluation.

E.g., `'(f(x))` means do not evaluate the expression  $f(x)$ . `'f(x)` (with the single quote applied to  $f$  instead of  $f(x)$ ) means return the noun form of  $f$  applied to  $[x]$ .

The single quote does not prevent simplification.

**'''** operator

The `''` (double single quotes) operator causes an extra evaluation to occur. E.g., `''%i4` will re-evaluate input line `%i4`. `''(f(x))` means evaluate the expression  $f(x)$  an extra time. `''f(x)` (with the double single quotes applied to  $f$  instead of  $f(x)$ ) means return the verb form of  $f$  applied to  $[x]$ .

### 4.2 Definitions for Command Line

**alias** (*new\_name\_1, old\_name\_1, ..., new\_name\_n, old\_name\_n*) Function  
 provides an alternate name for a (user or system) function, variable, array, etc. Any even number of arguments may be used.

**debugmode** Variable  
 Default value: `false`

When a Maxima error occurs, Maxima will start the debugger if `debugmode` is `true`. The user may enter commands to examine the call stack, set breakpoints, step through Maxima code, and so on. See [debugging](#) for a list of debugger commands.

Enabling `debugmode` will not catch Lisp errors.

**ev** (*expr, arg\_1, ..., arg\_n*) Function

Evaluates the expression *expr* in the environment specified by the arguments *arg\_1*, ..., *arg\_n*. The arguments are switches (Boolean flags), assignments, equations, and functions. `ev` returns the result (another expression) of the evaluation.

The evaluation is carried out in steps, as follows.

1. First the environment is set up by scanning the arguments which may be any or all of the following.
  - `simp` causes *expr* to be simplified regardless of the setting of the switch `simp` which inhibits simplification if `false`.
  - `noeval` suppresses the evaluation phase of `ev` (see step (4) below). This is useful in conjunction with the other switches and in causing *expr* to be resimplified without being reevaluated.
  - `expand` causes expansion.



- `expand (m, n)` causes expansion, setting the values of `maxposex` and `maxnegex` to `m` and `n` respectively.
- `detout` causes any matrix inverses computed in `expr` to have their determinant kept outside of the inverse rather than dividing through each element.
- `diff` causes all differentiations indicated in `expr` to be performed.
- `derivlist (x, y, z, ...)` causes only differentiations with respect to the indicated variables.
- `float` causes non-integral rational numbers to be converted to floating point.
- `numer` causes some mathematical functions (including exponentiation) with numerical arguments to be evaluated in floating point. It causes variables in `expr` which have been given numerals to be replaced by their values. It also sets the `float` switch on.
- `pred` causes predicates (expressions which evaluate to `true` or `false`) to be evaluated.
- `eval` causes an extra post-evaluation of `expr` to occur. (See step (5) below.)
- `A` where `A` is an atom declared to be an evaluation flag (see `evflag`) causes `A` to be bound to `true` during the evaluation of `expr`.
- `V: expression` (or alternately `V=expression`) causes `V` to be bound to the value of `expression` during the evaluation of `expr`. Note that if `V` is a Maxima option, then `expression` is used for its value during the evaluation of `expr`. If more than one argument to `ev` is of this type then the binding is done in parallel. If `V` is a non-atomic expression then a substitution rather than a binding is performed.
- `F` where `F`, a function name, has been declared to be an evaluation function (see `evfun`) causes `F` to be applied to `expr`.
- Any other function names (e.g., `sum`) cause evaluation of occurrences of those names in `expr` as though they were verbs.
- In addition a function occurring in `expr` (say `F(x)`) may be defined locally for the purpose of this evaluation of `expr` by giving `F(x) := expression` as an argument to `ev`.
- If an atom not mentioned above or a subscripted variable or subscripted expression was given as an argument, it is evaluated and if the result is an equation or assignment then the indicated binding or substitution is performed. If the result is a list then the members of the list are treated as if they were additional arguments given to `ev`. This permits a list of equations to be given (e.g. `[X=1, Y=A**2]`) or a list of names of equations (e.g., `[%t1, %t2]` where `%t1` and `%t2` are equations) such as that returned by `solve`.

The arguments of `ev` may be given in any order with the exception of substitution equations which are handled in sequence, left to right, and evaluation functions which are composed, e.g., `ev (expr, ratsimp, realpart)` is handled as `realpart (ratsimp (expr))`.

The `simp`, `numer`, `float`, and `pred` switches may also be set locally in a block, or globally in Maxima so that they will remain in effect until being reset.

- If *expr* is a canonical rational expression (CRE), then the expression returned by *ev* is also a CRE, provided the `numer` and `float` switches are not both `true`.
- During step (1), a list is made of the non-subscripted variables appearing on the left side of equations in the arguments or in the value of some arguments if the value is an equation. The variables (subscripted variables which do not have associated array functions as well as non-subscripted variables) in the expression *expr* are replaced by their global values, except for those appearing in this list. Usually, *expr* is just a label or % (as in %i2 in the example below), so this step simply retrieves the expression named by the label, so that *ev* may work on it.
  - If any substitutions are indicated by the arguments, they are carried out now.
  - The resulting expression is then re-evaluated (unless one of the arguments was `noeval`) and simplified according to the arguments. Note that any function calls in *expr* will be carried out after the variables in it are evaluated and that *ev*(*F*(*x*)) thus may behave like *F*(*ev*(*x*)).
  - If one of the arguments was `eval`, steps (3) and (4) are repeated.

## Examples

```
(%i1) sin(x) + cos(y) + (w+1)^2 + 'diff (sin(w), w);
                                d
(%o1)          cos(y) + sin(x) + -- (sin(w)) + (w + 1)
                                dw
(%i2) ev (% , sin, expand, diff, x=2, y=1);
                                2
(%o2)          cos(w) + w  + 2 w + cos(1) + 1.909297426825682
```

An alternate top level syntax has been provided for *ev*, whereby one may just type in its arguments, without the *ev*( ). That is, one may write simply

```
expr, arg_1, ..., arg_n
```

This is not permitted as part of another expression, e.g., in functions, blocks, etc.

Notice the parallel binding process in the following example.

```
(%i3) programmode: false;
(%o3)          false
(%i4) x+y, x: a+y, y: 2;
(%o4)          y + a + 2
(%i5) 2*x - 3*y = 3$
(%i6) -3*x + 2*y = -4$
(%i7) solve ([%o5, %o6]);
Solution

(%t7)          y = - -
                    1
                    5

(%t8)          x = -
                    6
                    5
(%o8)          [[%t7, %t8]]
(%i8) %o6, %o8;
```

```

(%o8)                                     - 4 = - 4
(%i9) x + 1/x > gamma (1/2);
(%o9)                                     1
                                     x + - > sqrt(%pi)
                                     x
(%i10) %, numer, x=1/2;
(%o10)                                     2.5 > 1.772453850905516
(%i11) %, pred;
(%o11)                                     true

```

**evflag**

property

Some Boolean flags have the `evflag` property. `ev` treats such flags specially. A flag with the `evflag` property will be bound to `true` during the execution of `ev` if it is mentioned in the call to `ev`. For example, `demoivre` and `ratfac` are bound to `true` during the call `ev (%, demoivre, ratfac)`.

The flags which have the `evflag` property are: `algebraic`, `cauchysum`, `demoivre`, `dotscrules`, `%emode`, `%enumer`, `exponentialize`, `exptisolate`, `factorflag`, `float`, `halfangles`, `infeval`, `isolate_wrt_times`, `keepfloat`, `letrat`, `listarith`, `logabs`, `logarc`, `logexpand`, `lognegint`, `lognumer`, `mipbranch`, `numer_pbranch`, `programmode`, `radexpand`, `ratalgdenom`, `ratfac`, `ratmx`, `ratsimpexpons`, `simp`, `simplsum`, `sumexpand`, and `trigexpand`.

The construct `:lisp (putprop '$foo t 'evflag)` gives the `evflag` property to the variable `foo`, so `foo` is bound to `true` during the call `ev (%, foo)`. Equivalently, `ev (%, foo:true)` has the same effect.

**evfun**

property

Some functions have the `evfun` property. `ev` treats such functions specially. A function with the `evfun` property will be applied during the execution of `ev` if it is mentioned in the call to `ev`. For example, `ratsimp` and `radcan` will be applied during the call `ev (%, ratsimp, radcan)`.

The functions which have the `evfun` property are: `bfloat`, `factor`, `fullratsimp`, `logcontract`, `polarform`, `radcan`, `ratexpand`, `ratsimp`, `rectform`, `rootscontract`, `trigexpand`, and `trigreduce`.

The construct `:lisp (putprop '$foo t 'evfun)` gives the `evfun` property to the function `foo`, so that `foo` is applied during the call `ev (%, foo)`. Equivalently, `ev (%, (ev (%)))` has the same effect.

**infeval**

special symbol

Enables "infinite evaluation" mode. `ev` repeatedly evaluates an expression until it stops changing. To prevent a variable, say `X`, from being evaluated away in this mode, simply include `X='X` as an argument to `ev`. Of course expressions such as `ev (X, X=X+1, infeval)` will generate an infinite loop.

<b>kill</b> ( <i>symbol_1, ..., symbol_n</i> )	Function
<b>kill</b> ( <i>labels</i> )	Function
<b>kill</b> ( <i>clabels, dlabels, elabels</i> )	Function
<b>kill</b> ( <i>n</i> )	Function
<b>kill</b> ( <i>[m, n]</i> )	Function
<b>kill</b> ( <i>values, functions, arrays, ...</i> )	Function
<b>kill</b> ( <i>all</i> )	Function
<b>kill</b> ( <i>allbut (symbol_1, ..., symbol_n)</i> )	Function

Removes all bindings (value, function, array, or rule) from the arguments *symbol\_1, ..., symbol\_n*. An argument may be a single array element or subscripted function.

Several special arguments are recognized. Different kinds of arguments may be combined, e.g., `kill (clabels, functions, allbut (foo, bar))`.

`kill (labels)` unbinds all input, output, and intermediate expression labels created so far. `kill (clabels)` unbinds only input labels which begin with the current value of `inchar`. Likewise, `kill (dlabels)` unbinds only output labels which begin with the current value of `outchar`, and `kill (elabels)` unbinds only intermediate expression labels which begin with the current value of `linechar`.

`kill (n)`, where *n* is an integer, unbinds the *n* most recent input and output labels.

`kill ([m, n])` unbinds input and output labels *m* through *n*.

`kill (infolist)`, where *infolist* is any item in `infolists` (such as `values`, `functions`, or `arrays`) unbinds all items in *infolist*. See also `infolists`.

`kill (all)` unbinds all items on all `infolists`.

`kill (allbut (symbol_1, ..., symbol_n))` unbinds all items on all `infolists` except for *symbol\_1, ..., symbol\_n*. `kill (allbut (infolist))` unbinds all items except for the ones on *infolist*, where *infolist* is `values`, `functions`, `arrays`, etc.

The memory taken up by a bound property is not released until all symbols are unbound from it. In particular, to release the memory taken up by the value of a symbol, one unbinds the output label which shows the bound value, as well as unbinding the symbol itself.

`kill` quotes its arguments. The double single quotes operator, `''`, defeats the quotation.

`kill (symbol)` unbinds all properties of *symbol*. In contrast, `remvalue`, `remfunction`, `remarray`, and `remrule` unbind a specific property.

`kill` always returns `done`, even if an argument has no binding.

<b>labels</b> ( <i>symbol</i> )	Function
<b>labels</b>	Variable

Returns the list of input, output, or intermediate expression labels which begin with *symbol*. Typically *symbol* is the value of `inchar`, `outchar`, or `linechar`. The label character may be given with or without a percent sign, so, for example, `i` and `%i` yield the same result.

If no labels begin with *symbol*, `labels` returns an empty list.

The function `labels` quotes its argument. The double single quotes operator `''` defeats quotation. For example, `labels (''inchar)` returns the input labels which begin with the current input label character.

The variable `labels` is the list of input, output, and intermediate expression labels, including all previous labels if `inchar`, `outchar`, or `linechar` were redefined.

By default, Maxima displays the result of each user input expression, giving the result an output label. The output display is suppressed by terminating the input with `$` (dollar sign) instead of `;` (semicolon). An output label is generated, but not displayed, and the label may be referenced in the same way as displayed output labels. See also `%`, `%%`, and `%th`.

Intermediate expression labels can be generated by some functions. The flag `programmode` controls whether `solve` and some other functions generate intermediate expression labels instead of returning a list of expressions. Some other functions, such as `ldisplay`, always generate intermediate expression labels.

`first (rest (labels ('inchar)))` returns the most recent input label.

See also `inchar`, `outchar`, `linechar`, and `infolists`.

**linenum** Variable

The line number of the current pair of input and output expressions.

**myoptions** Variable

Default value: `[]`

`myoptions` is the list of all options ever reset by the user, whether or not they get reset to their default value.

**nolabels** Variable

Default value: `false`

When `nolabels` is `true`, input and output labels are generated but not appended to `labels`, the list of all input and output labels. `kill (labels)` kills the labels on the `labels` list, but does not kill any labels generated since `nolabels` was assigned `true`. It seems likely this behavior is simply broken.

See also `batch`, `batchload`, and `labels`.

**optionset** Variable

Default value: `false`

When `optionset` is `true`, Maxima prints out a message whenever a Maxima option is reset. This is useful if the user is doubtful of the spelling of some option and wants to make sure that the variable he assigned a value to was truly an option variable.

**playback** `()` Function

**playback** `(n)` Function

**playback** `([m, n])` Function

**playback** `([m])` Function

**playback** `(input)` Function

**playback** `(slow)` Function

**playback** `(time)` Function

**playback** `(grind)` Function

Displays input, output, and intermediate expressions, without recomputing them.

`playback` only displays the expressions bound to labels; any other output (such as

text printed by `print` or `describe`, or error messages) is not displayed. See also `labels`.

`playback` quotes its arguments. The double single quotes operator, `''`, defeats quotation. `playback` always returns `done`.

`playback ()` (with no arguments) displays all input, output, and intermediate expressions generated so far. An output expression is displayed even if it was suppressed by the `$` terminator when it was originally computed.

`playback (n)` displays the most recent  $n$  expressions. Each input, output, and intermediate expression counts as one.

`playback ([m, n])` displays input, output, and intermediate expressions with numbers from  $m$  through  $n$ , inclusive.

`playback ([m])` is equivalent to `playback ([m, m])`; this usually prints one pair of input and output expressions.

`playback (input)` displays all input expressions generated so far.

`playback (slow)` pauses between expressions and waits for the user to press `enter`. This behavior is similar to `demo`. `playback (slow)` is useful in conjunction with `save` or `stringout` when creating a secondary-storage file in order to pick out useful expressions.

`playback (time)` displays the computation time for each expression.

`playback (grind)` displays input expressions in the same format as the `grind` function. Output expressions are not affected by the `grind` option. See `grind`.

Arguments may be combined, e.g., `playback ([5, 10], grind, time, slow)`.

<b>printprops</b> ( $a, i$ )	Function
<b>printprops</b> ( $[a_1, \dots, a_n], i$ )	Function
<b>printprops</b> ( $all, i$ )	Function

Displays the property with the indicator  $i$  associated with the atom  $a$ .  $a$  may also be a list of atoms or the atom `all` in which case all of the atoms with the given property will be used. For example, `printprops ([f, g], atvalue)`. `printprops` is for properties that cannot otherwise be displayed, i.e. for `atvalue`, `atomgrad`, `gradef`, and `matchdeclare`.

<b>prompt</b>	Variable
---------------	----------

Default value: `_`

`prompt` is the prompt symbol of the `demo` function, `playback (slow)` mode, and the Maxima break loop (as invoked by `break`).

<b>quit</b> ()	Function
----------------	----------

Terminates the Maxima session. Note that the function must be invoked as `quit()`; or `quit()$`, not `quit` by itself.

To stop a lengthy computation, type `control-C`. The default action is to return to the Maxima prompt. If `*debugger-hook*` is `nil`, `control-C` opens the Lisp debugger. See also `debugging`.

- remfunction** (*f<sub>1</sub>*, ..., *f<sub>n</sub>*) Function  
**remfunction** (*all*) Function  
 Removes the user defined functions *f<sub>1</sub>*, ..., *f<sub>n</sub>* from Maxima. **remfunction** (**all**) removes all functions.
- reset** () Function  
 Resets many global variables and options, and some other variables, to their default values.  
**reset** processes the variables on the Lisp list *\*variable-initial-values\**. The Lisp macro **defmvar** puts variables on this list (among other actions). Many, but not all, global variables and options are defined by **defmvar**, and some variables defined by **defmvar** are not global variables or options.
- showtime** Variable  
 Default value: **false**  
 When **showtime** is **true**, the computation time and elapsed time is printed with each output expression.  
 See also **time**, **timer**, and **playback**.
- sstatus** (*feature*, *package*) Function  
 Sets the status of *feature* in *package*. After **sstatus** (*feature*, *package*) is executed, **status** (*feature*, *package*) returns **true**. This can be useful for package writers, to keep track of what features they have loaded in.
- to\_lisp** () Function  
 Enters the Lisp system under Maxima. (**to-maxima**) returns to Maxima.
- values** Variable  
 Initial value: []  
**values** is a list of all bound user variables (not Maxima options or switches). The list comprises symbols bound by **:**, **::**, or **:=**.

## 5 Operators

### 5.1 nary

An nary operator is used to denote a function of any number of arguments, each of which is separated by an occurrence of the operator, e.g.  $A+B$  or  $A+B+C$ . The `nary("x")` function is a syntax extension function to declare `x` to be an nary operator. Functions may be declared to be nary. If `declare(j,nary);` is done, this tells the simplifier to simplify, e.g. `j(j(a,b),j(c,d))` to `j(a, b, c, d)`.

See also `syntax`.

### 5.2 nofix

`nofix` operators are used to denote functions of no arguments. The mere presence of such an operator in a command will cause the corresponding function to be evaluated. For example, when one types `"exit;"` to exit from a Maxima break, `"exit"` is behaving similar to a `nofix` operator. The function `nofix("x")` is a syntax extension function which declares `x` to be a `nofix` operator.

See also `syntax`.

### 5.3 operator

See `operators`.

### 5.4 postfix

`postfix` operators like the `prefix` variety denote functions of a single argument, but in this case the argument immediately precedes an occurrence of the operator in the input string, e.g. `3!`. The `postfix("x")` function is a syntax extension function to declare `x` to be a `postfix` operator.

See also `syntax`.

### 5.5 prefix

A `prefix` operator is one which signifies a function of one argument, which argument immediately follows an occurrence of the operator. `prefix("x")` is a syntax extension function to declare `x` to be a `prefix` operator.

See also `syntax`.



## 5.6 Definitions for Operators

"!"

operator

The factorial operator. For any complex number  $x$  (including integer, rational, and real numbers) except for negative integers,  $x!$  is defined as `gamma(x+1)`.

For an integer  $x$ ,  $x!$  simplifies to the product of the integers from 1 to  $x$  inclusive.  $0!$  simplifies to 1. For a floating point number  $x$ ,  $x!$  simplifies to the value of `gamma(x+1)`. For  $x$  equal to  $n/2$  where  $n$  is an odd integer,  $x!$  simplifies to a rational factor times `sqrt(%pi)` (since `gamma(1/2)` is equal to `sqrt(%pi)`). If  $x$  is anything else,  $x!$  is not simplified.

The variables `factlim`, `minfactorial`, and `factcomb` control the simplification of expressions containing factorials.

The functions `gamma`, `bffac`, and `cbffac` are varieties of the gamma function. `makegamma` substitutes `gamma` for factorials and related functions.

See also `binomial`.

- The factorial of an integer, half-integer, or floating point argument is simplified unless the operand is greater than `factlim`.

```
(%i1) factlim: 10$
(%i2) [0!, (7/2)!, 4.77!, 8!, 20!];
      105 sqrt(%pi)
(%o2) [1, -----, 81.44668037931193, 40320, 20!]
      16
```

- The factorial of a complex number, known constant, or general expression is not simplified. Even so it may be possible simplify the factorial after evaluating the operand.

```
(%i1) [(%i + 1)!, %pi!, %e!, (cos(1) + sin(1))!];
(%o1) [(%i + 1)!, %pi!, %e!, (sin(1) + cos(1))!]
(%i2) ev (%i, numer, %enumer);
(%o2) [(%i + 1)!, 7.188082728976031, 4.260820476357003,
                                           1.227580202486819]
```

- The factorial of an unbound symbol is not simplified.

```
(%i1) kill (foo)$
(%i2) foo!;
(%o2)                                     foo!
```

- Factorials are simplified, not evaluated. Thus  $x!$  may be replaced even in a quoted expression.

```
(%i1) '([0!, (7/2)!, 4.77!, 8!, 20!]);
      105 sqrt(%pi)
(%o1) [1, -----, 81.44668037931193, 40320, 20!]
      16
```

"!!"

operator

The double factorial operator.

For an integer, float, or rational number  $n$ ,  $n!!$  evaluates to the product  $n (n-2) (n-4) (n-6) \dots (n - 2 (k-1))$  where  $k$  is equal to `entier (n/2)`, that is, the largest integer less than or equal to  $n/2$ . Note that this definition does not coincide with other published definitions for arguments which are not integers.

For an even (or odd) integer  $n$ ,  $n!!$  evaluates to the product of all the consecutive even (or odd) integers from 2 (or 1) through  $n$  inclusive.

For an argument  $n$  which is not an integer, float, or rational,  $n!!$  yields a noun form `genfact (n, n/2, 2)`.

- "#"** operator  
The logical operator "Not equals".
- "."**  operator  
The dot operator, for matrix (non-commutative) multiplication. When "." is used in this way, spaces should be left on both sides of it, e.g. `A . B`. This distinguishes it plainly from a decimal point in a floating point number.  
See also `dot`, `dot0nscsimp`, `dot0simp`, `dot1simp`, `dotassoc`, `dotconstrules`, `dotdistrib`, `dotexptsimp`, `dotident`, and `dotscrules`.
- ":"**  operator  
The assignment operator. E.g. `A:3` sets the variable `A` to 3.
- "::"**  operator  
Assignment operator. `::` assigns the value of the expression on its right to the value of the quantity on its left, which must evaluate to an atomic variable or subscripted variable.
- "::="**  operator  
The `::="` is used instead of `:=` to indicate that what follows is a macro definition, rather than an ordinary functional definition. See `macros`.
- ":="**  operator  
The function definition operator. E.g. `f(x):=sin(x)` defines a function `f`.
- "="**  operator  
denotes an equation to Maxima. To the pattern matcher in Maxima it denotes a total relation that holds between two expressions if and only if the expressions are syntactically identical.
- and**  operator  
The logical conjunction operator. `and` is an n-ary infix operator; its operands are Boolean expressions, and its result is a Boolean value.  
`and` forces evaluation (like `is`) of one or more operands, and may force evaluation of all operands.

Operands are evaluated in the order in which they appear. **and** evaluates only as many of its operands as necessary to determine the result. If any operand is **false**, the result is **false** and no further operands are evaluated.

The global flag **prederror** governs the behavior of **and** when an evaluated operand cannot be determined to be **true** or **false**. **and** prints an error message when **prederror** is **true**. Otherwise, **and** returns **unknown**.

**and** is not commutative: **a and b** might not be equal to **b and a** due to the treatment of indeterminate operands.

**or** operator

The logical disjunction operator. **or** is an n-ary infix operator; its operands are Boolean expressions, and its result is a Boolean value.

**or** forces evaluation (like **is**) of one or more operands, and may force evaluation of all operands.

Operands are evaluated in the order in which they appear. **or** evaluates only as many of its operands as necessary to determine the result. If any operand is **true**, the result is **true** and no further operands are evaluated.

The global flag **prederror** governs the behavior of **or** when an evaluated operand cannot be determined to be **true** or **false**. **or** prints an error message when **prederror** is **true**. Otherwise, **or** returns **unknown**.

**or** is not commutative: **a or b** might not be equal to **b or a** due to the treatment of indeterminate operands.

**not** operator

The logical negation operator. **not** is a prefix operator; its operand is a Boolean expression, and its result is a Boolean value.

**not** forces evaluation (like **is**) of its operand.

The global flag **prederror** governs the behavior of **not** when its operand cannot be determined to be **true** or **false**. **not** prints an error message when **prederror** is **true**. Otherwise, **not** returns **unknown**.

**abs** (*expr*) Function

Returns the absolute value *expr*. If *expr* is complex, returns the complex modulus of *expr*.

**additive** special symbol

If **declare(f,additive)** has been executed, then:

(1) If **f** is univariate, whenever the simplifier encounters **f** applied to a sum, **f** will be distributed over that sum. I.e. **f(y+x)** will simplify to **f(y)+f(x)**.

(2) If **f** is a function of 2 or more arguments, additivity is defined as additivity in the first argument to **f**, as in the case of **sum** or **integrate**, i.e. **f(h(x)+g(x),x)** will simplify to **f(h(x),x)+f(g(x),x)**. This simplification does not occur when **f** is applied to expressions of the form **sum(x[i],i,lower-limit,upper-limit)**.

**allbut**

keyword

works with the `part` commands (i.e. `part`, `inpart`, `substpart`, `substinpart`, `dpart`, and `lpart`). For example,

```
(%i1) expr: e+d+c+b+a$
(%i2) part (expr, [2, 5]);
(%o2)                d + a
```

while

```
(%i3) part (expr, allbut (2, 5));
(%o3)                e + c + b
```

It also works with the `kill` command,

```
kill (allbut (name_1, ..., name_k))
```

will do a `kill (all)` except it will not kill the names specified. Note: `name_i` means a name such as function name such as `u`, `f`, `foo`, or `g`, not an infolist such as functions.

**antisymmetric**

declaration

If `declare(h,antisymmetric)` is done, this tells the simplifier that `h` is antisymmetric. E.g. `h(x,z,y)` will simplify to  $-h(x, y, z)$ . That is, it will give  $(-1)^n$  times the result given by `symmetric` or `commutative`, where  $n$  is the number of interchanges of two arguments necessary to convert it to that form.

**cabs** (*expr*)

Function

Returns the complex absolute value (the complex modulus) of *expr*.

**commutative**

declaration

If `declare(h,commutative)` is done, this tells the simplifier that `h` is a commutative function. E.g. `h(x,z,y)` will simplify to `h(x, y, z)`. This is the same as `symmetric`.

**entier** (*x*)

Function

Returns the largest integer less than or equal to *x* where *x* is numeric. `fix` (as in `fixnum`) is a synonym for this, so `fix(x)` is precisely the same.

**equal** (*expr\_1*, *expr\_2*)

Function

Used with an `is`, returns `true` (or `false`) if and only if *expr\_1* and *expr\_2* are equal (or not equal) for all possible values of their variables (as determined by `ratsimp`). Thus `is (equal ((x + 1)^2, x^2 + 2*x + 1))` returns `true` whereas if *x* is unbound `is ((x + 1)^2 = x^2 + 2*x + 1)` returns `false`. Note also that `is (rat(0)=0)` yields `false` but `is (equal (rat(0), 0))` yields `true`.

If a determination can't be made, then `is (equal (a, b))` returns a simplified but equivalent expression, whereas `is (a=b)` always returns either `true` or `false`.

All variables occurring in *expr\_1* and *expr\_2* are presumed to be real valued.

`ev (expr, pred)` is equivalent to `is (expr)`.

```
(%i1) is (x^2 >= 2*x - 1);
(%o1) true
(%i2) assume (a > 1);
(%o2) [a > 1]
(%i3) is (log (log (a+1) + 1) > 0 and a^2 + 1 > 2*a);
(%o3) true
```

**eval** operator  
 As an argument in a call to `ev (expr)`, `eval` causes an extra evaluation of `expr`. See `ev`.

**evenp** (`expr`) Function  
 Returns `true` if `expr` is an even integer. `false` is returned in all other cases.

**fix** (`x`) Function  
 A synonym for `entier (x)`.

**fullmap** (`f, expr_1, ...`) Function  
 Similar to `map`, but `fullmap` keeps mapping down all subexpressions until the main operators are no longer the same.  
`fullmap` is used by the Maxima simplifier for certain matrix manipulations; thus, Maxima sometimes generates an error message concerning `fullmap` even though `fullmap` was not explicitly called by the user.

```
(%i1) a + b*c$
(%i2) fullmap (g, %);
(%o2) g(b) g(c) + g(a)
(%i3) map (g, %th(2));
(%o3) g(b c) + g(a)
```

**fullmapl** (`f, list_1, ...`) Function  
 Similar to `fullmap`, but `fullmapl` only maps onto lists and matrices.

```
(%i1) fullmapl ("+", [3, [4, 5]], [[a, 1], [0, -1.5]]);
(%o1) [[a + 3, 4], [4, 3.5]]
```

**is** (`expr`) Function  
 Attempts to determine whether `expr` (which must evaluate to a predicate) is provable from the facts in the current data base. `is` returns `true` if the predicate is true for all values of its variables consistent with the data base and returns `false` if it is false for all such values. Otherwise, its action depends on the setting of the switch `prederror`. `is` errs out if the value of `prederror` is `true` and returns `unknown` if `prederror` is `false`.

**isqrt** (`x`) Function  
 Returns the "integer square root" of the absolute value of `x`, which is an integer.

- max** (*x\_1*, *x\_2*, ...) Function  
Returns the maximum of its arguments (or returns a simplified form if some of its arguments are non-numeric).
- min** (*x\_1*, *x\_2*, ...) Function  
Returns the minimum of its arguments (or returns a simplified form if some of its arguments are non-numeric).
- mod** (*p*) Function  
**mod** (*p*, *m*) Function  
Converts the polynomial *p* to a modular representation with respect to the current modulus which is the value of the variable `modulus`.  
`mod` (*p*, *m*) specifies a modulus *m* to be used instead of the current value of `modulus`.  
See `modulus`.
- oddp** (*expr*) Function  
is `true` if *expr* is an odd integer. `false` is returned in all other cases.
- pred** operator  
As an argument in a call to `ev` (*expr*), `pred` causes predicates (expressions which evaluate to `true` or `false`) to be evaluated. See `ev`.
- make\_random\_state** (*n*) Function  
**make\_random\_state** (*s*) Function  
**make\_random\_state** (*true*) Function  
**make\_random\_state** (*false*) Function  
A random state object represents the state of the random number generator. The state comprises 627 32-bit words.  
`make_random_state` (*n*) returns a new random state object created from an integer seed value equal to *n* modulo  $2^{32}$ . *n* may be negative.  
`make_random_state` (*s*) returns a copy of the random state *s*.  
`make_random_state` (`true`) returns a new random state object, using the current computer clock time as the seed.  
`make_random_state` (`false`) returns a copy of the current state of the random number generator.
- set\_random\_state** (*s*) Function  
Copies *s* to the random number generator state.  
`set_random_state` always returns `done`.
- random** (*x*) Function  
Returns a pseudorandom number. If *x* is an integer, `random` (*x*) returns an integer from 0 through *x* - 1 inclusive. If *x* is a floating point number, `random` (*x*) returns a nonnegative floating point number less than *x*. `random` complains with an error if *x* is neither an integer nor a float, or if *x* is not positive.

The functions `make_random_state` and `set_random_state` maintain the state of the random number generator.

The Maxima random number generator is an implementation of the Mersenne twister MT 19937.

Examples:

```
(%i1) s1: make_random_state (654321)$
(%i2) set_random_state (s1);
(%o2) done
(%i3) random (1000);
(%o3) 768
(%i4) random (9573684);
(%o4) 7657880
(%i5) random (2^75);
(%o5) 11804491615036831636390
(%i6) s2: make_random_state (false)$
(%i7) random (1.0);
(%o7) .2310127244107132
(%i8) random (10.0);
(%o8) 4.394553645870825
(%i9) random (100.0);
(%o9) 32.28666704056853
(%i10) set_random_state (s2);
(%o10) done
(%i11) random (1.0);
(%o11) .2310127244107132
(%i12) random (10.0);
(%o12) 4.394553645870825
(%i13) random (100.0);
(%o13) 32.28666704056853
```

**sign** (*expr*) Function  
 Attempts to determine the sign of *expr* on the basis of the facts in the current data base. It returns one of the following answers: **pos** (positive), **neg** (negative), **zero**, **pz** (positive or zero), **nz** (negative or zero), **pn** (positive or negative), or **pnz** (positive, negative, or zero, i.e. nothing known).

**signum** (*x*) Function  
 For numeric *x*, returns 0 if *x* is 0, otherwise returns -1 or +1 as *x* is less than or greater than 0, respectively.  
 If *x* is not numeric then a simplified but equivalent form is returned. For example, **signum**(-*x*) gives **-signum**(*x*).

**sort** (*list*, *p*) Function  
**sort** (*list*) Function  
 Sorts *list* according to a predicate *p* of two arguments, such as "<" or **orderlessp**.  
**sort** (*list*) sorts *list* according to Maxima's built-in ordering.  
*list* may contain numeric or nonnumeric items, or both.

- sqrt** (*x*) Function  
 The square root of *x*. It is represented internally by  $x^{(1/2)}$ . See also `rootscontract`.  
`radexpand` if `true` will cause *n*th roots of factors of a product which are powers of *n* to be pulled outside of the radical, e.g. `sqrt(16*x^2)` will become `4*x` only if `radexpand` is `true`.
- sqrtdispflag** Variable  
 Default value: `true`  
 When `sqrtdispflag` is `false`, causes `sqrt` to display with exponent `1/2`.
- sublis** (*list*, *expr*) Function  
 Makes multiple parallel substitutions into an expression.  
 The variable `sublis_apply_lambda` controls simplification after `sublis`.  
 Example:  

```
(%i1) sublis ([a=b, b=a], sin(a) + cos(b));
(%o1) sin(b) + cos(a)
```
- sublist** (*list*, *p*) Function  
 Returns the list of elements of *list* for which the predicate *p* returns `true`.  
 Example:  

```
(%i1) L: [1, 2, 3, 4, 5, 6]$
(%i2) sublist (L, evenp);
(%o2) [2, 4, 6]
```
- sublis\_apply\_lambda** Variable  
 default: `true` - controls whether `lambda`'s substituted are applied in simplification after `sublis` is used or whether you have to do an `ev` to get things to apply. `true` means do the application.
- subst** (*a*, *b*, *c*) Function  
 Substitutes *a* for *b* in *c*. *b* must be an atom or a complete subexpression of *c*. For example, `x+y+z` is a complete subexpression of `2*(x+y+z)/w` while `x+y` is not. When *b* does not have these characteristics, one may sometimes use `substpart` or `ratsubst` (see below). Alternatively, if *b* is of the form `e/f` then one could use `subst(a*f, e, c)` while if *b* is of the form `e^(1/f)` then one could use `subst(a^f, e, c)`. The `subst` command also discerns the  $x^y$  in  $x^{-y}$  so that `subst(a, sqrt(x), 1/sqrt(x))` yields `1/a`. *a* and *b* may also be operators of an expression enclosed in double-quotes " or they may be function names. If one wishes to substitute for the independent variable in derivative forms then the `at` function (see below) should be used.  
`subst` is an alias for `substitute`.  
`subst(eq_1, expr)` or `subst([eq_1, ..., eq_k], expr)` are other permissible forms. The `eq_i` are equations indicating substitutions to be made. For each equation, the right side will be substituted for the left in the expression *expr*.



exptsubst if true permits substitutions like  $y$  for  $e^x$  in  $e^{ax}$  to take place.

When opsubst is false, subst will not attempt to substitute into the operator of an expression. E.g. (opsubst: false, subst (x<sup>2</sup>, r, r+r[0])) will work.

Examples:

```
(%i1) subst (a, x+y, x + (x+y)^2 + y);
      2
      y + x + a
(%o1)
(%i2) subst (-%i, %i, a + b*%i);
(%o2)      a - %i b
```

For further examples, do `example (subst)`.

**substinpart** ( $x, expr, n_1, \dots, n_k$ ) Function

Similar to `substpart`, but `substinpart` works on the internal representation of `expr`.

```
(%i1) x . 'diff (f(x), x, 2);
      2
      D
(%o1)      x . --- (f(x))
      2
      dx
(%i2) substinpart (d^2, %, 2);
      2
      x . d
(%o2)
(%i3) substinpart (f1, f[1](x+1), 0);
(%o3)      f1(x + 1)
```

If the last argument to a part function is a list of indices then several subexpressions are picked out, each one corresponding to an index of the list. Thus

```
(%i1) part (x+y+z, [1, 3]);
(%o1)      z + x
```

`piece` holds the value of the last expression selected when using the part functions. It is set during the execution of the function and thus may be referred to in the function itself as shown below. If `partswitch` is set to `true` then `end` is returned when a selected part of an expression doesn't exist, otherwise an error message is given.

```
(%i1) expr: 27*y^3 + 54*x*y^2 + 36*x^2*y + y + 8*x^3 + x + 1;
      3      2      2      3
(%o1)      27 y + 54 x y + 36 x y + y + 8 x + x + 1
(%i2) part (expr, 2, [1, 3]);
      2
(%o2)      54 y
(%i3) sqrt (piece/54);
(%o3)      abs(y)
(%i4) substpart (factor (piece), expr, [1, 2, 3, 5]);
      3
(%o4)      (3 y + 2 x) + y + x + 1
(%i5) expr: 1/x + y/x - 1/z;
      1      y      1
(%o5)      - - + - + -
```

```
(%i6) substpart (xthru (piece), expr, [2, 3]);
(%o6)
      z  x  x
      y + 1  1
      ----- - -
           x  z
```

Also, setting the option `inflag` to `true` and calling `part` or `substpart` is the same as calling `inpart` or `substinpart`.

**substpart** (*x*, *expr*, *n\_1*, ..., *n\_k*) Function

Substitutes *x* for the subexpression picked out by the rest of the arguments as in `part`. It returns the new value of *expr*. *x* may be some operator to be substituted for an operator of *expr*. In some cases *x* needs to be enclosed in double-quotes " (e.g. `substpart ("+", a*b, 0)` yields `b + a`).

```
(%i1) 1/(x^2 + 2);
(%o1)
      1
      -----
      2
      x  + 2
(%i2) substpart (3/2, %, 2, 1, 2);
(%o2)
      1
      -----
      3/2
      x  + 2
(%i3) a*x + f (b, y);
(%o3)
      a x + f(b, y)
(%i4) substpart ("+", %, 1, 0);
(%o4)
      x + f(b, y) + a
```

Also, setting the option `inflag` to `true` and calling `part` or `substpart` is the same as calling `inpart` or `substinpart`.

**subvarp** (*expr*) Function

Returns `true` if *expr* is a subscripted variable, for example `a[i]`.

**symbolp** (*expr*) Function

Returns `true` if *expr* is a symbol, else `false`. In effect, `symbolp(x)` is equivalent to the predicate `atom(x)` and not `numberp(x)`.

**unorder** () Function

Disables the aliasing created by the last use of the ordering commands `ordergreat` and `orderless`. `ordergreat` and `orderless` may not be used more than one time each without calling `unorder`. See also `ordergreat` and `orderless`.

```
(%i1) unorder();
(%o1)
      []
(%i2) b*x + a^2;
(%o2)
      2
      b x + a
(%i3) ordergreat (a);
```

```

(%o3) done
(%i4) b*x + a^2;
(%o4) a^2 + b x
(%i5) %th(1) - %th(3);
(%o5) a^2 - a^2
(%i6) unorder();
(%o6) [a]

```

**vectorpotential** (*givencurl*)

Function

Returns the vector potential of a given curl vector, in the current coordinate system. **potentialzeroloc** has a similar role as for **potential**, but the order of the left-hand sides of the equations must be a cyclic permutation of the coordinate variables.

**xthru** (*expr*)

Function

Combines all terms of *expr* (which should be a sum) over a common denominator without expanding products and exponentiated sums as **ratsimp** does. **xthru** cancels common factors in the numerator and denominator of rational expressions but only if the factors are explicit.

Sometimes it is better to use **xthru** before **ratsimping** an expression in order to cause explicit factors of the gcd of the numerator and denominator to be canceled thus simplifying the expression to be **ratsimped**.

```

(%i1) ((x+2)^20 - 2*y)/(x+y)^20 + (x+y)^(-19) - x/(x+y)^20;
(%o1)
      1          (x + 2)  - 2 y          x
----- + ----- - -----
      19          20          20
      (y + x)      (y + x)      (y + x)
(%i2) xthru (%);
(%o2)
      20
      (x + 2)  - y
-----
      20
      (y + x)

```

**zeroequiv** (*expr, v*)

Function

Tests whether the expression *expr* in the variable *v* is equivalent to zero, returning true, false, or dontknow.

**zeroequiv** has these restrictions:

1. Do not use functions that Maxima does not know how to differentiate and evaluate.
2. If the expression has poles on the real line, there may be errors in the result (but this is unlikely to occur).
3. If the expression contains functions which are not solutions to first order differential equations (e.g. Bessel functions) there may be incorrect results.

4. The algorithm uses evaluation at randomly chosen points for carefully selected subexpressions. This is always a somewhat hazardous business, although the algorithm tries to minimize the potential for error.

For example `zeroequiv (sin(2*x) - 2*sin(x)*cos(x), x)` returns `true` and `zeroequiv (%e^x + x, x)` returns `false`. On the other hand `zeroequiv (log(a*b) - log(a) - log(b), a)` returns `dontknow` because of the presence of an extra parameter `b`.



## 6 Expressions

### 6.1 Introduction to Expressions

There are a number of reserved words which cannot be used as variable names. Their use would cause a possibly cryptic syntax error.

integrate	next	from	diff
in	at	limit	sum
for	and	elseif	then
else	do	or	if
unless	product	while	thru
step			

Most things in Maxima are expressions. A sequence of expressions can be made into an expression by separating them by commas and putting parentheses around them. This is similar to the **C** *comma expression*.

```
(%i1) x: 3$
(%i2) (x: x+1, x: x^2);
(%o2) 16
(%i3) (if (x > 17) then 2 else 4);
(%o3) 4
(%i4) (if (x > 17) then x: 2 else y: 4, y+x);
(%o4) 20
```

Even loops in Maxima are expressions, although the value they return is the not too useful done.

```
(%i1) y: (x: 1, for i from 1 thru 10 do (x: x*i))$
(%i2) y;
(%o2) done
```

whereas what you really want is probably to include a third term in the *comma expression* which actually gives back the value.

```
(%i3) y: (x: 1, for i from 1 thru 10 do (x: x*i), x)$
(%i4) y;
(%o4) 3628800
```

### 6.2 Assignment

There are two assignment operators in Maxima, `:` and `::`. E.g., `a: 3` sets the variable `a` to 3. `::` assigns the value of the expression on its right to the value of the quantity on its left, which must evaluate to an atomic variable or subscripted variable.

### 6.3 Complex

A complex expression is specified in Maxima by adding the real part of the expression to `%i` times the imaginary part. Thus the roots of the equation  $x^2 - 4x + 13 = 0$  are  $2 + 3\%i$  and  $2 - 3\%i$ . Note that simplification of products of complex expressions can be effected by expanding the product. Simplification of quotients, roots, and other functions

of complex expressions can usually be accomplished by using the `realpart`, `imagpart`, `rectform`, `polarform`, `abs`, `carg` functions.

## 6.4 Inequality

Maxima has the usual inequality operators:

```
less than: <
greater than: >
greater than or equal to: >=
less than or equal to: <=
```

## 6.5 Syntax

It is possible to add new operators to Maxima (infix, prefix, postfix, unary, or matchfix with given precedences), to remove existing operators, or to redefine the precedence of existing operators. While Maxima's syntax should be adequate for most ordinary applications, it is possible to define new operators or eliminate predefined ones that get in the user's way. The extension mechanism is rather straightforward and should be evident from the examples below.

```
(%i1) prefix ("ddx")$
(%i2) ddx y$ /* equivalent to "ddx"(y) */
(%i3) infix ("<-")$
(%i4) a <- ddx y$ /* equivalent to "<-"(a, "ddx"(y)) */
```

For each of the types of operator except `special`, there is a corresponding creation function that will give the lexeme specified the corresponding parsing properties. Thus `prefix ("ddx")` will make `ddx` a prefix operator just like `-` or `not`. Of course, certain extension functions require additional information such as the matching keyword for a matchfix operator. In addition, binding powers and parts of speech must be specified for all keywords defined. This is done by passing additional arguments to the extension functions. If a user does not specify these additional parameters, Maxima will assign default values. The six extension functions with binding powers and parts of speech defaults (enclosed in brackets) are summarized below.

```
prefix (operator, rbp[180], rpos[any], pos[any])
postfix (operator, lbp[180], lpos[any], pos[any])
infix (operator, lbp[180], rbp[180], lpos[any], rpos[any], pos[any])
nary (operator, bp[180], argpos[any], pos[any])
nofix (operator, pos[any])
matchfix (operator, match, argpos[any], pos[any])
```

The defaults have been provided so that a user who does not wish to concern himself with parts of speech or binding powers may simply omit those arguments to the extension functions. Thus the following are all equivalent.

```
prefix ("ddx", 180, any, any)$
prefix ("ddx", 180)$
prefix ("ddx")$
```

It is also possible to remove the syntax properties of an operator by using the functions `remove` or `kill`. Specifically, `remove ("ddx", op)` or `kill ("ddx")` will return `ddx` to operand status; but in the second case all the other properties of `ddx` will also be removed.

```
(%i1) prefix ("ddx", 180, any, any)$
(%i2) ddx yz;
(%o2)                                     yz + 4
(%i3) "ddx"(u) := u+4;
(%o3)                                     ddx u := u + 4
(%i4) ddx 8;
(%o4)                                     12
```

## 6.6 Definitions for Expressions

**at** (*expr*, *list*) Function

Evaluates *expr* (which may be any expression) with the variables assuming the values as specified for them in the list of equations or the single equation similar to that given to the `atvalue` function. If a subexpression depends on any of the variables in list but it hasn't had an `atvalue` specified and it can't be evaluated then a noun form of the `at` will be returned which will display in a two-dimensional form. `example ("at")` displays some examples of `at`.

**box** (*expr*) Function

**box** (*expr*, *label*) Function

Returns *expr* enclosed in a box. The box is actually part of the expression. `box (expr, label)` encloses *expr* in a labelled box. *label* is a name which will be truncated in display if it is too long.

`boxchar` is the character used to draw the box in this and in the `dpart` and `lpart` functions.

**boxchar** Variable

Default value: "

`boxchar` is the character used to draw the box in the `box` and in the `dpart` and `lpart` functions.

**carg** (*z*) Function

Returns the complex argument of *z*. The complex argument is an angle `theta` in  $[-\pi, \pi]$  such that  $r \exp(i\theta) = z$  where *r* is the magnitude of *z*.

See also `abs` (complex magnitude), `polarform`, `rectform`, `realpart`, and `imagpart`.

**constant** special operator

Makes *ai* a constant as is `%pi`.

**constantp** (*expr*) Function

Returns `true` if *expr* is a constant (i.e. composed of numbers and `%pi`, `%e`, `%i` or any variables bound to a constant or declared constant by `declare`) else `false`. Any function whose arguments are constant is also considered to be a constant.



**declare** (*a<sub>1</sub>*, *f<sub>1</sub>*, *a<sub>2</sub>*, *f<sub>2</sub>*, ...) Function

Assigns the atom *a<sub>i</sub>* the flag *f<sub>i</sub>*. The *a<sub>i</sub>*'s and *f<sub>i</sub>*'s may also be lists of atoms and flags respectively in which case each of the atoms gets all of the properties.

The possible flags and their meanings are:

`constant` makes *a<sub>i</sub>* a constant as is `%pi`.

`mainvar` makes *a<sub>i</sub>* a `mainvar`. The ordering scale for atoms: numbers < constants (e.g. `%e`, `%pi`) < scalars < other variables < mainvars.

`scalar` makes *a<sub>i</sub>* a scalar.

`nonscalar` makes *a<sub>i</sub>* behave as does a list or matrix with respect to the dot operator.

`noun` makes the function *a<sub>i</sub>* a noun so that it won't be evaluated automatically.

`evfun` makes *a<sub>i</sub>* known to the `ev` function so that it will get applied if its name is mentioned. See `evfun`.

`evflag` makes *a<sub>i</sub>* known to the `ev` function so that it will be bound to `true` during the execution of `ev` if it is mentioned. See `evflag`.

`bindtest` causes *a<sub>i</sub>* to signal an error if it ever is used in a computation unbound.

Maxima currently recognizes and uses the following features of objects:

`even`, `odd`, `integer`, `rational`, `irrational`, `real`, `imaginary`,  
and `complex`

The useful features of functions include:

`increasing`,  
`decreasing`, `oddfun` (odd function), `evenfun` (even function),  
`commutative` (or `symmetric`), `antisymmetric`, `lassociative` and  
`rassociative`

The *a<sub>i</sub>* and *f<sub>i</sub>* may also be lists of objects or features. `featurep` (`object`, `feature`) determines if an object has been declared to have *feature*. See also `features`.

**disolate** (*expr*, *x<sub>1</sub>*, ..., *x<sub>n</sub>*) Function

is similar to `isolate` (*expr*, *x*) except that it enables the user to isolate more than one variable simultaneously. This might be useful, for example, if one were attempting to change variables in a multiple integration, and that variable change involved two or more of the integration variables. This function is autoloaded from `'simplification/disol.mac'`. A demo is available by `demo("disol")$`.

**dispform** (*expr*) Function

Returns the external representation of *expr* with respect to its main operator. This should be useful in conjunction with `part` which also deals with the external representation. Suppose *expr* is `-A`. Then the internal representation of *expr* is `"*(-1,A)`, while the external representation is `-"(A)`. `dispform` (*expr*, `all`) converts the entire expression (not just the top-level) to external format. For example, if `expr: sin(sqrt(x))`, then `freeof(sqrt, expr)` and `freeof(sqrt, dispform(expr))` give `true`, while `freeof(sqrt, dispform(expr, all))` gives `false`.

**distrib** (*expr*)

Function

Distributes sums over products. It differs from `expand` in that it works at only the top level of an expression, i.e., it doesn't recurse and it is faster than `expand`. It differs from `multthru` in that it expands all sums at that level.

Examples:

```
(%i1) distrib ((a+b) * (c+d));
(%o1)          b d + a d + b c + a c
(%i2) multthru ((a+b) * (c+d));
(%o2)          (b + a) d + (b + a) c
(%i3) distrib (1/((a+b) * (c+d)));
(%o3)          1
              -----
              (b + a) (d + c)
(%i4) expand (1/((a+b) * (c+d)), 1, 0);
(%o4)          1
              -----
              b d + a d + b c + a c
```

**dpart** (*expr, n\_1, ..., n\_k*)

Function

Selects the same subexpression as `part`, but instead of just returning that subexpression as its value, it returns the whole expression with the selected subexpression displayed inside a box. The box is actually part of the expression.

```
(%i1) dpart (x+y/z^2, 1, 2, 1);
(%o1)          y
              ---- + x
              2
              ""
              "z"
              ""
```

**exp** (*x*)

Function

The exponential function. Instances of `exp (x)` in input are simplified to `%e^x`; `exp` does not appear in simplified expressions.

`demoivre` if `true` will cause `%e^(a + b %i)` to become `%e^(a (cos(b) + %i sin(b)))` if `b` is free of `%i`. See `demoivre`.

`%emode`, when `true`, causes `%e^(%pi %i x)` to be simplified. See `%emode`.

`%enumer`, when `true` will cause `%e` to be replaced by 2.718... whenever `numer` is `true`. See `%enumer`.

**%emode**

Variable

Default value: `true`

When `%emode` is `true`, `%e^(%pi %i x)` is simplified as follows: it is expressed as `cos (%pi x) + %i sin (%pi x)` if `x` is an integer or a multiple of 1/2, 1/3, 1/4, or 1/6, and thus further simplified. For other numerical `x`, it is expressed as `%e^(%pi %i y)` where `y` is `x - 2 k` for some integer `k` such that `abs(y) < 1`.

When `%emode` is `false`, no special simplification of `%e^(%pi %i x)` is carried out.

**%enumer** Variable

Default value: `false`

When `%enumer` is `true`, `%e` is replaced by its numeric value 2.718... whenever `numer` is `true`.

When `%enumer` is `false`, this substitution is carried out only if the exponent in `%ex` evaluates to a number.

See also `ev` and `numer`.

**exptisolate** Variable

Default value: `false`

When `exptisolate` is `true`, will cause `isolate (expr, var)` to examine exponents of atoms (like `%e`) which contain `var`.

**exptsubst** Variable

Default value: `false`

When `exptsubst` is `true`, permits substitutions such as `y` for `%ex` in `%e(a x)` to take place.

**freeof** (*x<sub>1</sub>, ..., x<sub>n</sub>, expr*) Function

`freeof (x1, expr)` returns `true` if no subexpression of `expr` is equal to `x1` or if `x1` occurs only as a dummy variable in `expr`, and returns `false` otherwise.

`freeof (x1, ..., xn, expr)` is equivalent to `freeof (x1, expr)` and ... and `freeof (xn, expr)`.

The arguments `x1, ..., xn` may be names of functions and variables, subscripted names, operators (enclosed in double quotes), or general expressions. `freeof` evaluates its arguments.

`freeof` operates only on `expr` as it stands (after simplification and evaluation) and does not attempt to determine if some equivalent expression would give a different result. In particular, simplification may yield an equivalent but different expression which comprises some different elements than the original form of `expr`.

A variable is a dummy variable in an expression if it has no binding outside of the expression. Dummy variables recognized by `freeof` are the index of a sum or product, the limit variable in `limit`, the integration variable in the definite integral form of `integrate`, the original variable in `laplace`, formal variables in `at` expressions, and arguments in `lambda` expressions. Local variables in `block` are not recognized by `freeof` as dummy variables; this is a bug.

The indefinite form of `integrate` is *not* free of its variable of integration.

- Arguments are names of functions, variables, subscripted names, operators, and expressions. `freeof (a, b, expr)` is equivalent to `freeof (a, expr)` and `freeof (b, expr)`.

```
(%i1) expr: z^3 * cos (a[1]) * b^(c+d);
              d + c 3
(%o1)          cos(a ) b      z
              1
```

```
(%i2) freeof (z, expr);
(%o2)                                     false
(%i3) freeof (cos, expr);
(%o3)                                     false
(%i4) freeof (a[1], expr);
(%o4)                                     false
(%i5) freeof (cos (a[1]), expr);
(%o5)                                     false
(%i6) freeof (b^(c+d), expr);
(%o6)                                     false
(%i7) freeof ("^", expr);
(%o7)                                     false
(%i8) freeof (w, sin, a[2], sin (a[2]), b*(c+d), expr);
(%o8)                                     true
```

- freeof evaluates its arguments.

```
(%i1) expr: (a+b)^5$
(%i2) c: a$
(%i3) freeof (c, expr);
(%o3)                                     false
```

- freeof does not consider equivalent expressions. Simplification may yield an equivalent but different expression.

```
(%i1) expr: (a+b)^5$
(%i2) expand (expr);
(%o2)      5      4      2 3      3 2      4      5
      b + 5 a b + 10 a b + 10 a b + 5 a b + a
(%i3) freeof (a+b, %);
(%o3)                                     true
(%i4) freeof (a+b, expr);
(%o4)                                     false
(%i5) exp (x);
(%o5)                                     x
                                             %e
(%i6) freeof (exp, exp (x));
(%o6)                                     true
```

- A summation or definite integral is free of its dummy variable. An indefinite integral is not free of its variable of integration.

```
(%i1) freeof (i, 'sum (f(i), i, 0, n));
(%o1)                                     true
(%i2) freeof (x, 'integrate (x^2, x, 0, 1));
(%o2)                                     true
(%i3) freeof (x, 'integrate (x^2, x));
(%o3)                                     false
```

**genfact** ( $x, y, z$ ) Function  
 Returns the generalized factorial, defined as  $x (x-z) (x-2z) \dots (x-(y-1)z)$ .  
 Thus, for integral  $x$ ,  $\text{genfact}(x, x, 1) = x!$  and  $\text{genfact}(x, x/2, 2) = x!!$ .

**imagpart** (*expr*) Function  
Returns the imaginary part of the expression *expr*.

**indices** (*expr*) Function  
Returns a list of two elements. The first is a list of the free indices in *expr* (those that occur only once); the second is the list of dummy indices in *expr* (those that occur exactly twice).

**infix** (*op*) Function

**infix** (*op*, *lbp*, *rbp*) Function

**infix** (*op*, *lbp*, *rbp*, *lpos*, *rpos*, *pos*) Function

Declares *op* to be an infix operator. An infix operator is a function of two arguments, with the name of the function written between the arguments. For example, the subtraction operator `-` is an infix operator.

`infix (op)` declares *op* to be an infix operator with default binding powers (left and right both equal to 180) and parts of speech (left and right both equal to `any`).

`infix (op, lbp, rbp)` declares *op* to be an infix operator with stated left and right binding powers and default parts of speech (left and right both equal to `any`).

`infix (op, lbp, rbp, lpos, rpos, pos)` declares *op* to be an infix operator with stated left and right binding powers and parts of speech.

The precedence of *op* with respect to other operators derives from the left and right binding powers of the operators in question. If the left and right binding powers of *op* are both greater the left and right binding powers of some other operator, then *op* takes precedence over the other operator. If the binding powers are not both greater or less, some more complicated relation holds.

The associativity of *op* depends on its binding powers. Greater left binding power (*lbp*) implies an instance of *op* is evaluated before other operators to its left in an expression, while greater right binding power (*rbp*) implies an instance of *op* is evaluated before other operators to its right in an expression. Thus greater *lbp* makes *op* right-associative, while greater *rbp* makes *op* left-associative. If *lbp* is equal to *rbp*, *op* is left-associative.

See also `Syntax`.

Examples:

- If the left and right binding powers of *op* are both greater the left and right binding powers of some other operator, then *op* takes precedence over the other operator.

```
(%i1) "@"(a, b) := sconcat("(", a, ",", b, ")")$
(%i2) :lisp (get '$+ 'lbp)
100
(%i2) :lisp (get '$+ 'rbp)
100
(%i2) infix ("@", 101, 101)$
(%i3) 1 + a@b + 2;
(%o3) (a,b) + 3
(%i4) infix ("@", 99, 99)$
```

```
(%i5) 1 + a@b + 2;
(%o5) (a+1,b+2)
```

- Greater *lbp* makes *op* right-associative, while greater *rbp* makes *op* left-associative.

```
(%i1) "@"(a, b) := sconcat("(", a, ",", b, ")")$
(%i2) infix ("@", 100, 99)$
(%i3) foo @ bar @ baz;
(%o3) (foo,(bar,baz))
(%i4) infix ("@", 100, 101)$
(%i5) foo @ bar @ baz;
(%o5) ((foo,bar),baz)
```

## inflag

Variable

Default value: false

When `inflag` is `true`, the functions for part extraction will look at the internal form of `expr`. Note that the simplifier re-orders expressions. Thus `first (x+y)` will be `x` if `inflag` is `true` and `Y` if `inflag` is `false`. (`first (y+x)` gives the same results). Also, setting `inflag` to `true` and calling `part/substpart` is the same as calling `inpart/substinpart`. Functions affected by the setting of `inflag` are: `part`, `substpart`, `first`, `rest`, `last`, `length`, the `for ... in` construct, `map`, `fullmap`, `maplist`, `reveal` and `pickapart`.

## inpart (expr, n.1, ..., n.k)

Function

is similar to `part` but works on the internal representation of the expression rather than the displayed form and thus may be faster since no formatting is done. Care should be taken with respect to the order of subexpressions in sums and products (since the order of variables in the internal form is often different from that in the displayed form) and in dealing with unary minus, subtraction, and division (since these operators are removed from the expression). `part (x+y, 0)` or `inpart (x+y, 0)` yield `+`, though in order to refer to the operator it must be enclosed in `"s`. For example `... if inpart (%o9,0) = "+" then ...`

Examples:

```
(%i1) x + y + w*z;
(%o1) w z + y + x
(%i2) inpart (%o1, 3, 2);
(%o2) z
(%i3) part (%th (2), 1, 2);
(%o3) z
(%i4) 'limit (f(x)^g(x+1), x, 0, minus);
(%o4) limit f(x)
      x -> 0-
      g(x + 1)
(%i5) inpart (%o4, 1, 2);
(%o5) g(x + 1)
```

**isolate** (*expr*, *x*) Function

Returns *expr* with subexpressions which are sums and which do not contain *var* replaced by intermediate expression labels (these being atomic symbols like %t1, %t2, ...). This is often useful to avoid unnecessary expansion of subexpressions which don't contain the variable of interest. Since the intermediate labels are bound to the subexpressions they can all be substituted back by evaluating the expression in which they occur.

`exptisolate` (default value: `false`) if `true` will cause `isolate` to examine exponents of atoms (like %e) which contain *var*.

`isolate_wrt_times` if `true`, then `isolate` will also isolate wrt products. See `isolate_wrt_times`.

Do `example (isolate)` for examples.

**isolate\_wrt\_times** Variable

Default value: `false`

When `isolate_wrt_times` is `true`, `isolate` will also isolate wrt products. E.g. compare both settings of the switch on

```
(%i1) isolate_wrt_times: true$
(%i2) isolate (expand ((a+b+c)^2), c);

(%t2)
                2 a

(%t3)
                2 b

(%t4)
                2          2
                b  + 2 a b + a

(%o4)
                2
                c  + %t3 c + %t2 c + %t4
(%i4) isolate_wrt_times: false$
(%i5) isolate (expand ((a+b+c)^2), c);

(%o5)
                2
                c  + 2 b c + 2 a c + %t4
```

**listconstvars** Variable

Default value: `false`

When `listconstvars` is `true`, it will cause `listofvars` to include %e, %pi, %i, and any variables declared constant in the list it returns if they appear in the expression `listofvars` is called on. The default is to omit these.

**listdummyvars** Variable

Default value: `true`

When `listdummyvars` is `false`, "dummy variables" in the expression will not be included in the list returned by `listofvars`. (The meaning of "dummy variables" is

as given in `freeof`. "Dummy variables" are mathematical things like the index of a sum or product, the limit variable, and the definite integration variable.) Example:

```
(%i1) listdummyvars: true$
(%i2) listofvars ('sum(f(i), i, 0, n));
(%o2)          [i, n]
(%i3) listdummyvars: false$
(%i4) listofvars ('sum(f(i), i, 0, n));
(%o4)          [n]
```

**listofvars** (*expr*) Function

Returns a list of the variables in *expr*.

`listconstvars` if `true` causes `listofvars` to include `%e`, `%pi`, `%i`, and any variables declared constant in the list it returns if they appear in *expr*. The default is to omit these.

```
(%i1) listofvars (f (x[1]+y) / g^(2+a));
(%o1)          [g, a, x , y]
              1
```

**lfreeof** (*list*, *expr*) Function

For each member *m* of *list*, calls `freeof` (*m*, *expr*). It returns `false` if any call to `freeof` does and `true` otherwise.

**lopow** (*expr*, *x*) Function

Returns the lowest exponent of *x* which explicitly appears in *expr*. Thus

```
(%i1) lopow ((x+y)^2 + (x+y)^a, x+y);
(%o1)          min(a, 2)
```

**lpart** (*label*, *expr*, *n*<sub>1</sub>, ..., *n*<sub>*k*</sub>) Function

is similar to `dpart` but uses a labelled box. A labelled box is similar to the one produced by `dpart` but it has a name in the top line.

**multthru** (*expr*) Function

**multthru** (*expr*<sub>1</sub>, *expr*<sub>2</sub>) Function

Multiplies a factor (which should be a sum) of *expr* by the other factors of *expr*. That is, *expr* is *f*<sub>1</sub> *f*<sub>2</sub> ... *f*<sub>*n*</sub> where at least one factor, say *f*<sub>*i*</sub>, is a sum of terms. Each term in that sum is multiplied by the other factors in the product. (Namely all the factors except *f*<sub>*i*</sub>). `multthru` does not expand exponentiated sums. This function is the fastest way to distribute products (commutative or noncommutative) over sums. Since quotients are represented as products `multthru` can be used to divide sums by products as well.

`multthru` (*expr*<sub>1</sub>, *expr*<sub>2</sub>) multiplies each term in *expr*<sub>2</sub> (which should be a sum or an equation) by *expr*<sub>1</sub>. If *expr*<sub>1</sub> is not itself a sum then this form is equivalent to `multthru` (*expr*<sub>1</sub>\**expr*<sub>2</sub>).

```
(%i1) x/(x-y)^2 - 1/(x-y) - f(x)/(x-y)^3;
(%o1)          - ---- + ----- - -----
              1          x          f(x)
```



```

          x - y      2      3
          (x - y)    (x - y)
(%i2) multthru ((x-y)^3, %);
          2
(%o2)      - (x - y) + x (x - y) - f(x)
(%i3) ratexpand (%);
          2
(%o3)      - y + x y - f(x)
(%i4) ((a+b)^10*s^2 + 2*a*b*s + (a*b)^2)/(a*b*s^2);
          10 2      2 2
          (b + a) s + 2 a b s + a b
(%o4)      -----
          2
          a b s
(%i5) multthru (%); /* note that this does not expand (b+a)^10 */
          10
          2 a b (b + a)
(%o5)      - + --- + -----
          s 2      a b
          s
(%i6) multthru (a.(b+c.(d+e)+f));
(%o6)      a . f + a . c . (e + d) + a . b
(%i7) expand (a.(b+c.(d+e)+f));
(%o7)      a . f + a . c . e + a . c . d + a . b

```

**nounify** (*f*)

Function

Returns the noun form of the function name *f*. This is needed if one wishes to refer to the name of a verb function as if it were a noun. Note that some verb functions will return their noun forms if they can't be evaluated for certain arguments. This is also the form returned if a function call is preceded by a quote.

**nterms** (*expr*)

Function

Returns the number of terms that *expr* would have if it were fully expanded out and no cancellations or combination of terms occurred. Note that expressions like `sin (expr)`, `sqrt (expr)`, `exp (expr)`, etc. count as just one term regardless of how many terms *expr* has (if it is a sum).

**op** (*expr*)

Function

Returns the operator of the expression, and functions the same way as `part (expr, 0)`. It observes the setting of the `inpart` flag.

**operatorp** (*expr*, *op*)

Function

**operatorp** (*expr*, [*op-1*, ..., *op-n*])

Function

`operatorp (expr, op)` returns `true` if *op* is equal to the operator of *expr*.

`operatorp (expr, [op-1, ..., op-n])` returns `true` if some element *op-1*, ..., *op-n* is equal to the operator of *expr*.

- optimize** (*expr*) Function  
 Returns an expression that produces the same value and side effects as *expr* but does so more efficiently by avoiding the recomputation of common subexpressions. **optimize** also has the side effect of "collapsing" its argument so that all common subexpressions are shared. Do `example (optimize)` for examples.
- optimprefix** Variable  
 Default value: %  
**optimprefix** is the prefix used for generated symbols by the **optimize** command.
- ordergreat** (*v\_1*, ..., *v\_n*) Function  
 Sets up aliases for the variables *v\_1*, ..., *v\_n* such that  $v_1 > v_2 > \dots > v_n$ , and  $v_n >$  any other variable not mentioned as an argument.  
 See also **orderless**.
- ordergreatp** (*expr\_1*, *expr\_2*) Function  
 Returns **true** if *expr\_2* precedes *expr\_1* in the ordering set up with the **ordergreat** function.
- orderless** (*v\_1*, ..., *v\_n*) Function  
 Sets up aliases for the variables *v\_1*, ..., *v\_n* such that  $v_1 < v_2 < \dots < v_n$ , and  $v_n <$  any other variable not mentioned as an argument.  
 Thus the complete ordering scale is: numerical constants < declared constants < declared scalars < first argument to **orderless** < ... < last argument to **orderless** < variables which begin with A < ... < variables which begin with Z < last argument to **ordergreat** < ... < first argument to **ordergreat** < declared **mainvars**.  
 See also **ordergreat** and **mainvar**.
- orderlessp** (*expr\_1*, *expr\_2*) Function  
 Returns **true** if *expr\_1* precedes *expr\_2* in the ordering set up by the **orderless** command.
- part** (*expr*, *n\_1*, ..., *n\_k*) Function  
 Returns parts of the displayed form of **expr**. It obtains the part of **expr** as specified by the indices *n\_1*, ..., *n\_k*. First part *n\_1* of **expr** is obtained, then part *n\_2* of that, etc. The result is part *n\_k* of ... part *n\_2* of part *n\_1* of **expr**.  
**part** can be used to obtain an element of a list, a row of a matrix, etc.  
 If the last argument to a **part** function is a list of indices then several subexpressions are picked out, each one corresponding to an index of the list. Thus **part** ( $x + y + z$ , [1, 3]) is  $z+x$ .  
**piece** holds the last expression selected when using the **part** functions. It is set during the execution of the function and thus may be referred to in the function itself as shown below.  
 If **partswitch** is set to **true** then **end** is returned when a selected part of an expression doesn't exist, otherwise an error message is given.

Example: `part (z+2*y, 2, 1)` yields 2.

`example (part)` displays additional examples.

### **partition** (*expr*, *x*)

Function

Returns a list of two expressions. They are (1) the factors of *expr* (if it is a product), the terms of *expr* (if it is a sum), or the list (if it is a list) which don't contain *var* and, (2) the factors, terms, or list which do.

```
(%i1) partition (2*a*x*f(x), x);
(%o1)          [2 a, x f(x)]
(%i2) partition (a+b, x);
(%o2)          [b + a, 0]
(%i3) partition ([a, b, f(a), c], a);
(%o3)          [[b, c], [a, f(a)]]
```

### **partswitch**

Variable

Default value: `false`

When `partswitch` is `true`, `end` is returned when a selected part of an expression doesn't exist, otherwise an error message is given.

### **pickapart** (*expr*, *n*)

Function

Assigns intermediate expression labels to subexpressions of *expr* at depth *n*, an integer. Subexpressions at greater or lesser depths are not assigned labels. `pickapart` returns an expression in terms of intermediate expressions equivalent to the original expression *expr*.

See also `part`, `dpart`, `lpart`, `inpart`, and `reveal`.

Examples:

```
(%i1) expr: (a+b)/2 + sin (x^2)/3 - log (1 + sqrt(x+1));
(%o1)          - log(sqrt(x + 1) + 1) +  $\frac{\sin(x^2)}{3}$  +  $\frac{b + a}{2}$ 
(%i2) pickapart (expr, 0);
(%t2)          - log(sqrt(x + 1) + 1) +  $\frac{\sin(x^2)}{3}$  +  $\frac{b + a}{2}$ 
(%o2)          %t2
(%i3) pickapart (expr, 1);
(%t3)          - log(sqrt(x + 1) + 1)
(%t4)           $\frac{\sin(x^2)}{\quad\quad\quad}$ 
```

```

3

(%t5)          b + a
              -----
                2

(%o5)          %t5 + %t4 + %t3
(%i5) pickapart (expr, 2);

(%t6)          log(sqrt(x + 1) + 1)

(%t7)          2
              sin(x )

(%t8)          b + a

(%o8)          %t8  %t7
              ---- + ---- - %t6
                2    3

(%i8) pickapart (expr, 3);

(%t9)          sqrt(x + 1) + 1

(%t10)         2
              x

(%o10)         b + a          sin(%t10)
              ----- - log(%t9) + -----
                2                3

(%i10) pickapart (expr, 4);

(%t11)         sqrt(x + 1)

(%o11)         2
              sin(x )  b + a
              ----- + ----- - log(%t11 + 1)
                3        2

(%i11) pickapart (expr, 5);

(%t12)         x + 1

(%o12)         2
              sin(x )  b + a
              ----- + ----- - log(sqrt(%t12) + 1)
                3        2

```

```
(%i12) pickapart (expr, 6);
```

```
(%o12)      2
            sin(x )  b + a
            ----- + ----- - log(sqrt(x + 1) + 1)
             3          2
```

**piece**

Variable

Holds the last expression selected when using the **part** functions. It is set during the execution of the function and thus may be referred to in the function itself.

**polarform** (*expr*)

Function

Returns an expression  $r \%e^{(i \text{ theta})}$  equivalent to *expr*, such that *r* and *theta* are purely real.

**powers** (*expr*, *x*)

Function

Gives the powers of *x* occurring in *expr*.

load (powers) loads this function.

**product** (*expr*, *i*, *i\_0*, *i\_1*)

Function

Returns the product of the values of *expr* as the index *i* varies from *i\_0* to *i\_1*. The evaluation is similar to that of **sum**.

If *i\_1* is one less than *i\_0*, the product is an "empty product" and **product** returns 1 rather than reporting an error. See also **prodhack**.

Maxima does not simplify products.

Example:

```
(%i1) product (x + i*(i+1)/2, i, 1, 4);
(%o1)      (x + 1) (x + 3) (x + 6) (x + 10)
```

**realpart** (*expr*)

Function

Returns the real part of *expr*. **realpart** and **imagpart** will work on expressions involving trigonometric and hyperbolic functions, as well as square root, logarithm, and exponentiation.

**rectform** (*expr*)

Function

Returns an expression  $a + b \%i$  equivalent to *expr*, such that *a* and *b* are purely real.

**rembox** (*expr*, *unlabelled*)

Function

**rembox** (*expr*, *label*)

Function

**rembox** (*expr*)

Function

Removes boxes from *expr*.

**rembox** (*expr*, *unlabelled*) removes all unlabelled boxes from *expr*.

**rembox** (*expr*, *label*) removes only boxes bearing *label*.

**rembox** (*expr*) removes all boxes, labelled and unlabelled.

**sum** (*expr*, *i*, *i\_0*, *i\_1*)

Function

Represents a summation of the values of *expr* as the index *i* varies from *i\_0* to *i\_1*. If the upper and lower limits differ by an integer then each term in the sum is evaluated and added together. Otherwise, if the `simplsum` is `true` the result is simplified. This simplification may sometimes be able to produce a closed form. Sums may be differentiated, added, subtracted, or multiplied with some automatic simplification being performed.

If `simplsum` is `false` or if `'sum` is used, the value is a sum noun form which is a representation of the sigma notation used in mathematics.

If *i\_1* is one less than *i\_0*, the sum is an "empty sum" and `sum` returns 0 rather than reporting an error. See also `sumhack`.

`cauchysum` when `true` causes the Cauchy product to be used when multiplying sums together rather than the usual product. In the Cauchy product the index of the inner summation is a function of the index of the outer one rather than varying independently.

`genindex` is the alphabetic prefix used to generate the next variable of summation.

`gensumnum` is the numeric suffix used to generate the next variable of summation. If it is set to `false` then the index will consist only of `genindex` with no numeric suffix.

See also `sumcontract`, `intosum`, `bashindices`, and `niceindices`.

`example (sum)` displays some examples.

**lsum** (*expr*, *i*, *list*)

Function

Represents the sum of *expr* for each element *i* in *list*.

```
(%i1) lsum (x^i, i, [1, 2, 7]);
              7      2
(%o1)          x  + x  + x
```

If the last element *list* argument does not evaluate, or does not evaluate to a Maxima list then the answer is left in noun form

```
(%i2) lsum (i^2, i, rootsof (x^3-1));
=====
          \      2
          >   i
          /
          =====
                          3
          i in rootsof(x  - 1)
```

**verb**

special symbol

`verb` is the opposite of "noun", i.e., a function form which "does something" ("action" - for most functions the usual case). E.g. `integrate` integrates a function, unless it is declared to be a "noun", in which case it represents the integral of the function.

See also `noun`, `nounify`, `verbify`.

**verbify** (*f*)

Function

Returns the function name *f* in its verb form.

See also `verb`, `noun`, and `nounify`.



## 7 Simplification

### 7.1 Definitions for Simplification

**askexp** Variable

When **asksign** is called, **askexp** is the expression **asksign** is testing.

At one time, it was possible for a user to inspect **askexp** by entering a Maxima break with control-A.

**askinteger** (*expr*, *integer*) Function

**askinteger** (*expr*) Function

**askinteger** (*expr*, *even*) Function

**askinteger** (*expr*, *odd*) Function

**askinteger** (*expr*, *integer*) attempts to determine from the **assume** database whether *expr* is an integer. **askinteger** prompts the user if it cannot tell otherwise, and attempt to install the information in the database if possible. **askinteger** (*expr*) is equivalent to **askinteger** (*expr*, *integer*).

**askinteger** (*expr*, *even*) and **askinteger** (*expr*, *odd*) likewise attempt to determine if *expr* is an even integer or odd integer, respectively.

**asksign** (*expr*) Function

First attempts to determine whether the specified expression is positive, negative, or zero. If it cannot, it asks the user the necessary questions to complete its deduction. The user's answer is recorded in the data base for the duration of the current computation. The return value of **asksign** is one of **pos**, **neg**, or **zero**.

**demoivre** (*expr*) Function

**demoivre** Variable

The function **demoivre** (*expr*) converts one expression without setting the global variable **demoivre**.

When the variable **demoivre** is **true**, complex exponentials are converted into equivalent expressions in terms of circular functions:  $\exp(a + b*i)$  simplifies to  $e^a * (\cos(b) + i*\sin(b))$  if *b* is free of *i*. *a* and *b* are not expanded.

The default value of **demoivre** is **false**.

**exponentialize** converts circular and hyperbolic functions to exponential form. **demoivre** and **exponentialize** cannot both be true at the same time.

**domain** Variable

Default value: **real**

When **domain** is set to **complex**, **sqrt** ( $x^2$ ) will remain **sqrt** ( $x^2$ ) instead of returning **abs**(*x*).



**expand** (*expr*) Function  
**expand** (*expr*, *p*, *n*) Function

Expand expression *expr*. Products of sums and exponentiated sums are multiplied out, numerators of rational expressions which are sums are split into their respective terms, and multiplication (commutative and non-commutative) are distributed over addition at all levels of *expr*.

For polynomials one should usually use `ratexpand` which uses a more efficient algorithm.

`maxnegex` and `maxposex` control the maximum negative and positive exponents, respectively, which will expand.

`expand` (*expr*, *p*, *n*) expands *expr*, using *p* for `maxposex` and *n* for `maxnegex`. This is useful in order to expand part but not all of an expression.

`expon` - the exponent of the largest negative power which is automatically expanded (independent of calls to `expand`). For example if `expon` is 4 then  $(x+1)^{-5}$  will not be automatically expanded.

`expop` - the highest positive exponent which is automatically expanded. Thus  $(x+1)^3$ , when typed, will be automatically expanded only if `expop` is greater than or equal to 3. If it is desired to have  $(x+1)^n$  expanded where *n* is greater than `expop` then executing `expand ((x+1)^n)` will work only if `maxposex` is not less than *n*.

The `expand` flag used with `ev` causes expansion.

The file `'simplification/facexp.mac'` contains several related functions (in particular `facsum`, `factorfacsum` and `collectterms`, which are autoloaded) and variables (`nextlayerfactor` and `facsum_combine`) that provide the user with the ability to structure expressions by controlled expansion. Brief function descriptions are available in `'simplification/facexp.usg'`. A demo is available by doing `demo("facexp")`.

**expandwrt** (*expr*, *x\_1*, ..., *x\_n*) Function

Expands expression *expr* with respect to the variables *x\_1*, ..., *x\_n*. All products involving the variables appear explicitly. The form returned will be free of products of sums of expressions that are not free of the variables. *x\_1*, ..., *x\_n* may be variables, operators, or expressions.

By default, denominators are not expanded, but this can be controlled by means of the switch `expandwrt_denom`.

This function is autoloaded from `'simplification/stopex.mac'`.

**expandwrt\_denom** Variable

Default value: `false`

`expandwrt_denom` controls the treatment of rational expressions by `expandwrt`. If `true`, then both the numerator and denominator of the expression will be expanded according to the arguments of `expandwrt`, but if `expandwrt_denom` is `false`, then only the numerator will be expanded in that way.

**expandwrt\_factored** (*expr*, *x\_1*, ..., *x\_n*) Function  
 is similar to `expandwrt`, but treats expressions that are products somewhat differently. `expandwrt_factored` expands only on those factors of `expr` that contain the variables *x\_1*, ..., *x\_n*.

This function is autoloaded from 'simplification/stopex.mac'.

**expon** Variable  
 Default value: 0  
`expon` is the exponent of the largest negative power which is automatically expanded (independent of calls to `expand`). For example, if `expon` is 4 then  $(x+1)^{-5}$  will not be automatically expanded.

**exponentialize** (*expr*) Function  
**exponentialize** Variable

The function `exponentialize (expr)` converts circular and hyperbolic functions in *expr* to exponentials, without setting the global variable `exponentialize`.

When the variable `exponentialize` is `true`, all circular and hyperbolic functions are converted to exponential form. The default value is `false`.

`demoivre` converts complex exponentials into circular functions. `exponentialize` and `demoivre` cannot both be true at the same time.

**expop** Variable  
 Default value: 0

`expop` is the highest positive exponent which is automatically expanded. Thus  $(x + 1)^3$ , when typed, will be automatically expanded only if `expop` is greater than or equal to 3. If it is desired to have  $(x + 1)^n$  expanded where *n* is greater than `expop` then executing `expand ((x + 1)^n)` will work only if `maxposex` is not less than *n*.

**factlim** Variable  
 Default value: -1

`factlim` specifies the highest factorial which is automatically expanded. If it is -1 then all integers are expanded.

**intosum** (*expr*) Function

Moves multiplicative factors outside a summation to inside. If the index is used in the outside expression, then the function tries to find a reasonable index, the same as it does for `sumcontract`. This is essentially the reverse idea of the `outative` property of summations, but note that it does not remove this property, it only bypasses it.

In some cases, a `scanmap (multthru, expr)` may be necessary before the `intosum`.

**lassociative** declaration

`declare (g, lassociative)` tells the Maxima simplifier that *g* is left-associative. E.g., `g (g (a, b), g (c, d))` will simplify to `g (g (a, b), c), d)`.

**linear** declaration

One of Maxima's operator properties. For univariate  $f$  so declared, "expansion"  $f(x + y)$  yields  $f(x) + f(y)$ ,  $f(a*x)$  yields  $a*f(x)$  takes place where  $a$  is a "constant". For functions of two or more arguments, "linearity" is defined to be as in the case of `sum` or `integrate`, i.e.,  $f(a*x + b, x)$  yields  $a*f(x, x) + b*f(1, x)$  for  $a$  and  $b$  free of  $x$ .

`linear` is equivalent to `additive` and `outative`. See also `opproperties`.

**mainvar** declaration

You may declare variables to be `mainvar`. The ordering scale for atoms is essentially: numbers < constants (e.g., `%e`, `%pi`) < scalars < other variables < mainvars. E.g., compare `expand((X+Y)^4)` with `(declare(x, mainvar), expand((x+y)^4))`. (Note: Care should be taken if you elect to use the above feature. E.g., if you subtract an expression in which  $x$  is a `mainvar` from one in which  $x$  isn't a `mainvar`, resimplification e.g. with `ev(expr, simp)` may be necessary if cancellation is to occur. Also, if you save an expression in which  $x$  is a `mainvar`, you probably should also save  $x$ .)

**maxapplydepth** Variable

Default value: 10000

`maxapplydepth` is the maximum depth to which `apply1` and `apply2` will delve.

**maxapplyheight** Variable

Default value: 10000

`maxapplyheight` is the maximum height to which `applyb1` will reach before giving up.

**maxnegex** Variable

Default value: 1000

`maxnegex` is the largest negative exponent which will be expanded by the `expand` command (see also `maxposex`).

**maxposex** Variable

Default value: 1000

`maxposex` is the largest exponent which will be expanded with the `expand` command (see also `maxnegex`).

**multiplicative** declaration

`declare(f, multiplicative)` tells the Maxima simplifier that  $f$  is multiplicative.

1. If  $f$  is univariate, whenever the simplifier encounters  $f$  applied to a product,  $f$  distributes over that product. E.g.,  $f(x*y)$  simplifies to  $f(x)*f(y)$ .
2. If  $f$  is a function of 2 or more arguments, multiplicativity is defined as multiplicativity in the first argument to  $f$ , e.g.,  $f(g(x) * h(x), x)$  simplifies to  $f(g(x), x) * f(h(x), x)$ .

This simplification does not occur when  $f$  is applied to expressions of the form `product(x[i], i, m, n)`.

**negdistrib**

Variable

Default value: `true`

When `negdistrib` is `true`, `-1` distributes over an expression. E.g.,  $-(x + y)$  becomes  $-y - x$ . Setting it to `false` will allow  $-(x + y)$  to be displayed like that. This is sometimes useful but be very careful: like the `simp` flag, this is one flag you do not want to set to `false` as a matter of course or necessarily for other than local use in your Maxima.

**negsumdispflag**

Variable

Default value: `true`

When `negsumdispflag` is `true`,  $x - y$  displays as  $x - y$  instead of as  $-y + x$ . Setting it to `false` causes the special check in display for the difference of two expressions to not be done. One application is that thus `a + %i*b` and `a - %i*b` may both be displayed the same way.

**noeval**

special symbol

`noeval` suppresses the evaluation phase of `ev`. This is useful in conjunction with other switches and in causing expressions to be resimplified without being reevaluated.

**noun**

declaration

`noun` is one of the options of the `declare` command. It makes a function so declared a "noun", meaning that it won't be evaluated automatically.

**noundisp**

Variable

Default value: `false`

When `noundisp` is `true`, nouns display with a single quote. This switch is always `true` when displaying function definitions.

**nouns**

special symbol

`nouns` is an `evflag`. When used as an option to the `ev` command, `nouns` converts all "noun" forms occurring in the expression being `ev`'d to "verbs", i.e., evaluates them. See also `noun`, `nounify`, `verb`, and `verbify`.

**numer**

special symbol

`numer` causes some mathematical functions (including exponentiation) with numerical arguments to be evaluated in floating point. It causes variables in `expr` which have been given numervals to be replaced by their values. It also sets the `float` switch on.

**numerval** (*x<sub>1</sub>*, *expr<sub>1</sub>*, ..., *var<sub>n</sub>*, *expr<sub>n</sub>*)

Function

Declares the variables `x1`, ..., `xn` to have numeric values equal to `expr1`, ..., `exprn`. The numeric value is evaluated and substituted for the variable in any expressions in which the variable occurs if the `numer` flag is `true`. See also `ev`.

The expressions `expr1`, ..., `exprn` can be any expressions, not necessarily numeric.

**opproperties**

Variable

`opproperties` is the list of the special operator properties recognized by the Maxima simplifier: `linear`, `additive`, `multiplicative`, `outative`, `evenfun`, `oddfun`, `commutative`, `symmetric`, `antisymmetric`, `nary`, `lassociative`, `rassociative`.

**opsubst**

Variable

Default value: `true`

When `opsubst` is `false`, `subst` does not attempt to substitute into the operator of an expression. E.g., `(opsubst: false, subst (x^2, r, r+r[0]))` will work.

**outative**

declaration

`declare (f, outative)` tells the Maxima simplifier that constant factors in the argument of `f` can be pulled out.

1. If `f` is univariate, whenever the simplifier encounters `f` applied to a product, that product will be partitioned into factors that are constant and factors that are not and the constant factors will be pulled out. E.g., `f(a*x)` will simplify to `a*f(x)` where `a` is a constant. Non-atomic constant factors will not be pulled out.
2. If `f` is a function of 2 or more arguments, outativity is defined as in the case of `sum` or `integrate`, i.e., `f(a*g(x), x)` will simplify to `a * f(g(x), x)` for a free of `x`.

`sum`, `integrate`, and `limit` are all outative.

**posfun**

declaration

`declare (f, posfun)` declares `f` to be a positive function. `is (f(x) > 0)` yields `true`.

**prodhack**

Variable

Default value: `false`

When `prodhack` is `true`, the identity `product (f(i), i, a, b) = 1/product (f(i), i, b+1, a-1)` is applied if `a` is greater than `b`. For example, `product (f(i), i, 3, 1)` yields `1/f(2)`.

**radcan** (*expr*)

Function

Simplifies *expr*, which can contain logs, exponentials, and radicals, by converting it into a form which is canonical over a large class of expressions and a given ordering of variables; that is, all functionally equivalent forms are mapped into a unique form. For a somewhat larger class of expressions, `radcan` produces a regular form. Two equivalent expressions in this class do not necessarily have the same appearance, but their difference can be simplified by `radcan` to zero.

For some expressions `radcan` is quite time consuming. This is the cost of exploring certain relationships among the components of the expression for simplifications based on factoring and partial-fraction expansions of exponents.

When `%e_to_numlog` is `true`, `%e^(r*log(expr))` simplifies to `expr^r` if `r` is a rational number.

When `radexpand` is `false`, certain transformations are inhibited. `radcan (sqrt (1-x))` remains `sqrt (1-x)` and is not simplified to `%i sqrt (x-1)`. `radcan (sqrt (x^2 - 2*x + 11))` remains `sqrt (x^2 - 2*x + 11)` and is not simplified to `x - 1`.

`example (radcan)` displays some examples.

## **radexpand**

Variable

Default value: `true`

`radexpand` controls some simplifications of radicals.

When `radexpand` is `all`, causes  $n$ th roots of factors of a product which are powers of  $n$  to be pulled outside of the radical. E.g. if `radexpand` is `all`, `sqrt (16*x^2)` simplifies to `4*x`.

More particularly, consider `sqrt (x^2)`.

- If `radexpand` is `all` or `assume (x > 0)` has been executed, `sqrt(x^2)` simplifies to `x`.
- If `radexpand` is `true` and `domain` is `real` (its default), `sqrt(x^2)` simplifies to `abs(x)`.
- If `radexpand` is `false`, or `radexpand` is `true` and `domain` is `complex`, `sqrt(x^2)` is not simplified.

Note that `domain` only matters when `radexpand` is `true`.

## **radsubstflag**

Variable

Default value: `false`

`radsubstflag`, if `true`, permits `ratsubst` to make substitutions such as `u` for `sqrt (x)` in `x`.

## **rassociative**

declaration

`declare (g, rassociative)` tells the Maxima simplifier that `g` is right-associative. E.g., `g(g(a, b), g(c, d))` simplifies to `g(a, g(b, g(c, d)))`.

## **scsimp** (*expr, rule\_1, ..., rule\_n*)

Function

Sequential Comparative Simplification (method due to Stoute). `scsimp` attempts to simplify `expr` according to the rules `rule_1, ..., rule_n`. If a smaller expression is obtained, the process repeats. Otherwise after all simplifications are tried, it returns the original answer.

`example (scsimp)` displays some examples.

## **simplsum**

Variable

Default value: `false`

When `simplsum` is `true`, the result of a `sum` is simplified. This simplification may sometimes be able to produce a closed form. If `simplsum` is `false` or if the quoted form `'sum` is used, the value is a sum noun form which is a representation of the sigma notation used in mathematics.

**sumcontract** (*expr*) Function

Combines all sums of an addition that have upper and lower bounds that differ by constants. The result is an expression containing one summation for each set of such summations added to all appropriate extra terms that had to be extracted to form this sum. **sumcontract** combines all compatible sums and uses one of the indices from one of the sums if it can, and then try to form a reasonable index if it cannot use any supplied.

It may be necessary to do an **intosum** (*expr*) before the **sumcontract**.

**sumexpand** Variable

Default value: **false**

When **sumexpand** is **true**, products of sums and exponentiated sums simplify to nested sums.

See also **cauchysum**.

Examples:

```
(%i1) sumexpand: true$
(%i2) sum (f (i), i, 0, m) * sum (g (j), j, 0, n);
      m      n
      ====  ====
      \      \
      >      >   f(i1) g(i2)
      /      /
      ====  ====
      i1 = 0 i2 = 0
(%i3) sum (f (i), i, 0, m)^2;
      m      m
      ====  ====
      \      \
      >      >   f(i3) f(i4)
      /      /
      ====  ====
      i3 = 0 i4 = 0
```

**sumhack** Variable

Default value: **false**

When **sumhack** is **true**, the identity  $\text{sum}(f(i), i, a, b) = -\text{sum}(f(i), i, b+1, a-1)$  is applied if  $a$  is greater than  $b$ . For example, **(sumhack: true, sum (f(i), i, 3, 1))** yields **-f(2)**.

**sumsplitfact** Variable

Default value: **true**

When **sumsplitfact** is **false**, **minfactorial** is applied after a **factcomb**.

**symmetric** declaration

**declare (h, symmetric)** tells the Maxima simplifier that **h** is a symmetric function.

E.g., **h (x, z, y)** simplifies to **h (x, y, z)**.

**commutative** is synonymous with **symmetric**.

**unknown** (*expr*)

Function

Returns `true` if and only if *expr* contains an operator or function not recognized by the Maxima simplifier.





## 8 Plotting

### 8.1 Definitions for Plotting

#### **in\_netmath**

Variable

Default value: `false`

When `in_netmath` is `true`, `plot3d` prints OpenMath output to the console if `plot_format` is `openmath`; otherwise `in_netmath` (even if `true`) has no effect.

`in_netmath` has no effect on `plot2d`.

#### **openplot\_curves** (*list, rest\_options*)

Function

Takes a list of curves such as

```
[x1, y1, x2, y2, ...], [u1, v1, u2, v2, ...], ..]
```

or

```
[[x1, y1], [x2, y2], ...], ...]
```

and plots them. This is similar to `xgraph_curves`, but uses the open plot routines. Additional symbol arguments may be given such as `"{xrange -3 4}"`. The following plots two curves, using big points, labeling the first one `jim` and the second one `jane`.

```
openplot_curves ([["{plotpoints 1} {pointsize 6} {label jim}
  {text {xaxislabel {joe is nice}}}],
  [1, 2, 3, 4, 5, 6, 7, 8],
  [{"label jane} {color pink }"], [3, 1, 4, 2, 5, 7]]);
```

Some other special keywords are `xfun`, `color`, `plotpoints`, `linecolors`, `pointsize`, `nolines`, `bargraph`, `labelposition`, `xaxislabel`, and `yaxislabel`.

#### **plot2d** (*expr, range, ..., options, ...*)

Function

#### **plot2d** (*parametric\_expr*)

Function

#### **plot2d** (*[expr<sub>1</sub>, ..., expr<sub>n</sub>], x\_range, y\_range*)

Function

#### **plot2d** (*[expr<sub>1</sub>, ..., expr<sub>n</sub>], x\_range*)

Function

#### **plot2d** (*expr, x\_range, y\_range*)

Function

#### **plot2d** (*expr, x\_range*)

Function

Displays a plot of one or more expressions as a function of one variable.

In all cases, *expr* is an expression to be plotted on the vertical axis as a function of one variable. *x\_range*, the range of the horizontal axis, is a list of the form [*variable*, *min*, *max*], where *variable* is a variable which appears in *expr*. *y\_range*, the range of the vertical axis, is a list of the form [*y*, *min*, *max*].

`plot2d (expr, x_range)` plots *expr* as a function of the variable named in *x\_range*, over the range specified in *x\_range*. If the vertical range is not otherwise specified by `set_plot_options`, it is chosen automatically. All options are assumed to have default values unless otherwise specified by `set_plot_options`.

`plot2d (expr, x_range, y_range)` plots *expr* as a function of the variable named in *x\_range*, over the range specified in *x\_range*. The vertical range is set to *y\_range*. All

options are assumed to have default values unless otherwise specified by `set_plot_options`.

`plot2d ([expr_1, ..., expr_n], x_range)` plots *expr\_1*, ..., *expr\_n* as a function of the variable named in *x\_range*, over the range specified in *x\_range*. If the vertical range is not otherwise specified by `set_plot_options`, it is chosen automatically. All options are assumed to have default values unless otherwise specified by `set_plot_options`.

`plot2d ([expr_1, ..., expr_n], x_range, y_range)` plots *expr\_1*, ..., *expr\_n* as a function of the variable named in *x\_range*, over the range specified in *x\_range*. The vertical range is set to *y\_range*. All options are assumed to have default values unless otherwise specified by `set_plot_options`.

Examples:

```
(%i1) plot2d (sin(x), [x, -5, 5])$
(%i2) plot2d (sec(x), [x, -2, 2], [y, -20, 20], [nticks, 200])$
```

Anywhere there may be an ordinary expression, there may be a parametric expression: *parametric\_expr* is a list of the form [*parametric*, *x\_expr*, *y\_expr*, *t\_range*, *options*]. Here *x\_expr* and *y\_expr* are expressions of 1 variable *var* which is the first element of the range *trange*. The plot is of the path traced out by the pair [*x\_expr*, *y\_expr*] as *var* varies in *trange*.

In the following example, we plot a circle, then we do the plot with only a few points used, so that we get a star, and finally we plot this together with an ordinary function of X.

Examples:

- Plot a circle with a parametric plot.

```
(%i1) plot2d ([parametric, cos(t), sin(t), [t, -%pi*2, %pi*2],
             [nticks, 80]])$
```

- Plot a star: join eight points on the circumference of a circle.

```
(%i2) plot2d ([parametric, cos(t), sin(t), [t, -%pi*2, %pi*2],
             [nticks, 8]])$
```

- Plot a cubic polynomial with an ordinary plot and a circle with a parametric plot.

```
(%i3) plot2d ([x^3 + 2, [parametric, cos(t), sin(t), [t, -5, 5],
             [nticks, 80]]], [x, -3, 3])$
```

See also `plot_options`, which describes plotting options and has more examples.

### **xgraph\_curves** (*list*)

Function

graphs the list of 'point sets' given in *list* by using `xgraph`.

A point set may be of the form

```
[x0, y0, x1, y1, x2, y2, ...]
```

or

```
[[x0, y0], [x1, y1], ...]
```

A point set may also contain symbols which give labels or other information.

```
xgraph_curves ([pt_set1, pt_set2, pt_set3]);
```

graph the three point sets as three curves.

```
pt_set: append (["NoLines: True", "LargePixels: true"], [x0, y0, x1, y1, ...])
```

would make the point set [and subsequent ones], have no lines between points, and to use large pixels. See the man page on xgraph for more options to specify.

```
pt_set: append ([concat ("\\", "x^2+y")], [x0, y0, x1, y1, ...]);
```

would make there be a "label" of "x<sup>2</sup>+y" for this particular point set. The " at the beginning is what tells xgraph this is a label.

```
pt_set: append ([concat ("TitleText: Sample Data")], [x0, ...])$
```

would make the main title of the plot be "Sample Data" instead of "Maxima Plot".

To make a bar graph with bars which are 0.2 units wide, and to plot two possibly different such bar graphs:

```
xgraph_curves ([append (["BarGraph: true", "NoLines: true", "BarWidth: .2"],
  create_list ([i - .2, i^2], i, 1, 3)),
  append (["BarGraph: true", "NoLines: true", "BarWidth: .2"],
  create_list ([i + .2, .7*i^2], i, 1, 3))]);
```

A temporary file 'xgraph-out' is used.

## plot\_options

Variable

Elements of this list state the default options for plotting. If an option is present in a `plot2d` or `plot3d` call, that value takes precedence over the default option. Otherwise, the value in `plot_options` is used. Default options are assigned by `set_plot_option`.

Each element of `plot_options` is a list of two or more items. The first item is the name of an option, and the remainder comprises the value or values assigned to the option. In some cases the, the assigned value is a list, which may comprise several items.

The plot options which are recognized by `plot2d` and `plot3d` are the following:

- Option: `plot_format` determines which plotting package is used by `plot2d` and `plot3d`.
  - Default value: `gnuplot` Gnuplot is the default, and most advanced, plotting package. It requires an external gnuplot installation.
  - Value: `mgnuplot` Mgnuplot is a Tk-based wrapper around gnuplot. It is included in the Maxima distribution. Mgnuplot offers a rudimentary GUI for gnuplot, but has fewer overall features than the plain gnuplot interface. Mgnuplot requires an external gnuplot installation and Tcl/Tk.
  - Value: `openmath` Openmath is a Tcl/Tk GUI plotting program. It is included in the Maxima distribution.
  - Value: `ps` Generates simple PostScript files directly from Maxima. Much more sophisticated PostScript output can be generated from gnuplot, by leaving the option `plot_format` unspecified (to accept the default), and setting the option `gnuplot_term` to `ps`.
- Option: `run_viewer` controls whether or not the appropriate viewer for the plot format should be run.

- Default value: `true` Execute the viewer program.
- Value: `false` Do not execute the viewer program.
- `gnuplot_term` Sets the output terminal type for gnuplot.
  - Default value: `default` Gnuplot output is displayed in a separate graphical window.
  - Value: `dumb` Gnuplot output is displayed in the Maxima console by an "ASCII art" approximation to graphics.
  - Value: `ps` Gnuplot generates commands in the PostScript page description language. If the option `gnuplot_out_file` is set to `filename`, gnuplot writes the PostScript commands to `filename`. Otherwise, the commands are printed to the Maxima console.
- Option: `gnuplot_out_file` Write gnuplot output to a file.
  - Default value: `false` No output file specified.
  - Value: `filename` Example: `[gnuplot_out_file, "myplot.ps"]` This example sends PostScript output to the file `myplot.ps` when used in conjunction with the PostScript gnuplot terminal.
- Option: `x` The default horizontal range.
 

`[x, - 3, 3]`

Sets the horizontal range to `[-3, 3]`.
- Option: `y` The default vertical range.
 

`[y, - 3, 3]`

Sets the vertical range to `[-3, 3]`.
- Option: `t` The default range for the parameter in parametric plots.
 

`[t, 0, 10]`

Sets the parametric variable range to `[0, 10]`.
- Option: `nticks` Initial number of points used by the adaptive plotting routine.
 

`[nticks, 20]`

The default for `nticks` is 10.
- Option: `adapt_depth` The maximum number of splittings used by the adaptive plotting routine.
 

`[adapt_depth, 5]`

The default for `adapt_depth` is 10.
- Option: `grid` Sets the number of grid points to use in the x- and y-directions for three-dimensional plotting.
 

`[grid, 50, 50]`

sets the grid to 50 by 50 points. The default grid is 30 by 30.
- Option: `transform_xy` Allows transformations to be applied to three-dimensional plots.
 

`[transform_xy, false]`

The default `transform_xy` is `false`. If it is not `false`, it should be the output of

```
make_transform ([x, y, z], f1(x, y, z), f2(x, y, z), f3(x, y, z))$█
```

The `polar_xy` transformation is built in. It gives the same transformation as

```
make_transform ([r, th, z], r*cos(th), r*sin(th), z)$
```

- Option: `colour_z` is specific to the `ps` plot format.

```
[colour_z, true]
```

The default value for `colour_z` is `false`.

- Option: `view_direction` Specific to the `ps` plot format.

```
[view_direction, 1, 1, 1]
```

The default `view_direction` is `[1, 1, 1]`.

There are several plot options specific to `gnuplot`. All of these options (except `gnuplot_pm3d`) are raw `gnuplot` commands, specified as strings. Refer to the `gnuplot` documentation for more details.

- Option: `gnuplot_pm3d` Controls the usage PM3D mode, which has advanced 3D features. PM3D is only available in `gnuplot` versions after 3.7. The default value for `gnuplot_pm3d` is `false`.

Example:

```
[gnuplot_pm3d, true]
```

- Option: `gnuplot_preamble` Inserts `gnuplot` commands before the plot is drawn. Any valid `gnuplot` commands may be used. Multiple commands should be separated with a semi-colon. The example shown produces a log scale plot. The default value for `gnuplot_preamble` is the empty string `""`.

Example:

```
[gnuplot_preamble, "set log y"]
```

- Option: `gnuplot_curve_titles` Controls the titles given in the plot key. The default value is `default`, which automatically sets the title of each curve to the function plotted. If not `default`, `gnuplot_curve_titles` should contain a list of strings. (To disable the plot key entirely, add `"set nokey"` to `gnuplot_preamble`.)

Example:

```
[gnuplot_curve_titles, ["my first function", "my second function"]]█
```

- Option: `gnuplot_curve_styles` A list of strings controlling the appearance of curves, i.e., color, width, dashing, etc., to be sent to the `gnuplot` plot command. The default value is `["with lines 3", "with lines 1", "with lines 2", "with lines 5", "with lines 4", "with lines 6", "with lines 7"]`, which cycles through different colors. See the `gnuplot` documentation for `plot` for more information.

Example:

```
[gnuplot_curve_styles, ["with lines 7", "with lines 2"]]
```

- Option: `gnuplot_default_term_command` The `gnuplot` command to set the terminal type for the default terminal. The default value is the empty string `""`, i.e., use `gnuplot`'s default.

Example:

```
[gnuplot_default_term_command, "set term x11"]
```

- Option: `gnuplot_dumb_term_command` The gnuplot command to set the terminal type for the dumb terminal. The default value is "set term dumb 79 22", which makes the text output 79 characters by 22 characters.

Example:

```
[gnuplot_dumb_term_command, "set term dumb 132 50"]
```

- Option: `gnuplot_ps_term_command` The gnuplot command to set the terminal type for the PostScript terminal. The default value is "set size 1.5, 1.5;set term postscript eps enhanced color solid 24", which sets the size to 1.5 times gnuplot's default, and the font size to 24, among other things. See the gnuplot documentation for `set term postscript` for more information.

Example:

```
[gnuplot_ps_term_command, "set term postscript eps enhanced color solid 18"]
```

Examples:

- Saves a plot of  $\sin(x)$  to the file `sin.eps`.  

```
plot2d (sin(x), [x, 0, 2*%pi], [gnuplot_term, ps], [gnuplot_out_file, "sin.eps"])
```
- Uses the `y` option to chop off singularities and the `gnuplot_preamble` option to put the key at the bottom of the plot instead of the top.  

```
plot2d ([gamma(x), 1/gamma(x)], [x, -4.5, 5], [y, -10, 10], [gnuplot_preamble,
```
- Uses a very complicated `gnuplot_preamble` to produce fancy x-axis labels. (Note that the `gnuplot_preamble` string must be entered without any line breaks.)  

```
my_preamble: "set xzeroaxis; set xtics ('-2pi' -6.283, '-3pi/2' -4.712, '-pi' -3.14159);
plot2d ([cos(x), sin(x), tan(x), cot(x)], [x, -2*%pi, 2*%pi],
[y, -2, 2], [gnuplot_preamble, my_preamble]);
```
- Uses a very complicated `gnuplot_preamble` to produce fancy x-axis labels, and produces PostScript output that takes advantage of the advanced text formatting available in gnuplot. (Note that the `gnuplot_preamble` string must be entered without any line breaks.)  

```
my_preamble: "set xzeroaxis; set xtics ('-2{/Symbol p}' -6.283, '-3{/Symbol p}' -4.712, '-1{/Symbol p}' -3.14159);
plot2d ([cos(x), sin(x), tan(x)], [x, -2*%pi, 2*%pi], [y, -2, 2],
[gnuplot_preamble, my_preamble], [gnuplot_term, ps], [gnuplot_out_file, "t"])
```
- A three-dimensional plot using the gnuplot `pm3d` terminal.  

```
plot3d (atan (-x^2 + y^3/4), [x, -4, 4], [y, -4, 4], [grid, 50, 50], [gnuplot_term,
```
- A three-dimensional plot without a mesh and with contours projected on the bottom plane.  

```
my_preamble: "set pm3d at s;unset surface;set contour;set cntrparam levels 20;
plot3d (atan (-x^2 + y^3/4), [x, -4, 4], [y, -4, 4], [grid, 50, 50],
[gnuplot_pm3d, true], [gnuplot_preamble, my_preamble])$
```
- A plot where the z-axis is represented by color only. (Note that the `gnuplot_preamble` string must be entered without any line breaks.)  

```
plot3d (cos (-x^2 + y^3/4), [x, -4, 4], [y, -4, 4],
[gnuplot_preamble, "set view map; unset surface"], [gnuplot_pm3d, true], [
```

**plot3d** (*expr*, *x\_range*, *y\_range*, ..., *options*, ...) Function  
**plot3d** (*expr\_1*, *expr\_2*, *expr\_3*, *x\_range*, *y\_range*, ..., *options*, ...) Function  
`plot3d (2^(-u^2+v^2), [u, -5, 5], [v, -7, 7]);`

plots  $z = 2^{-(u^2+v^2)}$  with  $u$  and  $v$  varying in  $[-5,5]$  and  $[-7,7]$  respectively, and with  $u$  on the  $x$  axis, and  $v$  on the  $y$  axis.

An example of the second pattern of arguments is

```
plot3d ([cos(x)*(3 + y*cos(x/2)), sin(x)*(3 + y*cos(x/2)), y*sin(x/2)],
        [x, -%pi, %pi], [y, -1, 1], ['grid, 50, 15]);
```

which plots a Moebius band, parametrized by the three expressions given as the first argument to `plot3d`. An additional optional argument `['grid, 50, 15]` gives the grid number of rectangles in the  $x$  direction and  $y$  direction.

This example shows a plot of the real part of  $z^{1/3}$ .

```
plot3d (r^.33*cos(th/3), [r, 0, 1], [th, 0, 6*%pi],
        ['grid, 12, 80], ['plot_format, ps],
        ['transform_xy, polar_to_xy], ['view_direction, 1, 1, 1.4],
        ['colour_z, true]);
```

Here the `view_direction` option indicates the direction from which we take a projection. We actually do this from infinitely far away, but parallel to the line from `view_direction` to the origin. This is currently only used in `ps` `plot_format`, since the other viewers allow interactive rotating of the object.

Another example is a Klein bottle:

```
expr_1: 5*cos(x)*(cos(x/2)*cos(y) + sin(x/2)*sin(2*y) + 3.0) - 10.0;
expr_2: -5*sin(x)*(cos(x/2)*cos(y) + sin(x/2)*sin(2*y) + 3.0);
expr_3: 5*(-sin(x/2)*cos(y) + cos(x/2)*sin(2*y));
```

```
plot3d ([expr_1, expr_2, expr_3], [x, -%pi, %pi], [y, -%pi, %pi], ['grid, 40,
```

or a torus

```
expr_1: cos(y)*(10.0+6*cos(x));
expr_2: sin(y)*(10.0+6*cos(x));
expr_3: -6*sin(x);
```

```
plot3d ([expr_1, expr_2, expr_3], [x, 0, 2*%pi], [y, 0, 2*%pi], ['grid, 40, 40,
```

We can output to `gnuplot` too:

```
plot3d (2^(x^2 - y^2), [x, -1, 1], [y, -2, 2], [plot_format, gnuplot]);
```

Sometimes you may need to define a function to plot the expression. All the arguments to `plot3d` are evaluated before being passed to `plot3d`, and so trying to make an expression which does just what you want may be difficult, and it is just easier to make a function.

```
M: matrix([1, 2, 3, 4], [1, 2, 3, 2], [1, 2, 3, 4], [1, 2, 3, 3])$
f(x, y) := float (M [?round(x), ?round(y)])$
plot3d (f, [x, 1, 4], [y, 1, 4], ['grid, 4, 4])$
```

See `plot_options` for more examples.



**make\_transform** (*vars, fx, fy, fz*) Function  
 Returns a function suitable for the transform function in plot3d. Use with the plot option `transform_xy`.

```
make_transform ([r, th, z], r*cos(th), r*sin(th), z)$
```

is a transformation to polar coordinates.

**plot2d\_ps** (*expr, range*) Function  
 Writes to `pstream` a sequence of PostScript commands which plot *expr* over *range*. *expr* is an expression. *range* is a list of the form `[x, min, max]` in which *x* is a variable which appears in *expr*.

See also `closeps`.

**closeps** () Function  
 This should usually be called at the end of a sequence of plotting commands. It closes the current output stream `pstream`, and sets it to `nil`. It also may be called at the start of a plot, to ensure `pstream` is closed if it was open. All commands which write to `pstream`, open it if necessary. `closeps` is separate from the other plotting commands, since we may want to plot 2 ranges or superimpose several plots, and so must keep the stream open.

**set\_plot\_option** (*option*) Function  
*option* is of the format of one of the elements of the `plot_options` list. Thus `set_plot_option ([grid, 30, 40])` would change the default grid used by `plot3d`. Note that if the symbol `grid` has a value, then you should quote it here: `set_plot_option (['grid, 30, 40])` so that the value will not be substituted.

**psdraw\_curve** (*ptlist*) Function  
 Draws a curve connecting the points in *ptlist*. The latter may be of the form `[x0, y0, x1, y1, ...]` or `[[x0, y0], [x1, y1], ...]`  
 The function `join` is handy for taking a list of x's and a list of y's and splicing them together.

`psdraw_curve` simply invokes the more primitive function `pscurve`. Here is the definition:

```
(defun $psdraw_curve (lis)
  (p "newpath")
  ($pscurve lis)
  (p "stroke"))
```

**pscom** (*cmd*) Function  
*cmd* is inserted in the PostScript file. Example:

```
pscom ("4.5 72 mul 5.5 72 mul translate 14 14 scale");
```

## 9 Input and Output

### 9.1 Introduction to Input and Output

### 9.2 Files

A file is simply an area on a particular storage device which contains data or text. Files on the disks are figuratively grouped into "directories". A directory is just a list of files. Commands which deal with files are: `save`, `load`, `loadfile`, `stringout`, `batch`, `demo`, `writefile`, `closefile`, and `appendfile`.

### 9.3 Definitions for Input and Output

**%** Variable  
 % is the output expression (e.g., %o1, %o2, %o3, ...) most recently computed by Maxima, whether or not it was displayed. See also %% and %th.

**%%** Variable  
 In a compound statement comprising two or more statements, %% is the value of the previous statement. For example,

```
block (integrate (x^5, x), ev (%%, x=2) - ev (%%, x=1));
block ([prev], prev: integrate (x^5, x), ev (prev, x=2) - ev (prev, x=1));
```

yield the same result, namely 21/2.

A compound statement may comprise other compound statements. Whether a statement be simple or compound, %% is the value of the previous statement. For example,

```
block (block (a^n, %%*42), %%/6)
```

yields  $7 \cdot a^n$ .

Within a compound statement, the value of %% may be inspected at a break prompt, which is opened by executing the `break` function. For example, at the break prompt opened by

```
block (a: 42, break ())$
```

entering %%; yields 42.

At the first statement in a compound statement, or outside of a compound statement, %% is undefined.

**%edispflag** Variable  
 Default value: `false`

When `%edispflag` is `true`, Maxima displays %e to a negative exponent as a quotient. For example,  $e^{-x}$  is displayed as  $1/e^x$ .

- %th** (*i*) Function  
 The value of the *i*'th previous output expression. That is, if the next expression to be computed is the *n*'th output, %th (*m*) is the (*n* - *m*)'th output.  
 %th is useful in `batch` files or for referring to a group of output expressions. For example,  

```
block (s: 0, for i:1 thru 10 do s: s + %th (i))$
```

 sets `s` to the sum of the last ten output expressions.
- "?"** special symbol  
 As prefix to a function or variable name, ? signifies that the name is a Lisp name, not a Maxima name. For example, ?round is a Lisp function.  
 The notation ? word (a question mark followed a word, separated by whitespace) is equivalent to `describe ("word")`.
- absboxchar** Variable  
 Default value: !  
 absboxchar is the character used to draw absolute value signs around expressions which are more than one line tall.
- appendfile** (*filename*) Function  
 Appends a console transcript to *filename*. appendfile is the same as writefile, except that the transcript file, if it exists, is always appended.  
 closefile closes the transcript file opened by appendfile or writefile.
- batch** (*filename*) Function  
 Reads Maxima expressions from *filename* and evaluates them. batch searches for *filename* in the list `file_search_maxima`. See `file_search`.  
*filename* comprises a sequence of Maxima expressions, each terminated with ; or \$.  
 The special variable % and the function %th refer to previous results within the file.  
 The file may include :lisp constructs. Spaces, tabs, and newlines in the file are ignored. A suitable input file may be created by a text editor or by the `stringout` function.  
 batch reads each input expression from *filename*, displays the input to the console, computes the corresponding output expression, and displays the output expression. Input labels are assigned to the input expressions and output labels are assigned to the output expressions. batch evaluates every input expression in the file unless there is an error. If user input is requested (by `asksign` or `askinteger`, for example) batch pauses to collect the requisite input and then continue.  
 It may be possible to halt batch by typing control-C at the console. The effect of control-C depends on the underlying Lisp implementation.  
 batch has several uses, such as to provide a reservoir for working command lines, to give error-free demonstrations, or to help organize one's thinking in solving complex problems.  
 batch evaluates its argument. batch has no return value.  
 See also `load`, `batchload`, and `demo`.

**batchload** (*filename*) Function

Reads Maxima expressions from *filename* and evaluates them, without displaying the input or output expressions and without assigning labels to output expressions. Printed output (such as produced by `print` or `describe`) is displayed, however.

The special variable `%` and the function `%th` refer to previous results from the interactive interpreter, not results within the file. The file cannot include `:lisp` constructs. `batchload` returns the path of *filename*, as a string. `batchload` evaluates its argument.

See also `batch` and `load`.

**closefile** () Function

Closes the transcript file opened by `writetfile` or `appendfile`.

**collapse** (*expr*) Function

Collapses *expr* by causing all of its common (i.e., equal) subexpressions to share (i.e., use the same cells), thereby saving space. (`collapse` is a subroutine used by the `optimize` command.) Thus, calling `collapse` may be useful after loading in a `save` file. You can collapse several expressions together by using `collapse ([expr_1, ..., expr_n])`. Similarly, you can collapse the elements of the array `A` by doing `collapse (listarray ('A))`.

**concat** (*arg\_1*, *arg\_2*, ...) Function

Concatenates its arguments. The arguments must evaluate to atoms. The return value is a symbol if the first argument is a symbol and a Maxima string otherwise.

`concat` evaluates its arguments. The single quote `'` prevents evaluation.

```
(%i1) y: 7$
(%i2) z: 88$
(%i3) concat (y, z/2);
(%o3)                                     744
(%i4) concat ('y, z/2);
(%o4)                                     y44
```

A symbol constructed by `concat` may be assigned a value and appear in expressions.

The `::` (double colon) assignment operator evaluates its left-hand side.

```
(%i5) a: concat ('y, z/2);
(%o5)                                     y44
(%i6) a:: 123;
(%o6)                                     123
(%i7) y44;
(%o7)                                     123
(%i8) b^a;
(%o8)                                     y44
(%o8)                                     b
(%i9) %, numer;
(%o9)                                     123
(%o9)                                     b
```

Note that although `concat (1, 2)` looks like a number, it is a Maxima string.

```
(%i10) concat (1, 2) + 3;
(%o10)          12 + 3
```

**sconcat** (*arg\_1*, *arg\_2*, ...) Function

Concatenates its arguments into a string. Unlike `concat`, the arguments do *not* need to be atoms.

The result is a Lisp string.

```
(%i1) sconcat ("xx[" , 3, "]" , expand ((x+y)^3));
(%o1)          xx[3]:y^3+3*x*y^2+3*x^2*y+x^3
```

**disp** (*expr\_1*, *expr\_2*, ...) Function

is like `display` but only the value of the arguments are displayed rather than equations. This is useful for complicated arguments which don't have names or where only the value of the argument is of interest and not the name.

**dispcn** (*tensor\_1*, *tensor\_2*, ...) Function

**dispcn** (*all*) Function

Displays the contraction properties of its arguments as were given to `defcon`. `dispcn` (*all*) displays all the contraction properties which were defined.

**display** (*expr\_1*, *expr\_2*, ...) Function

Displays equations whose left side is *expr\_i* unevaluated, and whose right side is the value of the expression centered on the line. This function is useful in blocks and `for` statements in order to have intermediate results displayed. The arguments to `display` are usually atoms, subscripted variables, or function calls. See also `disp`.

```
(%i1) display(B[1,2]);
              2
          B    = X - X
          1, 2
(%o1)          done
```

**display2d** Variable

Default value: `true`

When `display2d` is `false`, the console display is a string (1-dimensional) form rather than a display (2-dimensional) form.

**display\_format\_internal** Variable

Default value: `false`

When `display_format_internal` is `true`, expressions are displayed without being transformed in ways that hide the internal mathematical representation. The display then corresponds to what `inpart` returns rather than `part`.

Examples:

User	<code>part</code>	<code>inpart</code>
<code>a-b;</code>	<code>A - B</code>	<code>A + (- 1) B</code>

```

a/b;      A      - 1
          -      A B
          B

sqrt(x);  sqrt(X)  X
          1/2

X*4/3;    4 X      4
          ---      - X
          3        3

```

**dispterms** (*expr*)

Function

Displays *expr* in parts one below the other. That is, first the operator of *expr* is displayed, then each term in a sum, or factor in a product, or part of a more general expression is displayed separately. This is useful if *expr* is too large to be otherwise displayed. For example if P1, P2, ... are very large expressions then the display program may run out of storage space in trying to display P1 + P2 + ... all at once. However, `dispterms (P1 + P2 + ...)` displays P1, then below it P2, etc. When not using `dispterms`, if an exponential expression is too wide to be displayed as A^B it appears as `expt (A, B)` (or as `ncexpt (A, B)` in the case of A^B).

**error\_size**

Variable

Default value: 10

`error_size` modifies error messages according to the size of expressions which appear in them. If the size of an expression (as determined by the Lisp function `ERROR-SIZE`) is greater than `error_size`, the expression is replaced in the message by a symbol, and the symbol is assigned the expression. The symbols are taken from the list `error_syms`.

Otherwise, the expression is smaller than `error_size`, and the expression is displayed in the message.

See also `error` and `error_syms`.

Example:

The size of U, as determined by `ERROR-SIZE`, is 24.

```
(%i1) U: (C^D^E + B + A)/(cos(X-1) + 1)$
```

```
(%i2) error_size: 20$
```

```
(%i3) error ("Example expression is", U);
```

```
Example expression is errexp1
```

```
-- an error. Quitting. To debug this try debugmode(true);
```

```
(%i4) errexp1;
```

```
(%o4)
          E
          D
          C  + B + A
          -----
          cos(X - 1) + 1
```

```
(%i5) error_size: 30$

(%i6) error ("Example expression is", U);

      E
      D
      C  + B + A
Example expression is -----
                        cos(X - 1) + 1
-- an error.  Quitting.  To debug this try debugmode(true);
```

**error\_syms**

Variable

Default value: [errex1, errex2, errex3]

In error messages, expressions larger than `error_size` are replaced by symbols, and the symbols are set to the expressions. The symbols are taken from the list `error_syms`. The first too-large expression is replaced by `error_syms[1]`, the second by `error_syms[2]`, and so on.

If there are more too-large expressions than there are elements of `error_syms`, symbols are constructed automatically, with the  $n$ -th symbol equivalent to `concat('errex', n)`.

See also `error` and `error_size`.

**expt** (*a*, *b*)

Function

If an exponential expression is too wide to be displayed as  $a^b$  it appears as `expt (a, b)` (or as `ncxpt (a, b)` in the case of  $a^{b^c}$ ).

`expt` and `ncxpt` are not recognized in input.

**exptdispflag**

Variable

Default value: true

When `exptdispflag` is true, Maxima displays expressions with negative exponents using quotients, e.g.,  $X^{-1}$  as  $1/X$ .

**filename\_merge** (*path*, *filename*)

Function

Constructs a modified path from *path* and *filename*. If the final component of *path* is of the form `###.something`, the component is replaced with `filename.something`.

Otherwise, the final component is simply replaced by *filename*.

**file\_search** (*filename*)

Function

**file\_search** (*filename*, *pathlist*)

Function

`file_search` searches for the file *filename* and returns the path to the file (as a string) if it can be found; otherwise `file_search` returns false. `file_search (filename)` searches in the default search directories, which are specified by the `file_search_maxima`, `file_search_lisp`, and `file_search_demo` variables.

`file_search` first checks if the actual name passed exists, before attempting to match it to "wildcard" file search patterns. See `file_search_maxima` concerning file search patterns.

The argument *filename* can be a path and file name, or just a file name, or, if a file search directory includes a file search pattern, just the base of the file name (without an extension). For example,

```
file_search ("/home/wfs/special/zeta.mac");
file_search ("zeta.mac");
file_search ("zeta");
```

all find the same file, assuming the file exists and `/home/wfs/special/###.mac` is in `file_search_maxima`.

`file_search (filename, pathlist)` searches only in the directories specified by *pathlist*, which is a list of strings. The argument *pathlist* supersedes the default search directories, so if the path list is given, `file_search` searches only the ones specified, and not any of the default search directories. Even if there is only one directory in *pathlist*, it must still be given as a one-element list.

The user may modify the default search directories. See `file_search_maxima`.

`file_search` is invoked by `load` with `file_search_maxima` and `file_search_lisp` as the search directories.

<b>file_search_maxima</b>	Variable
<b>file_search_lisp</b>	Variable
<b>file_search_demo</b>	Variable

These variables specify lists of directories to be searched by `load`, `demo`, and some other Maxima functions. The default values of these variables name various directories in the Maxima installation.

The user can modify these variables, either to replace the default values or to append additional directories. For example,

```
file_search_maxima: ["/usr/local/foo/###.mac",
"/usr/local/bar/###.mac"]$
```

replaces the default value of `file_search_maxima`, while

```
file_search_maxima: append (file_search_maxima,
["/usr/local/foo/###.mac", "/usr/local/bar/###.mac"])$
```

appends two additional directories. It may be convenient to put such an expression in the file `maxima-init.mac` so that the file search path is assigned automatically when Maxima starts.

Multiple filename extensions and multiple paths can be specified by special “wildcard” constructions. The string `###` expands into the sought-after name, while a comma-separated list enclosed in curly braces `{foo,bar,baz}` expands into multiple strings. For example, supposing the sought-after name is `neumann`,

```
"/home/{wfs,gcj}/###.{lisp,mac}"
```

expands into `/home/wfs/neumann.lisp`, `/home/wfs/neumann.mac`, `/home/gcj/neumann.lisp`, and `/home/gcj/neumann.mac`.

<b>file_type (filename)</b>	Function
-----------------------------	----------

Returns a guess about the content of *filename*, based on the filename extension. *filename* need not refer to an actual file; no attempt is made to open the file and inspect the content.



The return value is a symbol, either `object`, `lisp`, or `maxima`. If the extension starts with `m` or `d`, `file_type` returns `maxima`. If the extension starts with `l`, `file_type` returns `lisp`. If none of the above, `file_type` returns `object`.

**grind** (*expr*)  
**grind**

Function  
 Variable

The function `grind` prints *expr* to the console in a form suitable for input to Maxima. `grind` always returns `done`.

See also `string`, which returns a string instead of printing its output. `grind` attempts to print the expression in a manner which makes it slightly easier to read than the output of `string`.

When the variable `grind` is `true`, the output of `string` and `stringout` has the same format as that of `grind`; otherwise no attempt is made to specially format the output of those functions. The default value of the variable `grind` is `false`.

`grind` can also be specified as an argument of `playback`. When `grind` is present, `playback` prints input expressions in the same format as the `grind` function. Otherwise, no attempt is made to specially format input expressions.

**ibase**

Variable

Default value: 10

Integers entered into Maxima are interpreted with respect to the base `ibase`.

`ibase` may be assigned any integer between 2 and 35 (decimal), inclusive. When `ibase` is greater than 10, the numerals comprise the decimal numerals 0 through 9 plus capital letters of the alphabet A, B, C, ..., as needed. The numerals for base 35, the largest acceptable base, comprise 0 through 9 and A through Y.

See also `obase`.

**inchar**

Variable

Default value: %i

`inchar` is the prefix of the labels of expressions entered by the user. Maxima automatically constructs a label for each input expression by concatenating `inchar` and `linenum`. `inchar` may be assigned any string or symbol, not necessarily a single character.

```
(%i1) inchar: "input";
(%o1)                                     input
(input1) expand ((a+b)^3);
(%o1)                                     3      2      2      3
      b + 3 a b + 3 a b + a
(input2)
```

See also `labels`.

**ldisp** (*expr\_1*, ..., *expr\_n*)

Function

Displays expressions *expr\_1*, ..., *expr\_n* to the console as printed output. `ldisp` assigns an intermediate expression label to each argument and returns the list of labels.

See also `disp`.

```
(%i1) e: (a+b)^3;
(%o1) (b + a)3
(%i2) f: expand (e);
(%o2) b3 + 3 a b2 + 3 a2 b + a3
(%i3) ldisp (e, f);
(%t3) (b + a)3
(%t4) b3 + 3 a b2 + 3 a2 b + a3
(%o4) [%t3, %t4]
(%i4) %t3;
(%o4) (b + a)3
(%i5) %t4;
(%o5) b3 + 3 a b2 + 3 a2 b + a3
```

**ldisplay** (*expr\_1*, ..., *expr\_n*)

Function

Displays expressions *expr\_1*, ..., *expr\_n* to the console as printed output. Each expression is printed as an equation of the form **lhs = rhs** in which **lhs** is one of the arguments of `ldisplay` and **rhs** is its value. Typically each argument is a variable. `ldisp` assigns an intermediate expression label to each equation and returns the list of labels.

See also `display`.

```
(%i1) e: (a+b)^3;
(%o1) (b + a)3
(%i2) f: expand (e);
(%o2) b3 + 3 a b2 + 3 a2 b + a3
(%i3) ldisplay (e, f);
(%t3) e = (b + a)3
(%t4) f = b3 + 3 a b2 + 3 a2 b + a3
(%o4) [%t3, %t4]
(%i4) %t3;
(%o4) e = (b + a)3
(%i5) %t4;
(%o5) f = b3 + 3 a b2 + 3 a2 b + a3
```

**linechar**

Variable

Default value: %t

**linechar** is the prefix of the labels of intermediate expressions generated by Maxima. Maxima constructs a label for each intermediate expression (if displayed) by concatenating **linechar** and **linenum**. **linechar** may be assigned any string or symbol, not necessarily a single character.

Intermediate expressions might or might not be displayed. See **programmode** and **labels**.

**linel**

Variable

Default value: 79

**linel** is the assumed width (in characters) of the console display for the purpose of displaying expressions. **linel** may be assigned any value by the user, although very small or very large values may be impractical. Text printed by built-in Maxima functions, such as error messages and the output of **describe**, is not affected by **linel**.

**load** (*filename*)

Function

Evaluates expressions in *filename*, thus bringing variables, functions, and other objects into Maxima. The binding of any existing object is clobbered by the binding recovered from *filename*. To find the file, **load** calls **file\_search** with **file\_search\_maxima** and **file\_search\_lisp** as the search directories. If **load** succeeds, it returns the name of the file. Otherwise **load** prints an error message.

**load** works equally well for Lisp code and Maxima code. Files created by **save**, **translate\_file**, and **compile\_file**, which create Lisp code, and **stringout**, which creates Maxima code, can all be processed by **load**. **load** calls **loadfile** to load Lisp files and **batchload** to load Maxima files.

See also **loadfile**, **batch**, **batchload**, and **demo**. **loadfile** processes Lisp files; **batch**, **batchload**, and **demo** process Maxima files.

See **file\_search** for more detail about the file search mechanism.

**load** evaluates its argument.

**loadfile** (*filename*)

Function

Evaluates Lisp expressions in *filename*. **loadfile** does not invoke **file\_search**, so *filename* must include the file extension and as much of the path as needed to find the file.

**loadfile** can process files created by **save**, **translate\_file**, and **compile\_file**. The user may find it more convenient to use **load** instead of **loadfile**.

**loadfile** quotes its argument, so *filename* must be a literal string, not a string variable. The double-single-quote operator defeats quotation.

**loadprint**

Variable

Default value: **true**

**loadprint** tells whether to print a message when a file is loaded.

- When `loadprint` is `true`, always print a message.
- When `loadprint` is `'loadfile`, print a message only if a file is loaded by the function `loadfile`.
- When `loadprint` is `'autoload`, print a message only if a file is automatically loaded. See `setup_autoload`.
- When `loadprint` is `false`, never print a message.

**obase**

Variable

Default value: 10

`obase` is the base for integers displayed by Maxima.

`obase` may be assigned any integer between 2 and 35 (decimal), inclusive. When `obase` is greater than 10, the numerals comprise the decimal numerals 0 through 9 plus capital letters of the alphabet A, B, C, ..., as needed. The numerals for base 35, the largest acceptable base, comprise 0 through 9, and A through Y.

See also `ibase`.**outchar**

Variable

Default value: `%o`

`outchar` is the prefix of the labels of expressions computed by Maxima. Maxima automatically constructs a label for each computed expression by concatenating `outchar` and `linenum`. `outchar` may be assigned any string or symbol, not necessarily a single character.

```
(%i1) outchar: "output";
(output1)                                     output
(%i2) expand ((a+b)^3);
(output2)          3      2      2      3
                  b  + 3 a b  + 3 a  b  + a
(%i3)
```

See also `labels`.**packagefile**

Variable

Default value: `false`

Package designers who use `save` or `translate` to create packages (files) for others to use may want to set `packagefile: true` to prevent information from being added to Maxima's information-lists (e.g. `values`, `functions`) except where necessary when the file is loaded in. In this way, the contents of the package will not get in the user's way when he adds his own data. Note that this will not solve the problem of possible name conflicts. Also note that the flag simply affects what is output to the package file. Setting the flag to `true` is also useful for creating Maxima init files.

**pformat**

Variable

Default value: `false`

When `pformat` is `true`, a ratio of integers is displayed with the solidus (forward slash) character, and an integer denominator `n` is displayed as a leading multiplicative term  $1/n$ .

```
(%i1) pformat: false$
(%i2) 2^16/7^3;
(%o2)
          65536
        -----
          343
(%i3) (a+b)/8;
(%o3)
          b + a
        -----
          8
(%i4) pformat: true$
(%i5) 2^16/7^3;
(%o5)
          65536/343
(%i6) (a+b)/8;
(%o6)
          1/8 (b + a)
```

**print** (*expr\_1*, ..., *expr\_n*) Function

Evaluates and displays *expr\_1*, ..., *expr\_n* one after another, from left to right, starting at the left edge of the console display.

The value returned by **print** is the value of its last argument. **print** does not generate intermediate expression labels.

See also **display**, **disp**, **ldisplay**, and **ldisp**. Those functions display one expression per line, while **print** attempts to display two or more expressions per line.

To display the contents of a file, see **printfile**.

```
(%i1) r: print ("(a+b)^3 is", expand ((a+b)^3), "log (a^10/b) is", radcan (log
          3      2      2      3
(a+b)^3 is b + 3 a b + 3 a b + a log (a^10/b) is
          10 log(a) - log(b)
(%i2) r;
(%o2)
          10 log(a) - log(b)
(%i3) disp ("(a+b)^3 is", expand ((a+b)^3), "log (a^10/b) is", radcan (log (a^
          (a+b)^3 is
          3      2      2      3
          b + 3 a b + 3 a b + a
          log (a^10/b) is
          10 log(a) - log(b)
```

**tcl\_output** (*list*, *i0*, *skip*) Function

**tcl\_output** (*list*, *i0*) Function

**tcl\_output** (*[list\_1*, ..., *list\_n]*, *i*) Function

Prints elements of a list enclosed by curly braces { }, suitable as part of a program in the Tcl/Tk language.

**tcl\_output** (*list*, *i0*, *skip*) prints *list*, beginning with element *i0* and printing elements *i0* + *skip*, *i0* + 2 *skip*, etc.

`tcl_output (list, i0)` is equivalent to `tcl_output (list, i0, 2)`.

`tcl_output ([list_1, ..., list_n], i)` prints the  $i$ 'th elements of `list_1, ..., list_n`.

Examples:

```
(%i1) tcl_output ([1, 2, 3, 4, 5, 6], 1, 3)$
{1.000000000    4.000000000
}
(%i2) tcl_output ([1, 2, 3, 4, 5, 6], 2, 3)$
{2.000000000    5.000000000
}
(%i3) tcl_output ([3/7, 5/9, 11/13, 13/17], 1)$
{((RAT SIMP) 3 7) ((RAT SIMP) 11 13)
}
(%i4) tcl_output ([x1, y1, x2, y2, x3, y3], 2)$
{$Y1 $Y2 $Y3
}
(%i5) tcl_output ([[1, 2, 3], [11, 22, 33]], 1)$
{SIMP 1.000000000    11.000000000
}
```

**read** (*expr\_1, ..., expr\_n*)

Function

Prints *expr\_1, ..., expr\_n*, then reads one expression from the console and returns the evaluated expression. The expression is terminated with a semicolon ; or dollar sign \$.

See also `readonly`.

```
(%i1) foo: 42$
(%i2) foo: read ("foo is", foo, " -- enter new value.")$
foo is 42 -- enter new value.
(a+b)^3;
(%i3) foo;
(%o3) (b + a)3
```

**readonly** (*expr\_1, ..., expr\_n*)

Function

Prints *expr\_1, ..., expr\_n*, then reads one expression from the console and returns the expression (without evaluation). The expression is terminated with a ; (semicolon) or \$ (dollar sign).

```
(%i1) aa: 7$
(%i2) foo: readonly ("Enter an expression:");
Enter an expression:
2^aa;
(%o2) aa2
```

```
(%i3) foo: read ("Enter an expression:");
Enter an expression:
2^aa;
(%o3)                                     128
```

See also read.

**reveal** (*expr*, *depth*) Function

Replaces parts of *expr* at the specified integer *depth* with descriptive summaries.

- Sums and differences are replaced by `sum(n)` where *n* is the number of operands of the sum.
- Products are replaced by `product(n)` where *n* is the number of operands of the sum.
- Exponentials are replaced by `expt`.
- Quotients are replaced by `quotient`.
- Unary negation is replaced by `negterm`.

When *depth* is greater than or equal to the maximum depth of *expr*, `reveal (expr, depth)` returns *expr* unmodified.

`reveal` evaluates its arguments. `reveal` returns the summarized expression.

Example:

```
(%i1) e: expand ((a - b)^2)/expand ((exp(a) + exp(b))^2);
                                     2          2
                                     b  - 2 a b + a
(%o1) -----
          b + a      2 b      2 a
2 %e      + %e      + %e
(%i2) reveal (e, 1);
(%o2) quotient
(%i3) reveal (e, 2);
          sum(3)
          -----
          sum(3)
(%i4) reveal (e, 3);
          expt + negterm + expt
(%o4) -----
          product(2) + expt + expt
(%i5) reveal (e, 4);
          2          2
          b  - product(3) + a
(%o5) -----
          product(2)      product(2)
2 expt + %e      + %e
(%i6) reveal (e, 5);
          2          2
          b  - 2 a b + a
(%o6) -----
          sum(2)      2 b      2 a
```

```

                2 %e      + %e      + %e
(%i7) reveal (e, 6);
                2          2
                b  - 2 a b + a
(%o7)  -----
                b + a      2 b      2 a
2 %e      + %e      + %e

```

**rmxchar**

Variable

Default value: ]

**rmxchar** is the character drawn on the right-hand side of a matrix.

See also **lmxchar**.

**save** (*filename*, *name\_1*, *name\_2*, *name\_3*, ...)

Function

**save** (*filename*, *values*, *functions*, *labels*, ...)

Function

**save** (*filename*, [*m*, *n*])

Function

**save** (*filename*, *name\_1=expr\_1*, ...)

Function

**save** (*filename*, *all*)

Function

Stores the current values of *name\_1*, *name\_2*, *name\_3*, ..., in *filename*. The arguments must be names of variables, functions, or other objects. **save** returns *filename*.

**save** stores data in the form of Lisp expressions. The data stored by **save** may be recovered by **load** (*filename*). The effect of executing **save** when *filename* already exists depends on the underlying Lisp implementation; the file may be clobbered, or **save** may complain with an error message.

The special form **save** (*filename*, *values*, *functions*, *labels*, ...) stores the items named by *values*, *functions*, *labels*, etc. The names may be any specified by the variable *infolists*. *values* comprises all user-defined variables.

The special form **save** (*filename*, [*m*, *n*]) stores the values of input and output labels *m* through *n*. Note that *m* and *n* must be literal integers or double-quoted symbols. Input and output labels may also be stored one by one, e.g., **save** ("foo.1", %i42, %o42). **save** (*filename*, *labels*) stores all input and output labels. When the stored labels are recovered, they clobber existing labels.

The special form **save** (*filename*, *name\_1=expr\_1*, *name\_2=expr\_2*, ...) stores the values of *expr\_1*, *expr\_2*, ..., with names *name\_1*, *name\_2*, .... It is useful to apply this form to input and output labels, e.g., **save** ("foo.1", aa=%o88). The right-hand side of the equality in this form may be any expression, which is evaluated. This form does not introduce the new names into the current Maxima environment, but only stores them in *filename*.

These special forms and the general form of **save** may be mixed at will. For example, **save** (*filename*, aa, bb, cc=42, functions, [11, 17]).

The special form **save** (*filename*, *all*) stores the current state of Maxima. This includes all user-defined variables, functions, arrays, etc., as well as some automatically defined items. The saved items include system variables, such as *file\_search\_maxima* or *showtime*, if they have been assigned new values by the user; see *myoptions*.



**save** quotes its arguments. *filename* must be a string, not a string variable. The first and last labels to save, if specified, must be integers. The double quote operator evaluates a string variable to its string value, e.g., `s: "foo.1"$ save ('s, all)$`, and integer variables to their integer values, e.g., `m: 5$ n: 12$ save ("foo.1", ['m, 'n])$`.

**savedef** Variable

Default value: `true`

When **savedef** is `true`, the Maxima version of a user function is preserved when the function is translated. This permits the definition to be displayed by `dispfun` and allows the function to be edited.

When **savedef** is `false`, the names of translated functions are removed from the `functions` list.

**show** (*expr*) Function

Displays *expr* with the indexed objects in it shown having covariant indices as subscripts, contravariant indices as superscripts. The derivative indices are displayed as subscripts, separated from the covariant indices by a comma.

**showratvars** (*expr*) Function

Returns a list of the canonical rational expression (CRE) variables in expression *expr*.

See also `ratvars`.

**stardisp** Variable

Default value: `false`

When **stardisp** is `true`, multiplication is displayed with an asterisk `*` between operands.

**string** (*expr*) Function

Converts *expr* to Maxima's linear notation just as if it had been typed in.

The return value of **string** is a string, and thus it cannot be used in a computation.

**stringout** (*filename, expr\_1, expr\_2, expr\_3, ...*) Function

**stringout** (*filename, [m, n]*) Function

**stringout** (*filename, input*) Function

**stringout** (*filename, functions*) Function

**stringout** (*filename, values*) Function

**stringout** writes expressions to a file in the same form the expressions would be typed for input. The file can then be used as input for the `batch` or `demo` commands, and it may be edited for any purpose. **stringout** can be executed while `writefile` is in progress.

The general form of **stringout** writes the values of one or more expressions to the output file. Note that if an expression is a variable, only the value of the variable is written and not the name of the variable. As a useful special case, the expressions may be input labels (`%i1, %i2, %i3, ...`) or output labels (`%o1, %o2, %o3, ...`).

If `grind` is `true`, `stringout` formats the output using the `grind` format. Otherwise the `string` format is used. See `grind` and `string`.

The special form `stringout (filename, [m, n])` writes the values of input labels `m` through `n`, inclusive.

The special form `stringout (filename, input)` writes all input labels to the file.

The special form `stringout (filename, functions)` writes all user-defined functions (named by the global list `functions`) to the file.

The special form `stringout (filename, values)` writes all user-assigned variables (named by the global list `values`) to the file. Each variable is printed as an assignment statement, with the name of the variable, a colon, and its value. Note that the general form of `stringout` does not print variables as assignment statements.

<b>tex</b> ( <i>expr</i> )	Function
<b>tex</b> ( <i>expr</i> , <i>filename</i> )	Function
<b>tex</b> ( <i>label</i> , <i>filename</i> )	Function

In the case of a label, a left-equation-number is produced. In case a file-name is supplied, the output is appended to the file.

```
(%i1) integrate (1/(1+x^3), x);
              2      2      2      2      2      2      2      2      2      2
              log(x  - x + 1)  atan(-----)  log(x + 1)
              6              sqrt(3)  3
(%o1)  - ----- + ----- + -----
              6              sqrt(3)  3

(%i2) tex (%o1);
$$$$-{\log \left(x^2-x+1\right)}\over{6}}+{\{\arctan \left({2\,x-1}\right)}\over{\sqrt{3}}}\over{\sqrt{3}}+{\log \left(x+1\right)}\over{3}}\leqno{\tt (\%o1)}$$$$
(%o2)  (\%o1)
(%i3) tex (integrate (sin(x), x));
$$$$-\cos x$$$$
(%o3)  false
(%i4) tex (%o1, "foo.tex");
(%o4)  (\%o1)
```

<b>system</b> ( <i>command</i> )	Function
----------------------------------	----------

Executes *command* as a separate process. The command is passed to the default shell for execution. `system` is not supported by all operating systems, but generally exists in Unix and Unix-like environments.

Supposing `_hist.out` is a list of frequencies which you wish to plot as a bar graph using `xgraph`.

```
(%i1) (with_stdout("_hist.out",
      for i:1 thru length(hist) do (
        print(i,hist[i]))),
      system("xgraph -bar -brw .7 -nl < _hist.out"));
```

In order to make the plot be done in the background (returning control to Maxima) and remove the temporary file after it is done do:

```
system("(xgraph -bar -brw .7 -nl < _hist.out; rm -f _hist.out)&")
```

**ttyoff**

Variable

Default value: `false`

When `ttyoff` is `true`, output expressions are not displayed. Output expressions are still computed and assigned labels. See `labels`.

Text printed by built-in Maxima functions, such as error messages and the output of `describe`, is not affected by `ttyoff`.

**with\_stdout** (*filename*, *expr\_1*, *expr\_2*, *expr\_3*, ...)

Macro

Opens *filename* and then evaluates *expr\_1*, *expr\_2*, *expr\_3*, .... The values of the arguments are not stored in *filename*, but any printed output generated by evaluating the arguments (from `print`, `display`, `disp`, or `grind`, for example) goes to *filename* instead of the console.

`with_stdout` returns the value of its final argument.

See also `writefile`.

```
(%i1) with_stdout ("tmp.out", for i:5 thru 10 do print (i, "! yields", i!))$
(%i2) printfile ("tmp.out")$
5 ! yields 120
6 ! yields 720
7 ! yields 5040
8 ! yields 40320
9 ! yields 362880
10 ! yields 3628800
```

**writefile** (*filename*)

Function

Begins writing a transcript of the Maxima session to *filename*. All interaction between the user and Maxima is then recorded in this file, just as it appears on the console.

As the transcript is printed in the console output format, it cannot be reloaded into Maxima. To make a file containing expressions which can be reloaded, see `save` and `stringout`. `save` stores expressions in Lisp form, while `stringout` stores expressions in Maxima form.

The effect of executing `writefile` when *filename* already exists depends on the underlying Lisp implementation; the transcript file may be clobbered, or the file may be appended. `appendfile` always appends to the transcript file.

It may be convenient to execute `playback` after `writefile` to save the display of previous interactions. As `playback` displays only the input and output variables (`%i1`, `%o1`, etc.), any output generated by a print statement in a function (as opposed to a return value) is not displayed by `playback`.

`closefile` closes the transcript file opened by `writefile` or `appendfile`.

## 10 Floating Point

### 10.1 Definitions for Floating Point

**bfac** (*expr*, *n*) Function  
 Bigfloat version of the factorial (shifted gamma) function. The second argument is how many digits to retain and return, it's a good idea to request a couple of extra.  
`load ("bfac")` loads this function.

**algepsilon** Variable  
 Default value:  $10^{-8}$   
`algepsilon` is used by `algsys`.

**bfloat** (*expr*) Function  
 Converts all numbers and functions of numbers to bigfloat numbers. Setting `fpprec` to *n*, sets the bigfloat precision to *n* digits.  
 When `float2bf` is `false` a warning message is printed when a floating point number is converted into a bigfloat number (since this may lead to loss of precision).

**bfloatp** (*expr*) Function  
 Returns `true` if *expr* is a bigfloat number, otherwise `false`.

**bfpsi** (*n*, *z*, *fpprec*) Function  
**bfpsi0** (*z*, *fpprec*) Function  
`bfpsi` is the polygamma function of real argument *z* and integer order *n*. `bfpsi0` is the digamma function. `bfpsi0` (*z*, *fpprec*) is equivalent to `bfpsi` (0, *z*, *fpprec*).  
 These functions return bigfloat values. *fpprec* is the bigfloat precision of the return value.  
`load ("bfac")` loads these functions.

**bftorat** Variable  
 Default value: `false`  
`bftorat` controls the conversion of bfloats to rational numbers. When `bftorat` is `false`, `ratepsilon` will be used to control the conversion (this results in relatively small rational numbers). When `bftorat` is `true`, the rational number generated will accurately represent the bfloat.

**bftrunc** Variable  
 Default value: `true`  
`bftrunc` causes trailing zeroes in non-zero bigfloat numbers not to be displayed. Thus, if `bftrunc` is `false`, `bfloat` (1) displays as `1.000000000000000B0`. Otherwise, this is displayed as `1.0B0`.

- cbffac** (*z*, *fpprec*) Function  
 Complex bigfloat factorial.  
`load ("bfffac")` loads this function.
- float** (*expr*) Function  
 Converts integers, rational numbers and bigfloats in *expr* to floating point numbers.  
 It is also an **evflag**, **float** causes non-integral rational numbers and bigfloat numbers to be converted to floating point.
- float2bf** Variable  
 Default value: **false**  
 When **float2bf** is **false**, a warning message is printed when a floating point number is converted into a bigfloat number (since this may lead to loss of precision).
- floatnump** (*expr*) Function  
 Returns **true** if *expr* is a floating point number, otherwise **false**.
- fpprec** Variable  
 Default value: 16  
**fpprec** is the Maxima floating point precision. **fpprec** can be set to an integer representing the desired precision.
- fpprintprec** Variable  
 Default value: 0  
**fpprintprec** is the number of digits to print when printing a bigfloat number, making it possible to compute with a large number of digits of precision, but have the answer printed out with a smaller number of digits.  
 When **fpprintprec** is 0, or greater than or equal to **fpprec**, then the value of **fpprec** controls the number of digits used for printing.  
 When **fpprintprec** has a value between 2 and **fpprec** - 1, then it controls the number of digits used. (The minimal number of digits used is 2, one to the left of the point and one to the right.  
 The value 1 for **fpprintprec** is illegal.
- ?round** (*x*) Lisp function  
**?round** (*x*, *divisor*) Lisp function  
 Round the floating point *x* to the nearest integer. The argument must be an ordinary float, not a bigfloat. The ? beginning the name indicates this is a Lisp function.  

```
(%i1) ?round (-2.8);
(%o1) -3
```
- ?truncate** (*x*) Lisp function  
**?truncate** (*x*, *divisor*) Lisp function  
 Truncate the floating point *x* towards 0, to become an integer. The argument must be an ordinary float, not a bigfloat. The ? beginning the name indicates this is a Lisp function.

```
(%i1) ?truncate (-2.8);  
(%o1) - 2  
(%i2) ?truncate (2.4);  
(%o2) 2  
(%i3) ?truncate (2.8);  
(%o3) 2
```



# 11 Contexts

## 11.1 Definitions for Contexts

**activate** (*context\_1, ..., context\_n*) Function

Activates the contexts *context\_1, ..., context\_n*. The facts in these contexts are then available to make deductions and retrieve information. The facts in these contexts are not listed by `facts ()`.

The variable `activecontexts` is the list of contexts which are active by way of the `activate` function.

**activecontexts** Variable

Default value: `[]`

`activecontexts` is a list of the contexts which are active by way of the `activate` function, as opposed to being active because they are subcontexts of the current context.

**assume** (*pred\_1, ..., pred\_n*) Function

Adds predicates *pred\_1, ..., pred\_n* to the current database, after checking for redundancy and inconsistency. If the predicates are consistent and non-redundant, they are added to the data base; if inconsistent or redundant, no action is taken.

`assume` returns a list whose elements are the predicates added to the database and the atoms `redundant` or `inconsistent` where applicable.

**assumescalar** Variable

Default value: `true`

`assumescalar` helps govern whether expressions `expr` for which `nonscalarp (expr)` is `false` are assumed to behave like scalars for certain transformations.

Let `expr` represent any expression other than a list or a matrix, and let `[1, 2, 3]` represent any list or matrix. Then `expr . [1, 2, 3]` yields `[expr, 2 expr, 3 expr]` if `assumescalar` is `true`, or `scalarp (expr)` is `true`, or `constantp (expr)` is `true`.

If `assumescalar` is `true`, such expressions will behave like scalars only for commutative operators, but not for noncommutative multiplication ..

When `assumescalar` is `false`, such expressions will behave like non-scalars.

When `assumescalar` is `all`, such expressions will behave like scalars for all the operators listed above.

**assume\_pos** Variable

Default value: `false`

When `assume_pos` is `true` and the sign of a parameter `x` cannot be determined from the `assume` database or other considerations, `sign` and `asksign (x)` return `true`. This may forestall some automatically-generated `asksign` queries, such as may arise from `integrate` or other computations.



By default, a parameter is  $x$  such that `symbolp (x)` or `subvarp (x)`. The class of expressions considered parameters can be modified to some extent via the variable `assume_pos_pred`.

`sign` and `asksign` attempt to deduce the sign of expressions from the sign of operands within the expression. For example, if `a` and `b` are both positive, then `a + b` is also positive.

However, there is no way to bypass all `asksign` queries. In particular, when the `asksign` argument is a difference `x - y` or a logarithm `log(x)`, `asksign` always requests an input from the user, even when `assume_pos` is `true` and `assume_pos_pred` is a function which returns `true` for all arguments.

### `assume_pos_pred`

Variable

Default value: `false`

When `assume_pos_pred` is assigned the name of a function or a lambda expression of one argument  $x$ , that function is called to determine whether  $x$  is considered a parameter for the purpose of `assume_pos`. `assume_pos_pred` is ignored when `assume_pos` is `false`.

The `assume_pos_pred` function is called by `sign` and `asksign` with an argument  $x$  which is either an atom, a subscripted variable, or a function call expression. If the `assume_pos_pred` function returns `true`,  $x$  is considered a parameter for the purpose of `assume_pos`.

By default, a parameter is  $x$  such that `symbolp (x)` or `subvarp (x)`.

See also `assume` and `assume_pos`.

Examples:

```
(%i1) assume_pos: true$
(%i2) assume_pos_pred: symbolp$
(%i3) sign (a);
(%o3)                                     pos
(%i4) sign (a[1]);
(%o4)                                     pnz
(%i5) assume_pos_pred: lambda ([x], display (x), true)$
(%i6) asksign (a);
                                     x = a
(%o6)                                     pos
(%i7) asksign (a[1]);
                                     x = a
                                     1
(%o7)                                     pos
(%i8) asksign (foo (a));
                                     x = foo(a)
(%o8)                                     pos
(%i9) asksign (foo (a) + bar (b));
                                     x = foo(a)
```

```

                                x = bar(b)

(%o9)                                pos
(%i10) asksign (log (a));          x = a

Is a - 1 positive, negative, or zero?

p;
(%o10)                                pos
(%i11) asksign (a - b);          x = a

                                x = b

                                x = a

                                x = b

Is b - a positive, negative, or zero?

p;
(%o11)                                neg

```

**context**

Variable

Default value: `initial`

`context` names the collection of facts maintained by `assume` and `forget`. `assume` adds facts to the collection named by `context`, while `forget` removes facts.

Binding `context` to a name `foo` changes the current context to `foo`. If the specified context `foo` does not yet exist, it is created automatically by a call to `newcontext`. The specified context is activated automatically.

See `context` for a general description of the context mechanism.

**contexts**

Variable

Default value: `[initial, global]`

`contexts` is a list of the contexts which currently exist, including the currently active context.

The context mechanism makes it possible for a user to bind together and name a selected portion of his database, called a context. Once this is done, the user can have Maxima assume or forget large numbers of facts merely by activating or deactivating their context.

Any symbolic atom can be a context, and the facts contained in that context will be retained in storage until destroyed one by one by calling `forget` or destroyed as a whole by calling `kill` to destroy the context to which they belong.

Contexts exist in a hierarchy, with the root always being the context `global`, which contains information about Maxima that some functions need. When in a given

context, all the facts in that context are "active" (meaning that they are used in deductions and retrievals) as are all the facts in any context which is a subcontext of the active context.

When a fresh Maxima is started up, the user is in a context called `initial`, which has `global` as a subcontext.

See also `facts`, `newcontext`, `supcontext`, `killcontext`, `activate`, `deactivate`, `assume`, and `forget`.

**deactivate** (*context\_1, ..., context\_n*) Function  
Deactivates the specified contexts *context\_1, ..., context\_n*.

**facts** (*item*) Function  
**facts** () Function

If *item* is the name of a context, `facts (item)` returns a list of the facts in the specified context.

If *item* is not the name of a context, `facts (item)` returns a list of the facts known about *item* in the current context. Facts that are active, but in a different context, are not listed.

`facts ()` (i.e., without an argument) lists the current context.

**features** declaration

Maxima recognizes certain mathematical properties of functions and variables. These are called "features".

`declare (x, foo)` gives the property *foo* to the function or variable *x*.

`declare (foo, feature)` declares a new feature *foo*. For example, `declare ([red, green, blue], feature)` declares three new features, `red`, `green`, and `blue`.

The predicate `featurep (x, foo)` returns `true` if *x* has the *foo* property, and `false` otherwise.

The infolist `features` is a list of known features. These are `integer`, `noninteger`, `even`, `odd`, `rational`, `irrational`, `real`, `imaginary`, `complex`, `analytic`, `increasing`, `decreasing`, `oddfun`, `evenfun`, `posfun`, `commutative`, `lassociative`, `rassociative`, `symmetric`, and `antisymmetric`, plus any user-defined features.

`features` is a list of mathematical features. There is also a list of non-mathematical, system-dependent features. See `status`.

**forget** (*pred\_1, ..., pred\_n*) Function  
**forget** (*L*) Function

Removes predicates established by `assume`. The predicates may be expressions equivalent to (but not necessarily identical to) those previously assumed.

`forget (L)`, where *L* is a list of predicates, forgets each item on the list.

**killcontext** (*context\_1, ..., context\_n*) Function  
Kills the contexts *context\_1, ..., context\_n*.

If one of the contexts is the current context, the new current context will become the first available subcontext of the current context which has not been killed. If the first available unkillable context is `global` then `initial` is used instead. If the `initial` context is killed, a new, empty `initial` context is created.

`killcontext` refuses to kill a context which is currently active, either because it is a subcontext of the current context, or by use of the function `activate`.

**newcontext** (*name*) Function

Creates a new, empty context, called *name*, which has `global` as its only subcontext. The newly-created context becomes the currently active context.

**supcontext** (*name, context*) Function

**supcontext** (*name*) Function

Creates a new context, called *name*, which has *context* as a subcontext. *context* must exist.

If *context* is not specified, the current context is assumed.



## 12 Polynomials

### 12.1 Introduction to Polynomials

Polynomials are stored in Maxima either in General Form or as Canonical Rational Expressions (CRE) form. The latter is a standard form, and is used internally by operations such as `factor`, `ratsimp`, and so on.

Canonical Rational Expressions constitute a kind of representation which is especially suitable for expanded polynomials and rational functions (as well as for partially factored polynomials and rational functions when `RATFAC` is set to `true`). In this CRE form an ordering of variables (from most to least main) is assumed for each expression. Polynomials are represented recursively by a list consisting of the main variable followed by a series of pairs of expressions, one for each term of the polynomial. The first member of each pair is the exponent of the main variable in that term and the second member is the coefficient of that term which could be a number or a polynomial in another variable again represented in this form. Thus the principal part of the CRE form of  $3X^2-1$  is  $(X\ 2\ 3\ 0\ -1)$  and that of  $2XY+X-3$  is  $(Y\ 1\ (X\ 1\ 2)\ 0\ (X\ 1\ 1\ 0\ -3))$  assuming  $Y$  is the main variable, and is  $(X\ 1\ (Y\ 1\ 2\ 0\ 1)\ 0\ -3)$  assuming  $X$  is the main variable. "Main"-ness is usually determined by reverse alphabetical order. The "variables" of a CRE expression needn't be atomic. In fact any subexpression whose main operator is not `+`, `-`, `*`, `/` or `^` with integer power will be considered a "variable" of the expression (in CRE form) in which it occurs. For example the CRE variables of the expression  $X+\text{SIN}(X+1)+2\sqrt{X}+1$  are  $X$ ,  $\text{SQRT}(X)$ , and  $\text{SIN}(X+1)$ . If the user does not specify an ordering of variables by using the `RATVARS` function Maxima will choose an alphabetic one. In general, CRE's represent rational expressions, that is, ratios of polynomials, where the numerator and denominator have no common factors, and the denominator is positive. The internal form is essentially a pair of polynomials (the numerator and denominator) preceded by the variable ordering list. If an expression to be displayed is in CRE form or if it contains any subexpressions in CRE form, the symbol `/R/` will follow the line label. See the `RAT` function for converting an expression to CRE form. An extended CRE form is used for the representation of Taylor series. The notion of a rational expression is extended so that the exponents of the variables can be positive or negative rational numbers rather than just positive integers and the coefficients can themselves be rational expressions as described above rather than just polynomials. These are represented internally by a recursive polynomial form which is similar to and is a generalization of CRE form, but carries additional information such as the degree of truncation. As with CRE form, the symbol `/T/` follows the line label of such expressions.

### 12.2 Definitions for Polynomials

#### algebraic

Variable

Default value: `false`

`algebraic` must be set to `true` in order for the simplification of algebraic integers to take effect.

**berlefact**

Variable

Default value: `true`

When `berlefact` is `false` then the Kronecker factoring algorithm will be used otherwise the Berlekamp algorithm, which is the default, will be used.

**bezout** (*p1*, *p2*, *x*)

Function

an alternative to the `resultant` command. It returns a matrix. `determinant` of this matrix is the desired resultant.

**bothcoef** (*expr*, *x*)

Function

Returns a list whose first member is the coefficient of *x* in *expr* (as found by `ratcoef` if *expr* is in CRE form otherwise by `coeff`) and whose second member is the remaining part of *expr*. That is, `[A, B]` where  $expr = A*x + B$ .

Example:

```
(%i1) islinear (expr, x) := block ([c],
      c: bothcoef (rat (expr, x), x),
      is (freeof (x, c) and c[1] # 0))$
(%i2) islinear ((r^2 - (x - r)^2)/x, x);
(%o2) true
```

**coeff** (*expr*, *x*, *n*)

Function

Returns the coefficient of  $x^n$  in *expr*. *n* may be omitted if it is 1. *x* may be an atom, or complete subexpression of *expr* e.g., `sin(x)`, `a[i+1]`, `x + y`, etc. (In the last case the expression `(x + y)` should occur in *expr*). Sometimes it may be necessary to expand or factor *expr* in order to make  $x^n$  explicit. This is not done automatically by `coeff`.

Examples:

```
(%i1) coeff (2*a*tan(x) + tan(x) + b = 5*tan(x) + 3, tan(x));
(%o1) 2 a + 1 = 5
(%i2) coeff (y + x*e^x + 1, x, 0);
(%o2) y + 1
```

**combine** (*expr*)

Function

Simplifies the sum *expr* by combining terms with the same denominator into a single term.

**content** (*p\_1*, *x\_1*, ..., *x\_n*)

Function

Returns a list whose first element is the greatest common divisor of the coefficients of the terms of the polynomial *p\_1* in the variable *x\_n* (this is the content) and whose second element is the polynomial *p\_1* divided by the content.

Examples:

```
(%i1) content (2*x*y + 4*x^2*y^2, y);
(%o1) [2 x, 2 x y + y]
```

**denom** (*expr*) Function  
Returns the denominator of the rational expression *expr*.

**divide** (*p\_1*, *p\_2*, *x\_1*, ..., *x\_n*) Function  
computes the quotient and remainder of the polynomial *p\_1* divided by the polynomial *p\_2*, in a main polynomial variable, *x\_n*. The other variables are as in the **ratvars** function. The result is a list whose first element is the quotient and whose second element is the remainder.

Examples:

```
(%i1) divide (x + y, x - y, x);
(%o1)          [1, 2 y]
(%i2) divide (x + y, x - y);
(%o2)          [- 1, 2 x]
```

Note that *y* is the main variable in the second example.

**eliminate** (*[eqn\_1, ..., eqn\_n]*, *[x\_1, ..., x\_k]*) Function  
Eliminates variables from equations (or expressions assumed equal to zero) by taking successive resultants. This returns a list of  $n - k$  expressions with the  $k$  variables  $x_1, \dots, x_k$  eliminated. First  $x_1$  is eliminated yielding  $n - 1$  expressions, then  $x_2$  is eliminated, etc. If  $k = n$  then a single expression in a list is returned free of the variables  $x_1, \dots, x_k$ . In this case **solve** is called to solve the last resultant for the last variable.

Example:

```
(%i1) expr1: 2*x^2 + y*x + z;
(%o1)          z + x y + 2 x^2
(%i2) expr2: 3*x + 5*y - z - 1;
(%o2)          - z + 5 y + 3 x - 1
(%i3) expr3: z^2 + x - y^2 + 5;
(%o3)          z^2 - y^2 + x + 5
(%i4) eliminate ([expr3, expr2, expr1], [y, z]);
(%o4) [7425 x^8 - 1170 x^7 + 1299 x^6 + 12076 x^5 + 22887 x^4
      - 5154 x^3 - 1291 x^2 + 7688 x + 15376]
```

**ezgcd** (*p\_1*, *p\_2*, *p\_3*, ...) Function  
Returns a list whose first element is the g.c.d of the polynomials *p\_1*, *p\_2*, *p\_3*, ... and whose remaining elements are the polynomials divided by the g.c.d. This always uses the **ezgcd** algorithm.

**facexpand** Variable  
Default value: **true**  
**facexpand** controls whether the irreducible factors returned by **factor** are in expanded (the default) or recursive (normal CRE) form.



**factcomb** (*expr*)

Function

Tries to combine the coefficients of factorials in *expr* with the factorials themselves by converting, for example,  $(n + 1)*n!$  into  $(n + 1)!$ .

`sumsplitfact` if set to `false` will cause `minfactorial` to be applied after a `factcomb`.

**factor** (*expr*)

Function

Factors the expression *expr*, containing any number of variables or functions, into factors irreducible over the integers. `factor (expr, p)` factors *expr* over the field of integers with an element adjoined whose minimum polynomial is *p*.

`factorflag` if `false` suppresses the factoring of integer factors of rational expressions.

`dontfactor` may be set to a list of variables with respect to which factoring is not to occur. (It is initially empty). Factoring also will not take place with respect to any variables which are less important (using the variable ordering assumed for CRE form) than those on the `dontfactor` list.

`savefactors` if `true` causes the factors of an expression which is a product of factors to be saved by certain functions in order to speed up later factorizations of expressions containing some of the same factors.

`berlefact` if `false` then the Kronecker factoring algorithm will be used otherwise the Berlekamp algorithm, which is the default, will be used.

`intfaclim` is the largest divisor which will be tried when factoring a bignum integer. If set to `false` (this is the case when the user calls `factor` explicitly), or if the integer is a fixnum (i.e. fits in one machine word), complete factorization of the integer will be attempted. The user's setting of `intfaclim` is used for internal calls to `factor`. Thus, `intfaclim` may be reset to prevent Maxima from taking an inordinately long time factoring large integers.

Examples:

```
(%i1) factor (2^63 - 1);
```

```
2
```

```
(%o1) 7 73 127 337 92737 649657
```

```
(%i2) factor (-8*y - 4*x + z^2*(2*y + x));
```

```
(%o2) (2 y + x) (z - 2) (z + 2)
```

```
(%i3) -1 - 2*x - x^2 + y^2 + 2*x*y^2 + x^2*y^2;
```

```
2 2 2 2 2
```

```
(%o3) x y + 2 x y + y - x - 2 x - 1
```

```
(%i4) block ([dontfactor: [x]], factor (%/36/(1 + 2*y + y^2)));
```

```
2
```

```
(x + 2 x + 1) (y - 1)
```

```
(%o4) -----
```

```
36 (y + 1)
```

```
(%i5) factor (1 + %e^(3*x));
```

```
x 2 x x
```

```
(%o5) (%e + 1) (%e - %e + 1)
```

```
(%i6) factor (1 + x^4, a^2 - 2);
```

```
2
```

```
2
```

```
(%o6) (x - a x + 1) (x + a x + 1)
```

```
(%i7) factor (-y^2*z^2 - x*z^2 + x^2*y^2 + x^3);
```

```

(%o7)          2
      - (y + x) (z - x) (z + x)
(%i8) (2 + x)/(3 + x)/(b + x)/(c + x)^2;
          x + 2
(%o8) -----
          2
      (x + 3) (x + b) (x + c)
(%i9) ratsimp (%);
          4          3
(%o9) (x + 2)/(x + (2 c + b + 3) x
      2          2          2          2
+ (c + (2 b + 6) c + 3 b) x + ((b + 3) c + 6 b c) x + 3 b c )
(%i10) partfrac (% , x);
          2          4          3
(%o10) - (c - 4 c - b + 6)/((c + (- 2 b - 6) c
      2          2          2          2
+ (b + 12 b + 9) c + (- 6 b - 18 b) c + 9 b ) (x + c))
      c - 2
-----
      2          2
(c + (- b - 3) c + 3 b) (x + c)
+ -----
      b - 2
      ((b - 3) c + (6 b - 2 b ) c + b - 3 b ) (x + b)
- -----
      1
      ((b - 3) c + (18 - 6 b) c + 9 b - 27) (x + 3)
(%i11) map ('factor, %);
          2          c - 2
(%o11) - ----- - -----
          2          2          2          2
      (c - 3) (c - b) (x + c) (c - 3) (c - b) (x + c)
+ ----- - -----
          b - 2          1
          2          2
      (b - 3) (c - b) (x + b) (b - 3) (c - 3) (x + 3)
(%i12) ratsimp ((x^5 - 1)/(x - 1));
          4          3          2
(%o12) x + x + x + x + 1
(%i13) subst (a, x, %);

```

```

(%o13)          4 3 2
              a + a + a + a + 1
(%i14) factor (%th(2), %);
(%o14) (x - a) (x - a) (x - a) (x + a + a + a + 1)
(%i15) factor (1 + x^12);
(%o15)          4      8 4
              (x + 1) (x - x + 1)
(%i16) factor (1 + x^99);
(%o16) (x + 1) (x - x + 1) (x - x + 1)

          10 9 8 7 6 5 4 3 2
(x - x + x - x + x - x + x - x + x - x + 1)

          20 19 17 16 14 13 11 10 9 7 6
(x + x - x - x + x + x - x - x - x + x + x

          4 3      60 57 51 48 42 39 33
- x - x + x + 1) (x + x - x - x + x + x - x

          30 27 21 18 12 9 3
- x - x + x + x - x - x + x + 1)

```

**factorflag**

Variable

Default value: false

When `factorflag` is false, suppresses the factoring of integer factors of rational expressions.

**factorout** (*expr*, *x*<sub>1</sub>, *x*<sub>2</sub>, ...)

Function

Rearranges the sum *expr* into a sum of terms of the form *f* (*x*<sub>1</sub>, *x*<sub>2</sub>, ...) \* *g* where *g* is a product of expressions not containing any *x*<sub>*i*</sub> and *f* is factored.

**factorsum** (*expr*)

Function

Tries to group terms in factors of *expr* which are sums into groups of terms such that their sum is factorable. `factorsum` can recover the result of `expand ((x + y)^2 + (z + w)^2)` but it can't recover `expand ((x + 1)^2 + (x + y)^2)` because the terms have variables in common.

Example:

```

(%i1) expand ((x + 1)*((u + v)^2 + a*(w + z)^2));
(%o1) a x z + a z + 2 a w x z + 2 a w z + a w x + v x
      2      2      2      2
      + 2 u v x + u x + a w + v + 2 u v + u
(%i2) factorsum (%);
(%o2) (x + 1) (a (z + w) + (v + u) )
      2      2

```

**fasttimes** (*p*<sub>1</sub>, *p*<sub>2</sub>)

Function

Returns the product of the polynomials *p*<sub>1</sub> and *p*<sub>2</sub> by using a special algorithm for multiplication of polynomials. *p*<sub>1</sub> and *p*<sub>2</sub> should be multivariate, dense, and nearly the same size. Classical multiplication is of order *n*<sub>1</sub> *n*<sub>2</sub> where *n*<sub>1</sub> is the degree of *p*<sub>1</sub> and *n*<sub>2</sub> is the degree of *p*<sub>2</sub>. **fasttimes** is of order  $\max(n_1, n_2)^{1.585}$ .

**fullratsimp** (*expr*)

Function

**fullratsimp** repeatedly applies **ratsimp** followed by non-rational simplification to an expression until no further change occurs, and returns the result.

When non-rational expressions are involved, one call to **ratsimp** followed as is usual by non-rational ("general") simplification may not be sufficient to return a simplified result. Sometimes, more than one such call may be necessary. **fullratsimp** makes this process convenient.

**fullratsimp** (*expr*, *x*<sub>1</sub>, ..., *x*<sub>*n*</sub>) takes one or more arguments similar to **ratsimp** and **rat**.

Example:

```
(%i1) expr: (x^(a/2) + 1)^2*(x^(a/2) - 1)^2/(x^a - 1);
```

```

              a/2      2      a/2      2
          (x      - 1) (x      + 1)
(%o1)  -----

```

```

              a
              x  - 1
(%i2) ratsimp (expr);
```

```

          2 a      a
          x  - 2 x  + 1
(%o2)  -----

```

```

              a
              x  - 1
(%i3) fullratsimp (expr);
```

```

              a
              x  - 1
(%o3)
```

```
(%i4) rat (expr);
```

```

          a/2 4      a/2 2
          (x  ) - 2 (x  ) + 1
(%o4) /R/  -----

```

```

              a
              x  - 1

```

**fullratsubst** (*a*, *b*, *c*)

Function

is the same as **ratsubst** except that it calls itself recursively on its result until that result stops changing. This function is useful when the replacement expression and the replaced expression have one or more variables in common.

**fullratsubst** will also accept its arguments in the format of **lratsubst**. That is, the first argument may be a single substitution equation or a list of such equations, while the second argument is the expression being processed.

**load** ("lrats") loads **fullratsubst** and **lratsubst**.

Examples:

- ```
(%i1) load ("lrats")$
```
- `subst` can carry out multiple substitutions. `lratsubst` is analogous to `subst`.

```
(%i2) subst ([a = b, c = d], a + c);
(%o2)          d + b
(%i3) lratsubst ([a^2 = b, c^2 = d], (a + e)*c*(a + c));
(%o3)          (d + a c) e + a d + b c
```
  - If only one substitution is desired, then a single equation may be given as first argument.

```
(%i4) lratsubst (a^2 = b, a^3);
(%o4)          a b
```
  - `fullratsubst` is equivalent to `ratsubst` except that it recurses until its result stops changing.

```
(%i5) ratsubst (b*a, a^2, a^3);
(%o5)          2
          a b
(%i6) fullratsubst (b*a, a^2, a^3);
(%o6)          2
          a b
```
  - `fullratsubst` also accepts a list of equations or a single equation as first argument.

```
(%i7) fullratsubst ([a^2 = b, b^2 = c, c^2 = a], a^3*b*c);
(%o7)          b
(%i8) fullratsubst (a^2 = b*a, a^3);
(%o8)          2
          a b
```
  - `fullratsubst` may cause an indefinite recursion.

```
(%i9) errcatch (fullratsubst (b*a^2, a^2, a^3));

*** - Lisp stack overflow. RESET
```

**gcd** ( $p_1, p_2, x_1, \dots$ )

Function

Returns the greatest common divisor of  $p_1$  and  $p_2$ . The flag `gcd` determines which algorithm is employed. Setting `gcd` to `ez`, `eez`, `subres`, `red`, or `smod` selects the `ezgcd`, New `eez gcd`, `subresultant prs`, `reduced`, or modular algorithm, respectively. If `gcd false` then `GCD(p1,p2,var)` will always return 1 for all `var`. Many functions (e.g. `ratsimp`, `factor`, etc.) cause `gcd`'s to be taken implicitly. For homogeneous polynomials it is recommended that `gcd` equal to `subres` be used. To take the `gcd` when an algebraic is present, e.g. `GCD(X^2-2*SQRT(2)*X+2,X-SQRT(2))`; , `algebraic` must be `true` and `gcd` must not be `ez`. `subres` is a new algorithm, and people who have been using the `red` setting should probably change it to `subres`.

The `gcd` flag, default: `subres`, if `false` will also prevent the greatest common divisor from being taken when expressions are converted to canonical rational expression (CRE) form. This will sometimes speed the calculation if `gcds` are not required.

**gcdex** (*f*, *g*)

Function

**gcdex** (*f*, *g*, *x*)

Function

Returns a list [*a*, *b*, *u*] where *u* is the greatest common divisor (gcd) of *f* and *g*, and *u* is equal to  $a f + b g$ . The arguments *f* and *g* should be univariate polynomials, or else polynomials in *x* a supplied **main** variable since we need to be in a principal ideal domain for this to work. The gcd means the gcd regarding *f* and *g* as univariate polynomials with coefficients being rational functions in the other variables.

**gcdex** implements the Euclidean algorithm, where we have a sequence of  $L[i]$ : [*a*[*i*], *b*[*i*], *r*[*i*]] which are all perpendicular to [*f*, *g*, -1] and the next one is built as if  $q = \text{quotient}(r[i]/r[i+1])$  then  $L[i+2]: L[i] - q L[i+1]$ , and it terminates at  $L[i+1]$  when the remainder  $r[i+2]$  is zero.

```
(%i1) gcdex (x^2 + 1, x^3 + 4);
              2
              x  + 4 x - 1  x + 4
(%o1)/R/      [- -----, -----, 1]
              17          17
(%i2) % . [x^2 + 1, x^3 + 4, -1];
(%o2)/R/      0
```

Note that the gcd in the following is 1 since we work in  $k(y)[x]$ , not the  $y+1$  we would expect in  $k[y, x]$ .

```
(%i1) gcdex (x*(y + 1), y^2 - 1, x);
              1
(%o1)/R/      [0, -----, 1]
              2
              y  - 1
```

**gcfactor** (*n*)

Function

Factors the Gaussian integer *n* over the Gaussian integers, i.e., numbers of the form  $a + b \%i$  where *a* and *b* are rational integers (i.e., ordinary integers). Factors are normalized by making *a* and *b* non-negative.

**gfactor** (*expr*)

Function

Factors the polynomial *expr* over the Gaussian integers (that is, the integers with the imaginary unit  $\%i$  adjoined). This is like **factor** (*expr*,  $a^2+1$ ) where *a* is  $\%i$ .

Example:

```
(%i1) gfactor (x^4 - 1);
(%o1)      (x - 1) (x + 1) (x - \%i) (x + \%i)
```

**gfactorsum** (*expr*)

Function

is similar to **factorsum** but applies **gfactor** instead of **factor**.

**hipow** (*expr*, *x*)

Function

Returns the highest explicit exponent of *x* in *expr*. *x* may be a variable or a general expression. If *x* does not appear in *expr*, **hipow** returns 0.

`hipow` does not consider expressions equivalent to `expr`. In particular, `hipow` does not expand `expr`, so `hipow (expr, x)` and `hipow (expand (expr, x))` may yield different results.

Examples:

```
(%i1) hipow (y^3 * x^2 + x * y^4, x);
(%o1)          2
(%i2) hipow ((x + y)^5, x);
(%o2)          1
(%i3) hipow (expand ((x + y)^5), x);
(%o3)          5
(%i4) hipow ((x + y)^5, x + y);
(%o4)          5
(%i5) hipow (expand ((x + y)^5), x + y);
(%o5)          0
```

### **intfaclim**

Variable

Default value: 1000

`intfaclim` is the largest divisor which will be tried when factoring a bignum integer.

When `intfaclim` is `false` (this is the case when the user calls `factor` explicitly), or if the integer is a fixnum (i.e., fits in one machine word), factors of any size are considered. `intfaclim` is set to `false` when factors are computed in `divsum`, `totient`, and `primep`.

Internal calls to `factor` respect the user-specified value of `intfaclim`. Setting `intfaclim` to a smaller value may reduce the time spent factoring large integers.

### **keepfloat**

Variable

Default value: `false`

When `keepfloat` is `true`, prevents floating point numbers from being rationalized when expressions which contain them are converted to canonical rational expression (CRE) form.

### **lratsubst** (*L*, *expr*)

Function

is analogous to `subst` (*L*, *expr*) except that it uses `ratsubst` instead of `subst`.

The first argument of `lratsubst` is an equation or a list of equations identical in format to that accepted by `subst`. The substitutions are made in the order given by the list of equations, that is, from left to right.

`load ("lrats")` loads `fullratsubst` and `lratsubst`.

Examples:

```
(%i1) load ("lrats")$
```

- `subst` can carry out multiple substitutions. `lratsubst` is analogous to `subst`.

```
(%i2) subst ([a = b, c = d], a + c);
(%o2)          d + b
(%i3) lratsubst ([a^2 = b, c^2 = d], (a + e)*c*(a + c));
(%o3)          (d + a c) e + a d + b c
```

- If only one substitution is desired, then a single equation may be given as first argument.

```
(%i4) lratsubst (a^2 = b, a^3);
(%o4) a b
```

**modulus**

Variable

Default value: `false`

When `modulus` is a positive number  $p$ , operations on rational numbers (as returned by `rat` and related functions) are carried out modulo  $p$ , using the so-called "balanced" modulus system in which  $n$  modulo  $p$  is defined as an integer  $k$  in  $[-(p-1)/2, \dots, 0, \dots, (p-1)/2]$  when  $p$  is odd, or  $[-(p/2 - 1), \dots, 0, \dots, p/2]$  when  $p$  is even, such that  $a p + k$  equals  $n$  for some integer  $a$ .

If `expr` is already in canonical rational expression (CRE) form when `modulus` is reset, then you may need to re-rat `expr`, e.g., `expr: rat (ratdisrep (expr))`, in order to get correct results.

Typically `modulus` is set to a prime number. If `modulus` is set to a positive non-prime integer, this setting is accepted, but a warning message is displayed. Maxima will allow zero or a negative integer to be assigned to `modulus`, although it is not clear if that has any useful consequences.

**num** (`expr`)

Function

Returns the numerator of `expr` if it is a ratio. If `expr` is not a ratio, `expr` is returned. `num` evaluates its argument.

**quotient** (`p_1`, `p_2`)

Function

**quotient** (`p_1`, `p_2`, `x_1`, ..., `x_n`)

Function

Returns the polynomial `p_1` divided by the polynomial `p_2`. The arguments `x_1`, ..., `x_n` are interpreted as in `ratvars`.

`quotient` returns the first element of the two-element list returned by `divide`.

**rat** (`expr`)

Function

**rat** (`expr`, `x_1`, ..., `x_n`)

Function

Converts `expr` to canonical rational expression (CRE) form by expanding and combining all terms over a common denominator and cancelling out the greatest common divisor of the numerator and denominator, as well as converting floating point numbers to rational numbers within a tolerance of `ratepsilon`. The variables are ordered according to the `x_1`, ..., `x_n`, if specified, as in `ratvars`.

`rat` does not generally simplify functions other than addition `+`, subtraction `-`, multiplication `*`, division `/`, and exponentiation to an integer power, whereas `ratsimp` does handle those cases. Note that atoms (numbers and variables) in CRE form are not the same as they are in the general form. For example, `rat(x) - x` yields `rat(0)` which has a different internal representation than `0`.

When `ratfac` is `true`, `rat` yields a partially factored form for CRE. During rational operations the expression is maintained as fully factored as possible without an actual call to the factor package. This should always save space and may save some



time in some computations. The numerator and denominator are still made relatively prime (e.g. `rat ((x^2 - 1)^4/(x + 1)^2)` yields  $(x - 1)^4 (x + 1)^2$ ), but the factors within each part may not be relatively prime.

`ratprint` if `false` suppresses the printout of the message informing the user of the conversion of floating point numbers to rational numbers.

`keepfloat` if `true` prevents floating point numbers from being converted to rational numbers.

See also `ratexpand` and `ratsimp`.

Examples:

```
(%i1) ((x - 2*y)^4/(x^2 - 4*y^2)^2 + 1)*(y + a)*(2*y + x) / (4*y^2 + x^2);
```

$$\frac{(y + a)(2y + x)\left(\frac{(x - 2y)^4}{(x^2 - 4y^2)^2} + 1\right)}{4y^2 + x^2}$$

```
(%o1)
```

```
(%i2) rat (%o1, y, a, x);
```

$$\frac{2a + 2y}{x + 2y}$$

```
(%o2)/R/
```

### **rataldenom**

Variable

Default value: `true`

When `rataldenom` is `true`, allows rationalization of denominators with respect to radicals to take effect. `rataldenom` has an effect only when canonical rational expressions (CRE) are used in algebraic mode.

### **ratcoef** (*expr*, *x*, *n*)

Function

### **ratcoef** (*expr*, *x*)

Function

Returns the coefficient of the expression  $x^n$  in the expression *expr*. If omitted, *n* is assumed to be 1.

The return value is free (except possibly in a non-rational sense) of the variables in *x*. If no coefficient of this type exists, 0 is returned.

`ratcoef` expands and rationally simplifies its first argument and thus it may produce answers different from those of `coeff` which is purely syntactic. Thus `RATCOEF((X+1)/Y+X,X)` returns  $(Y+1)/Y$  whereas `coeff` returns 1.

`ratcoef` (*expr*, *x*, 0), viewing *expr* as a sum, returns a sum of those terms which do not contain *x*. Therefore if *x* occurs to any negative powers, `ratcoef` should not be used.

Since *expr* is rationally simplified before it is examined, coefficients may not appear quite the way they were envisioned.

Example:

```
(%i1) s: a*x + b*x + 5$
(%i2) ratcoef (s, a + b);
(%o2) x
```

**ratdenom** (*expr*)

Function

Returns the denominator of *expr*, after coercing *expr* to a canonical rational expression (CRE). The return value is a CRE.

*expr* is coerced to a CRE by `rat` if it is not already a CRE. This conversion may change the form of *expr* by putting all terms over a common denominator.

`denom` is similar, but returns an ordinary expression instead of a CRE. Also, `denom` does not attempt to place all terms over a common denominator, and thus some expressions which are considered ratios by `ratdenom` are not considered ratios by `denom`.

**ratdenomdivide**

Variable

Default value: `true`

When `ratdenomdivide` is `true`, `ratexpand` expands a ratio in which the numerator is a sum into a sum of ratios, all having a common denominator. Otherwise, `ratexpand` collapses a sum of ratios into a single ratio, the numerator of which is the sum of the numerators of each ratio.

Examples:

```
(%i1) expr: (x^2 + x + 1)/(y^2 + 7);
(%o1)

$$\frac{x^2 + x + 1}{y^2 + 7}$$

(%i2) ratdenomdivide: true$
(%i3) ratexpand (expr);
(%o3)

$$\frac{x^2}{y^2 + 7} + \frac{x}{y^2 + 7} + \frac{1}{y^2 + 7}$$

(%i4) ratdenomdivide: false$
(%i5) ratexpand (expr);
(%o5)

$$\frac{x^2 + x + 1}{y^2 + 7}$$

(%i6) expr2: a^2/(b^2 + 3) + b/(b^2 + 3);
(%o6)

$$\frac{b}{b^2 + 3} + \frac{a}{b^2 + 3}$$

```

```
(%i7) ratexpand (expr2);
```

```
(%o7)
          2
        b + a
        -----
          2
        b  + 3
```

**ratdiff** (*expr*, *x*) Function

Differentiates the rational expression *expr* with respect to *x*. *expr* must be a ratio of polynomials or a polynomial in *x*. The argument *x* may be a variable or a subexpression of *expr*.

The result is equivalent to `diff`, although perhaps in a different form. `ratdiff` may be faster than `diff`, for rational expressions.

`ratdiff` returns a canonical rational expression (CRE) if *expr* is a CRE. Otherwise, `ratdiff` returns a general expression.

`ratdiff` considers only the dependence of *expr* on *x*, and ignores any dependencies established by `depends`.

Example:

```
(%i1) expr: (4*x^3 + 10*x - 11)/(x^5 + 5);
```

```
(%o1)
          3
        4 x  + 10 x - 11
        -----
          5
          x  + 5
```

```
(%i2) ratdiff (expr, x);
```

```
(%o2)
          7          5          4          2
        8 x  + 40 x  - 55 x  - 60 x  - 50
        -----
          10          5
          x  + 10 x  + 25
```

```
(%i3) expr: f(x)^3 - f(x)^2 + 7;
```

```
(%o3)
          3          2
        f (x) - f (x) + 7
```

```
(%i4) ratdiff (expr, f(x));
```

```
(%o4)
          2
        3 f (x) - 2 f(x)
```

```
(%i5) expr: (a + b)^3 + (a + b)^2;
```

```
(%o5)
          3          2
        (b + a)  + (b + a)
```

```
(%i6) ratdiff (expr, a + b);
```

```
(%o6)
          2          2
        3 b  + (6 a + 2) b + 3 a  + 2 a
```

**ratdisrep** (*expr*) Function

Returns its argument as a general expression. If *expr* is a general expression, it is returned unchanged.

Typically `ratdisrep` is called to convert a canonical rational expression (CRE) into a general expression. This is sometimes convenient if one wishes to stop the "contagion", or use rational functions in non-rational contexts.

See also `totaldisrep`.

### **ratepsilon**

Variable

Default value: 2.0e-8

`ratepsilon` is the tolerance used in the conversion of floating point numbers to rational numbers.

### **ratexpand** (*expr*)

Function

### **ratexpand**

Variable

Expands *expr* by multiplying out products of sums and exponentiated sums, combining fractions over a common denominator, cancelling the greatest common divisor of the numerator and denominator, then splitting the numerator (if a sum) into its respective terms divided by the denominator.

The return value of `ratexpand` is a general expression, even if *expr* is a canonical rational expression (CRE).

The switch `ratexpand` if `true` will cause CRE expressions to be fully expanded when they are converted back to general form or displayed, while if it is `false` then they will be put into a recursive form. See also `ratsimp`.

When `ratdenomdivide` is `true`, `ratexpand` expands a ratio in which the numerator is a sum into a sum of ratios, all having a common denominator. Otherwise, `ratexpand` collapses a sum of ratios into a single ratio, the numerator of which is the sum of the numerators of each ratio.

When `keepfloat` is `true`, prevents floating point numbers from being rationalized when expressions which contain them are converted to canonical rational expression (CRE) form.

Examples:

```
(%i1) ratexpand ((2*x - 3*y)^3);
(%o1)          3      2      2      3
      - 27 y  + 54 x y  - 36 x  y + 8 x
(%i2) expr: (x - 1)/(x + 1)^2 + 1/(x - 1);
(%o2)          x - 1      1
      ----- + -----
              2      x - 1
              (x + 1)
(%i3) expand (expr);
(%o3)          x          1          1
      ----- - ----- + -----
              2          2          x - 1
              x  + 2 x + 1  x  + 2 x + 1
(%i4) ratexpand (expr);
(%o4)          2      2
              2 x          2
      ----- + -----
```

$$\frac{x^3 + x^2 - x - 1}{x^3 + x^2 - x - 1}$$

**ratfac**

Variable

Default value: `false`

When `ratfac` is `true`, canonical rational expressions (CRE) are manipulated in a partially factored form.

During rational operations the expression is maintained as fully factored as possible without calling `factor`. This should always save space and may save time in some computations. The numerator and denominator are made relatively prime, for example `rat ((x^2 - 1)^4/(x + 1)^2)` yields  $(x - 1)^4 (x + 1)^2$ , but the factors within each part may not be relatively prime.

In the `ctensr` (Component Tensor Manipulation) package, Ricci, Einstein, Riemann, and Weyl tensors and the scalar curvature are factored automatically when `ratfac` is `true`. *ratfac should only be set for cases where the tensorial components are known to consist of few terms.*

The `ratfac` and `ratweight` schemes are incompatible and may not both be used at the same time.

**ratnumer** (*expr*)

Function

Returns the numerator of *expr*, after coercing *expr* to a canonical rational expression (CRE). The return value is a CRE.

*expr* is coerced to a CRE by `rat` if it is not already a CRE. This conversion may change the form of *expr* by putting all terms over a common denominator.

`num` is similar, but returns an ordinary expression instead of a CRE. Also, `num` does not attempt to place all terms over a common denominator, and thus some expressions which are considered ratios by `ratnumer` are not considered ratios by `num`.

**ratnump** (*expr*)

Function

Returns `true` if *expr* is a literal integer or ratio of literal integers, otherwise `false`.

**ratp** (*expr*)

Function

Returns `true` if *expr* is a canonical rational expression (CRE) or extended CRE, otherwise `false`.

CRE are created by `rat` and related functions. Extended CRE are created by `taylor` and related functions.

**ratprint**

Variable

Default value: `true`

When `ratprint` is `true`, a message informing the user of the conversion of floating point numbers to rational numbers is displayed.

**ratsimp** (*expr*) Function  
**ratsimp** (*expr*, *x\_1*, ..., *x\_n*) Function

Simplifies the expression *expr* and all of its subexpressions, including the arguments to non-rational functions. The result is returned as the quotient of two polynomials in a recursive form, that is, the coefficients of the main variable are polynomials in the other variables. Variables may include non-rational functions (e.g., `sin (x^2 + 1)`) and the arguments to any such functions are also rationally simplified.

`ratsimp (expr, x_1, ..., x_n)` enables rational simplification with the specification of variable ordering as in `ratvars`.

When `ratsimpexpons` is `true`, `ratsimp` is applied to the exponents of expressions during simplification.

See also `ratexpand`. Note that `ratsimp` is affected by some of the flags which affect `ratexpand`.

Examples:

```
(%i1) sin (x/(x^2 + x)) = exp ((log(x) + 1)^2 - log(x)^2);
                                2      2
                                (log(x) + 1) - log (x)
(%o1)      sin(-----) = %e
              2
              x  + x
(%i2) ratsimp (%);
(%o2)      sin(-----) = %e x
              1      2
              x + 1
(%i3) ((x - 1)^(3/2) - (x + 1)*sqrt(x - 1))/sqrt((x - 1)*(x + 1));
              3/2
              (x - 1) - sqrt(x - 1) (x + 1)
(%o3)      -----
              sqrt((x - 1) (x + 1))
(%i4) ratsimp (%);
              2 sqrt(x - 1)
(%o4)      - -----
              2
              sqrt(x - 1)
(%i5) x^(a + 1/a), ratsimpexpons: true;
              2
              a + 1
              -----
              a
(%o5)      x
```

**ratsimpexpons** Variable

Default value: `false`

When `ratsimpexpons` is `true`, `ratsimp` is applied to the exponents of expressions during simplification.

**ratsubst** (*a*, *b*, *c*) Function

Substitutes *a* for *b* in *c* and returns the resulting expression. *b* may be a sum, product, power, etc.

**ratsubst** knows something of the meaning of expressions whereas **subst** does a purely syntactic substitution. Thus **subst** (*a*, *x* + *y*, *x* + *y* + *z*) returns *x* + *y* + *z* whereas **ratsubst** returns *z* + *a*.

When **radsubstflag** is true, **ratsubst** makes substitutions for radicals in expressions which don't explicitly contain them.

Examples:

```
(%i1) ratsubst (a, x*y^2, x^4*y^3 + x^4*y^8);
              3      4
(%o1)          a x y + a
(%i2) cos(x)^4 + cos(x)^3 + cos(x)^2 + cos(x) + 1;
              4      3      2
(%o2)    cos (x) + cos (x) + cos (x) + cos(x) + 1
(%i3) ratsubst (1 - sin(x)^2, cos(x)^2, %);
              4      2      2
(%o3)    sin (x) - 3 sin (x) + cos(x) (2 - sin (x)) + 3
(%i4) ratsubst (1 - cos(x)^2, sin(x)^2, sin(x)^4);
              4      2
(%o4)          cos (x) - 2 cos (x) + 1
(%i5) radsubstflag: false$
(%i6) ratsubst (u, sqrt(x), x);
(%o6)          x
(%i7) radsubstflag: true$
(%i8) ratsubst (u, sqrt(x), x);
              2
(%o8)          u
```

**ratvars** (*x<sub>1</sub>*, ..., *x<sub>n</sub>*) Function

**ratvars** () Function

**ratvars** Variable

Declares main variables *x<sub>1</sub>*, ..., *x<sub>n</sub>* for rational expressions. *x<sub>n</sub>*, if present in a rational expression, is considered the main variable. Otherwise, *x<sub>[n-1]</sub>* is considered the main variable if present, and so on through the preceding variables to *x<sub>1</sub>*, which is considered the main variable only if none of the succeeding variables are present.

If a variable in a rational expression is not present in the **ratvars** list, it is given a lower priority than *x<sub>1</sub>*.

The arguments to **ratvars** can be either variables or non-rational functions such as **sin**(*x*).

The variable **ratvars** is a list of the arguments of the function **ratvars** when it was called most recently. Each call to the function **ratvars** resets the list. **ratvars** () clears the list.

**ratweight** ( $x_1, w_1, \dots, x_n, w_n$ ) Function  
**ratweight** () Function

Assigns a weight  $w_i$  to the variable  $x_i$ . This causes a term to be replaced by 0 if its weight exceeds the value of the variable `ratwtlvl` (default yields no truncation). The weight of a term is the sum of the products of the weight of a variable in the term times its power. For example, the weight of  $3 x_1^2 x_2$  is  $2 w_1 + w_2$ . Truncation according to `ratwtlvl` is carried out only when multiplying or exponentiating canonical rational expressions (CRE).

`ratweight` () returns the cumulative list of weight assignments.

Note: The `ratfac` and `ratweight` schemes are incompatible and may not both be used at the same time.

Examples:

```
(%i1) ratweight (a, 1, b, 1);
(%o1) [a, 1, b, 1]
(%i2) expr1: rat(a + b + 1)$
(%i3) expr1^2;
(%o3)/R/      2          2
      b  + (2 a + 2) b + a  + 2 a + 1
(%i4) ratwtlvl: 1$
(%i5) expr1^2;
(%o5)/R/      2 b + 2 a + 1
```

**ratweights** Variable

Default value: []

`ratweights` is the list of weights assigned by `ratweight`. The list is cumulative: each call to `ratweight` places additional items in the list.

`kill (ratweights)` and `save (ratweights)` both work as expected.

**ratwtlvl** Variable

Default value: `false`

`ratwtlvl` is used in combination with the `ratweight` function to control the truncation of canonical rational expressions (CRE). For the default value of `false`, no truncation occurs.

**remainder** ( $p_1, p_2$ ) Function

**remainder** ( $p_1, p_2, x_1, \dots, x_n$ ) Function

Returns the remainder of the polynomial  $p_1$  divided by the polynomial  $p_2$ . The arguments  $x_1, \dots, x_n$  are interpreted as in `ratvars`.

`remainder` returns the second element of the two-element list returned by `divide`.

**resultant** ( $p_1, p_2, x$ ) Function

**resultant** Variable

Computes the resultant of the two polynomials  $p_1$  and  $p_2$ , eliminating the variable  $x$ . The resultant is a determinant of the coefficients of  $x$  in  $p_1$  and  $p_2$ , which equals zero if and only if  $p_1$  and  $p_2$  have a non-constant factor in common.



If  $p_1$  or  $p_2$  can be factored, it may be desirable to call `factor` before calling `resultant`.

The variable `resultant` controls which algorithm will be used to compute the resultant. `subres` for subresultant prs, `mod` for modular resultant algorithm, and `red` for reduced prs. On most problems `subres` should be best. On some large degree univariate or bivariate problems `mod` may be better.

The function `bezout` takes the same arguments as `resultant` and returns a matrix. The determinant of the return value is the desired resultant.

### `savefactors`

Variable

Default value: `false`

When `savefactors` is `true`, causes the factors of an expression which is a product of factors to be saved by certain functions in order to speed up later factorizations of expressions containing some of the same factors.

### `sqfr` (*expr*)

Function

is similar to `factor` except that the polynomial factors are "square-free." That is, they have factors only of degree one. This algorithm, which is also used by the first stage of `factor`, utilizes the fact that a polynomial has in common with its  $n$ 'th derivative all its factors of degree greater than  $n$ . Thus by taking greatest common divisors with the polynomial of the derivatives with respect to each variable in the polynomial, all factors of degree greater than 1 can be found.

Example:

```
(%i1) sqfr (4*x^4 + 4*x^3 - 3*x^2 - 4*x - 1);
              2      2
(%o1)          (2 x + 1) (x - 1)
```

### `tellrat` ( $p_1, \dots, p_n$ )

Function

### `tellrat` ()

Function

Adds to the ring of algebraic integers known to Maxima the elements which are the solutions of the polynomials  $p_1, \dots, p_n$ . Each argument  $p_i$  is a polynomial with integer coefficients.

`tellrat (x)` effectively means substitute 0 for  $x$  in rational functions.

`tellrat ()` returns a list of the current substitutions.

`algebraic` must be set to `true` in order for the simplification of algebraic integers to take effect.

Maxima initially knows about the imaginary unit `%i` and all roots of integers.

There is a command `untellrat` which takes kernels and removes `tellrat` properties.

When `tellrat`'ing a multivariate polynomial, e.g., `tellrat (x^2 - y^2)`, there would be an ambiguity as to whether to substitute  $y^2$  for  $x^2$  or vice versa. Maxima picks a particular ordering, but if the user wants to specify which, e.g. `tellrat (y^2 = x^2)` provides a syntax which says replace  $y^2$  by  $x^2$ .

Examples:

```

(%i1) 10*(%i + 1)/(%i + 3^(1/3));
(%o1)

$$\frac{10 (\%i + 1)}{\sqrt[3]{\%i + 3}}$$

(%i2) ev (ratdisrep (rat(%)), algebraic);
(%o2)

$$(4 \sqrt[3]{3} - 2 \sqrt[3]{3} - 4) \%i + 2 \sqrt[3]{3} + 4 \sqrt[3]{3} - 2$$

(%i3) tellrat (1 + a + a^2);
(%o3)

$$[a^2 + a + 1]$$

(%i4) 1/(a*sqrt(2) - 1) + a/(sqrt(3) + sqrt(2));
(%o4)

$$\frac{1}{\sqrt{2} a - 1} + \frac{a}{\sqrt{3} + \sqrt{2}}$$

(%i5) ev (ratdisrep (rat(%)), algebraic);
(%o5)

$$\frac{(7 \sqrt{3} - 10 \sqrt{2} + 2) a - 2 \sqrt{2} - 1}{7}$$

(%i6) tellrat (y^2 = x^2);
(%o6)

$$[y^2 - x^2, a^2 + a + 1]$$


```

**totaldisrep** (*expr*)

Function

Converts every subexpression of *expr* from canonical rational expressions (CRE) to general form and returns the result. If *expr* is itself in CRE form then **totaldisrep** is identical to **ratdisrep**.

**totaldisrep** may be useful for ratdisrepping expressions such as equations, lists, matrices, etc., which have some subexpressions in CRE form.

**untellrat** (*x\_1, ..., x\_n*)

Function

Removes **tellrat** properties from *x\_1, ..., x\_n*.



## 13 Constants

### 13.1 Definitions for Constants

|                                                                         |          |
|-------------------------------------------------------------------------|----------|
| <b>%e</b>                                                               | Constant |
| - The base of natural logarithms, $e$ , is represented in Maxima as %e. |          |
| <b>false</b>                                                            | Constant |
| - the Boolean constant, false. (NIL in Lisp)                            |          |
| <b>%inf</b>                                                             | Constant |
| - real positive infinity.                                               |          |
| <b>%infinity</b>                                                        | Constant |
| - complex infinity.                                                     |          |
| <b>%minf</b>                                                            | Constant |
| - real minus infinity.                                                  |          |
| <b>pi</b>                                                               | Constant |
| - "pi" is represented in Maxima as %pi.                                 |          |
| <b>true</b>                                                             | Constant |
| - the Boolean constant, true. (T in Lisp)                               |          |



## 14 Logarithms

### 14.1 Definitions for Logarithms

**%e\_to\_numlog** option variable

Default value: `false`

When `true`,  $r$  some rational number, and  $x$  some expression,  $e^{(r \cdot \log(x))}$  will be simplified into  $x^r$ . It should be noted that the `radcan` command also does this transformation, and more complicated transformations of this ilk as well. The `logcontract` command "contracts" expressions containing `log`.

**log (x)** Function

Represents the natural logarithm of  $x$ .

Simplification and evaluation of logarithms is governed by several global flags:

`logexpand` - causes  $\log(a^b)$  to become  $b \cdot \log(a)$ . If it is set to `all`,  $\log(a \cdot b)$  will also simplify to  $\log(a) + \log(b)$ . If it is set to `super`, then  $\log(a/b)$  will also simplify to  $\log(a) - \log(b)$  for rational numbers  $a/b$ ,  $a \neq 1$ . ( $\log(1/b)$ , for  $b$  integer, always simplifies.) If it is set to `false`, all of these simplifications will be turned off.

`logsimp` - if `false` then no simplification of  $e$  to a power containing `log`'s is done.

`lognumber` - if `true` then negative floating point arguments to `log` will always be converted to their absolute value before the `log` is taken. If `number` is also `true`, then negative integer arguments to `log` will also be converted to their absolute value.

`lognegint` - if `true` implements the rule  $\log(-n) \rightarrow \log(n) + i \cdot \pi$  for  $n$  a positive integer.

`%e_to_numlog` - when `true`,  $r$  some rational number, and  $x$  some expression,  $e^{(r \cdot \log(x))}$  will be simplified into  $x^r$ . It should be noted that the `radcan` command also does this transformation, and more complicated transformations of this ilk as well. The `logcontract` command "contracts" expressions containing `log`.

**logabs** option variable

Default value: `false`

When doing indefinite integration where logs are generated, e.g. `integrate(1/x,x)`, the answer is given in terms of  $\log(\text{abs}(\dots))$  if `logabs` is `true`, but in terms of  $\log(\dots)$  if `logabs` is `false`. For definite integration, the `logabs:true` setting is used, because here "evaluation" of the indefinite integral at the endpoints is often needed.

**logarc** option variable

Default value: `false`

If `true` will cause the inverse circular and hyperbolic functions to be converted into logarithmic form. `logarc(exp)` will cause this conversion for a particular expression `exp` without setting the switch or having to re-evaluate the expression with `ev`.

**logconcoeffp** option variable

Default value: `false`

Controls which coefficients are contracted when using `logcontract`. It may be set to the name of a predicate function of one argument. E.g. if you like to generate SQRTs, you can do `logconcoeffp: 'logconfun$ logconfun(m):=featurep(m,integer) or ratnump(m)$`. Then `logcontract(1/2*log(x))`; will give `log(sqrt(x))`.

**logcontract** (*expr*) Function

Recursively scans the expression *expr*, transforming subexpressions of the form  $a_1 \log(b_1) + a_2 \log(b_2) + c$  into `log(ratsimp(b1^a1 * b2^a2)) + c`

(%i1) `2*(a*log(x) + 2*a*log(y))$`

(%i2) `logcontract(%);`

(%o2) 
$$a \log(x^2 y^4)$$

If you do `declare(n,integer)`; then `logcontract(2*a*n*log(x))`; gives `a*log(x^(2*n))`. The coefficients that "contract" in this manner are those such as the 2 and the n here which satisfy `featurep(coeff,integer)`. The user can control which coefficients are contracted by setting the option `logconcoeffp` to the name of a predicate function of one argument. E.g. if you like to generate SQRTs, you can do `logconcoeffp: 'logconfun$ logconfun(m):=featurep(m,integer) or ratnump(m)$`. Then `logcontract(1/2*log(x))`; will give `log(sqrt(x))`.

**logexpand** option variable

Default value: `true`

Causes `log(a^b)` to become `b*log(a)`. If it is set to `all`, `log(a*b)` will also simplify to `log(a)+log(b)`. If it is set to `super`, then `log(a/b)` will also simplify to `log(a)-log(b)` for rational numbers `a/b`, `a#1`. (`log(1/b)`, for integer `b`, always simplifies.) If it is set to `false`, all of these simplifications will be turned off.

**lognegint** option variable

Default value: `false`

If `true` implements the rule `log(-n) -> log(n)+%i*%pi` for `n` a positive integer.

**lognumer** option variable

Default value: `false`

If `true` then negative floating point arguments to `log` will always be converted to their absolute value before the `log` is taken. If `numer` is also `true`, then negative integer arguments to `log` will also be converted to their absolute value.

**logsimp** option variable

Default value: `true`

If `false` then no simplification of `%e` to a power containing `log`'s is done.

**plog** (*x*) Function

Represents the principal branch of the complex-valued natural logarithm with `-%pi < carg(x) <= +%pi`.

## 15 Trigonometric

### 15.1 Introduction to Trigonometric

Maxima has many trigonometric functions defined. Not all trigonometric identities are programmed, but it is possible for the user to add many of them using the pattern matching capabilities of the system. The trigonometric functions defined in Maxima are: `acos`, `acosh`, `acot`, `acoth`, `acsc`, `acsch`, `asec`, `asech`, `asin`, `asinh`, `atan`, `atanh`, `cos`, `cosh`, `cot`, `coth`, `csc`, `csch`, `sec`, `sech`, `sin`, `sinh`, `tan`, and `tanh`. There are a number of commands especially for handling trigonometric functions, see `trigexpand`, `trigreduce`, and the switch `trigsign`. Two share packages extend the simplification rules built into Maxima, `ntrig` and `atrig1`. Do `describe(command)` for details.

### 15.2 Definitions for Trigonometric

|                                                     |          |
|-----------------------------------------------------|----------|
| <b>acos</b> ( $x$ )<br>- Arc Cosine.                | Function |
| <b>acosh</b> ( $x$ )<br>- Hyperbolic Arc Cosine.    | Function |
| <b>acot</b> ( $x$ )<br>- Arc Cotangent.             | Function |
| <b>acoth</b> ( $x$ )<br>- Hyperbolic Arc Cotangent. | Function |
| <b>acsc</b> ( $x$ )<br>- Arc Cosecant.              | Function |
| <b>acsch</b> ( $x$ )<br>- Hyperbolic Arc Cosecant.  | Function |
| <b>asec</b> ( $x$ )<br>- Arc Secant.                | Function |
| <b>asech</b> ( $x$ )<br>- Hyperbolic Arc Secant.    | Function |
| <b>asin</b> ( $x$ )<br>- Arc Sine.                  | Function |
| <b>asinh</b> ( $x$ )<br>- Hyperbolic Arc Sine.      | Function |



|                                                                                                                                                                                                                                                                                                                                                                              |                 |
|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----------------|
| <b>atan</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                          | Function        |
| - Arc Tangent.                                                                                                                                                                                                                                                                                                                                                               |                 |
| <b>atan2</b> ( $y, x$ )                                                                                                                                                                                                                                                                                                                                                      | Function        |
| - yields the value of <b>atan</b> ( $y/x$ ) in the interval $-\pi$ to $\pi$ .                                                                                                                                                                                                                                                                                                |                 |
| <b>atanh</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                         | Function        |
| - Hyperbolic Arc Tangent.                                                                                                                                                                                                                                                                                                                                                    |                 |
| <b>atrig1</b>                                                                                                                                                                                                                                                                                                                                                                | Package         |
| The <b>atrig1</b> package contains several additional simplification rules for inverse trigonometric functions. Together with rules already known to Maxima, the following angles are fully implemented: $0$ , $\pi/6$ , $\pi/4$ , $\pi/3$ , and $\pi/2$ . Corresponding angles in the other three quadrants are also available. Do <code>load(atrig1)</code> ; to use them. |                 |
| <b>cos</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                           | Function        |
| - Cosine.                                                                                                                                                                                                                                                                                                                                                                    |                 |
| <b>cosh</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                          | Function        |
| - Hyperbolic Cosine.                                                                                                                                                                                                                                                                                                                                                         |                 |
| <b>cot</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                           | Function        |
| - Cotangent.                                                                                                                                                                                                                                                                                                                                                                 |                 |
| <b>coth</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                          | Function        |
| - Hyperbolic Cotangent.                                                                                                                                                                                                                                                                                                                                                      |                 |
| <b>csc</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                           | Function        |
| - Cosecant.                                                                                                                                                                                                                                                                                                                                                                  |                 |
| <b>csch</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                          | Function        |
| - Hyperbolic Cosecant.                                                                                                                                                                                                                                                                                                                                                       |                 |
| <b>halfangles</b>                                                                                                                                                                                                                                                                                                                                                            | option variable |
| Default value: <code>false</code>                                                                                                                                                                                                                                                                                                                                            |                 |
| When <b>halfangles</b> is <code>true</code> , half-angles are simplified away.                                                                                                                                                                                                                                                                                               |                 |
| <b>ntrig</b>                                                                                                                                                                                                                                                                                                                                                                 | Package         |
| The <b>ntrig</b> package contains a set of simplification rules that are used to simplify trigonometric function whose arguments are of the form $f(n\pi/10)$ where $f$ is any of the functions <b>sin</b> , <b>cos</b> , <b>tan</b> , <b>csc</b> , <b>sec</b> and <b>cot</b> .                                                                                              |                 |
| <b>sec</b> ( $x$ )                                                                                                                                                                                                                                                                                                                                                           | Function        |
| - Secant.                                                                                                                                                                                                                                                                                                                                                                    |                 |

|                                                   |          |
|---------------------------------------------------|----------|
| <b>sech</b> ( <i>x</i> )<br>- Hyperbolic Secant.  | Function |
| <b>sin</b> ( <i>x</i> )<br>- Sine.                | Function |
| <b>sinh</b> ( <i>x</i> )<br>- Hyperbolic Sine.    | Function |
| <b>tan</b> ( <i>x</i> )<br>- Tangent.             | Function |
| <b>tanh</b> ( <i>x</i> )<br>- Hyperbolic Tangent. | Function |

**trigexpand** (*expr*) Function

Expands trigonometric and hyperbolic functions of sums of angles and of multiple angles occurring in *expr*. For best results, *expr* should be expanded. To enhance user control of simplification, this function expands only one level at a time, expanding sums of angles or multiple angles. To obtain full expansion into sines and cosines immediately, set the switch `trigexpand: true`.

`trigexpand` is governed by the following global flags:

`trigexpand`

If `true` causes expansion of all expressions containing `sin`'s and `cos`'s occurring subsequently.

`halfangles`

If `true` causes half-angles to be simplified away.

`trigexpandplus`

Controls the "sum" rule for `trigexpand`, expansion of sums (e.g. `sin(x + y)`) will take place only if `trigexpandplus` is `true`.

`trigexpandtimes`

Controls the "product" rule for `trigexpand`, expansion of products (e.g. `sin(2 x)`) will take place only if `trigexpandtimes` is `true`.

Examples:

```
(%i1) x+sin(3*x)/sin(x),trigexpand=true,expand;
          2          2
(%o1)      - sin (x) + 3 cos (x) + x
(%i2) trigexpand(sin(10*x+y));
(%o2)      cos(10 x) sin(y) + sin(10 x) cos(y)
```

**trigexpandplus** option variable

Default value: `true`

`trigexpandplus` controls the "sum" rule for `trigexpand`. Thus, when the `trigexpand` command is used or the `trigexpand` switch set to `true`, expansion of sums (e.g. `sin(x+y)`) will take place only if `trigexpandplus` is `true`.

**trigexpandtimes**

option variable

Default value: true

`trigexpandtimes` controls the "product" rule for `trigexpand`. Thus, when the `trigexpand` command is used or the `trigexpand` switch set to `true`, expansion of products (e.g.  $\sin(2*x)$ ) will take place only if `trigexpandtimes` is `true`.

**triginverses**

option variable

Default value: all

`triginverses` controls the simplification of the composition of trigonometric and hyperbolic functions with their inverse functions.

If `all`, both e.g.  $\text{atan}(\tan(x))$  and  $\tan(\text{atan}(x))$  simplify to  $x$ .

If `true`, the  $\text{arcfun}(\text{fun}(x))$  simplification is turned off.

If `false`, both the  $\text{arcfun}(\text{fun}(x))$  and  $\text{fun}(\text{arcfun}(x))$  simplifications are turned off.

**trigreduce** (*expr*, *x*)

Function

**trigreduce** (*expr*)

Function

Combines products and powers of trigonometric and hyperbolic sin's and cos's of  $x$  into those of multiples of  $x$ . It also tries to eliminate these functions when they occur in denominators. If  $x$  is omitted then all variables in *expr* are used.

See also `poissimp`.

```
(%i1) trigreduce(-sin(x)^2+3*cos(x)^2+x);
              cos(2 x)   cos(2 x)   1       1
(%o1)         ----- + 3 (----- + -) + x - -
              2           2         2       2
```

The trigonometric simplification routines will use declared information in some simple cases. Declarations about variables are used as follows, e.g.

```
(%i1) declare(j, integer, e, even, o, odd)$
(%i2) sin(x + (e + 1/2)*%pi);
(%o2)                                     cos(x)
(%i3) sin(x + (o + 1/2)*%pi);
(%o3)                                     - cos(x)
```

**trigsign**

option variable

Default value: true

When `trigsign` is `true`, it permits simplification of negative arguments to trigonometric functions. E.g.,  $\sin(-x)$  will become  $-\sin(x)$  only if `trigsign` is `true`.

**trigsimp** (*expr*)

Function

Employs the identities  $\sin(x)^2 + \cos(x)^2 = 1$  and  $\cosh(x)^2 - \sinh(x)^2 = 1$  to simplify expressions containing `tan`, `sec`, etc., to `sin`, `cos`, `sinh`, `cosh` so that further simplification may be obtained by using `trigreduce` on the result.

`demo ("trgsmp.dem")` displays some examples of `trigsimp`.

See also `trigsum`.

**trigrat** (*expr*)

Function

Gives a canonical simplified quasilinear form of a trigonometrical expression; *expr* is a rational fraction of several **sin**, **cos** or **tan**, the arguments of them are linear forms in some variables (or kernels) and  $\%pi/n$  ( $n$  integer) with integer coefficients. The result is a simplified fraction with numerator and denominator linear in **sin** and **cos**. Thus **trigrat** linearize always when it is possible.

```
(%i1) trigrat(sin(3*a)/sin(a+%pi/3));
(%o1)          sqrt(3) sin(2 a) + cos(2 a) - 1
```

The following example is taken from Davenport, Siret, and Tournier, *Calcul Formel*, Masson (or in English, Addison-Wesley), section 1.5.5, Morley theorem.

```
(%i1) c: %pi/3 - a - b;
(%o1)          - b - a + ----
                    %pi
                    3
(%i2) bc: sin(a)*sin(3*c)/sin(a+b);
(%o2)          sin(a) sin(3 b + 3 a)
                    -----
                    sin(b + a)
(%i3) ba: bc, c=a, a=c$
(%i4) ac2: ba^2 + bc^2 - 2*bc*ba*cos(b);
(%o4)          sin (a) sin (3 b + 3 a)
                    -----
                    2
                    sin (b + a)

                    %pi
                    2 sin(a) sin(3 a) cos(b) sin(b + a - ----) sin(3 b + 3 a)
                    3
-----
                    %pi
                    sin(a - ----) sin(b + a)
                    3

                    2      2      %pi
                    sin (3 a) sin (b + a - ----)
                    3
+ -----
                    2      %pi
                    sin (a - ----)
                    3
(%i5) trigrat (ac2);
(%o5) - (sqrt(3) sin(4 b + 4 a) - cos(4 b + 4 a)
- 2 sqrt(3) sin(4 b + 2 a) + 2 cos(4 b + 2 a)
- 2 sqrt(3) sin(2 b + 4 a) + 2 cos(2 b + 4 a)
```

$$\begin{aligned} &+ 4 \sqrt{3} \sin(2 b + 2 a) - 8 \cos(2 b + 2 a) - 4 \cos(2 b - 2 a) \\ &+ \sqrt{3} \sin(4 b) - \cos(4 b) - 2 \sqrt{3} \sin(2 b) + 10 \cos(2 b) \\ &+ \sqrt{3} \sin(4 a) - \cos(4 a) - 2 \sqrt{3} \sin(2 a) + 10 \cos(2 a) \\ &- 9)/4 \end{aligned}$$

## 16 Special Functions

### 16.1 Introduction to Special Functions

### 16.2 `specint`

`hypgeo` is a package for handling Laplace transforms of special functions. `hyp` is a package for handling generalized Hypergeometric functions.

`specint` attempts to compute the definite integral (over the range from zero to infinity) of an expression containing special functions. When the integrand contains a factor `exp (-s t)`, the result is a Laplace transform.

The syntax is as follows:

```
specint (exp (-s*t) * expr, t);
```

where  $t$  is the variable of integration and  $expr$  is an expression containing special functions.

If `specint` cannot compute the integral, the return value may contain various Lisp symbols, including `other-defint-to-follow-negttest`, `other-lt-exponential-to-follow`, `product-of-y-with-nofract-indices`, etc.; this is a bug.

Special function notation follows:

|                                     |                                                       |
|-------------------------------------|-------------------------------------------------------|
| <code>bessel_j (index, expr)</code> | Bessel function, 1st kind                             |
| <code>bessel_y (index, expr)</code> | Bessel function, 2nd kind                             |
| <code>bessel_i (index, expr)</code> | Modified Bessel function, 1st kind                    |
| <code>bessel_k (index, expr)</code> | Modified Bessel function, 2nd kind                    |
| <code>%he[n] (z)</code>             | Hermite polynomial (Nota bene: he, not h. See A&S 22) |
| <code>%p[u,v] (z)</code>            | Legendre function                                     |
| <code>%q[u,v] (z)</code>            | Legendre function, 2nd kind                           |
| <code>hstruve[n] (z)</code>         | Struve H function                                     |
| <code>lstruve[n] (z)</code>         | Struve L function                                     |
| <code>%f[p,q] ([], [], expr)</code> | Generalized Hypergeometric function                   |
| <code>gamma()</code>                | Gamma function                                        |
| <code>gammagreek(a,z)</code>        | Incomplete gamma function                             |
| <code>gammaincomplete(a,z)</code>   | Tail of incomplete gamma function                     |
| <code>slommel</code>                |                                                       |
| <code>%m[u,k] (z)</code>            | Whittaker function, 1st kind                          |
| <code>%w[u,k] (z)</code>            | Whittaker function, 2nd kind                          |
| <code>erfc (z)</code>               | Complement of the erf function                        |
| <code>ei (z)</code>                 | Exponential integral (?)                              |
| <code>kelliptic (z)</code>          | Complete elliptic integral of the first kind (K)      |
| <code>%d [n] (z)</code>             | Parabolic cylinder function                           |

`demo ("hypgeo")` displays several examples of Laplace transforms computed by `specint`.

This is a work in progress. Some of the function names may change.

## 16.3 Definitions for Special Functions

**airy** ( $x$ ) Function

The Airy function  $Ai$ . If the argument  $x$  is a number, the numerical value of **airy** ( $x$ ) is returned. Otherwise, an unevaluated expression **airy** ( $x$ ) is returned.

The Airy equation  $\text{diff}(y(x), x, 2) - x y(x) = 0$  has two linearly independent solutions, named **ai** and **bi**. This equation is very popular as an approximation to more complicated problems in many mathematical physics settings.

`load("airy")` loads the functions **ai**, **bi**, **dai**, and **dbi**.

The **airy** package contains routines to compute **ai** and **bi** and their derivatives **dai** and **dbi**. The result is a floating point number if the argument is a number, and an unevaluated expression otherwise.

An error occurs if the argument is large enough to cause an overflow in the exponentials, or a loss of accuracy in **sin** or **cos**. This makes the range of validity about -2800 to  $10^{38}$  for **ai** and **dai**, and -2800 to 25 for **bi** and **dbi**.

These derivative rules are known to Maxima:

- $\text{diff}(\text{ai}(x), x)$  yields  $\text{dai}(x)$ ,
- $\text{diff}(\text{dai}(x), x)$  yields  $x \text{ai}(x)$ ,
- $\text{diff}(\text{bi}(x), x)$  yields  $\text{dbi}(x)$ ,
- $\text{diff}(\text{dbi}(x), x)$  yields  $x \text{bi}(x)$ .

Function values are computed from the convergent Taylor series for  $\text{abs}(x) < 3$ , and from the asymptotic expansions for  $x < -3$  or  $x > 3$  as needed. This results in only very minor numerical discrepancies at  $x = 3$  and  $x = -3$ . For details, see Abramowitz and Stegun, *Handbook of Mathematical Functions*, Section 10.4 and Table 10.11.

`ev(taylor(ai(x), x, 0, 9), infeval)` yields a floating point Taylor expansions of the function **ai**. A similar expression can be constructed for **bi**.

**asympa** Function

**asympa** is a package for asymptotic analysis. The package contains simplification functions for asymptotic analysis, including the “big O” and “little o” functions that are widely used in complexity analysis and numerical analysis.

`load("asympa")` loads this package.

**bessel** ( $z, a$ ) Function

The Bessel function of the first kind.

This function is deprecated. Write **bessel\_j** ( $z, a$ ) instead.

**bessel\_j** ( $v, z$ ) Function

The Bessel function of the first kind of order  $v$  and argument  $z$ .

**bessel\_j** computes the array **besselarray** such that  $\text{besselarray}[i] = \text{bessel\_j}[i + v - \text{int}(v)](z)$  for  $i$  from zero to  $\text{int}(v)$ .

**bessel\_j** is defined as

$$\sum_{k=0}^{\infty} \frac{(-1)^k \left(\frac{z}{2}\right)^{v+2k}}{k! \Gamma(v+k+1)}$$

although the infinite series is not used for computations.

**bessel\_y** (*v*, *z*) Function

The Bessel function of the second kind of order *v* and argument *z*.

**bessel\_y** computes the array **besselarray** such that **besselarray** [*i*] = **bessel\_y** [*i* + *v* - **int**(*v*)] (*z*) for *i* from zero to **int**(*v*).

**bessel\_y** is defined as

$$\frac{\cos(\pi v) J_v(z) - J_{-v}(z)}{\sin(\pi v)}$$

when *v* is not an integer. When *v* is an integer *n*, the limit as *v* approaches *n* is taken.

**bessel\_i** (*v*, *z*) Function

The modified Bessel function of the first kind of order *v* and argument *z*.

**bessel\_i** computes the array **besselarray** such that **besselarray** [*i*] = **bessel\_i** [*i* + *v* - **int**(*v*)] (*z*) for *i* from zero to **int**(*v*).

**bessel\_i** is defined as

$$\sum_{k=0}^{\infty} \frac{1}{k! \Gamma(v+k+1)} \left(\frac{z}{2}\right)^{v+2k}$$

although the infinite series is not used for computations.

**bessel\_k** (*v*, *z*) Function

The modified Bessel function of the second kind of order *v* and argument *z*.

**bessel\_k** computes the array **besselarray** such that **besselarray** [*i*] = **bessel\_k** [*i* + *v* - **int**(*v*)] (*z*) for *i* from zero to **int**(*v*).

**bessel\_k** is defined as

$$\frac{\pi \csc(\pi v) (I_{-v}(z) - I_v(z))}{2}$$

when *v* is not an integer. If *v* is an integer *n*, then the limit as *v* approaches *n* is taken.

**besselexpand** Variable

Default value: **false**

Controls expansion of the Bessel functions when the order is half of an odd integer. In this case, the Bessel functions can be expanded in terms of other elementary functions. When **besselexpand** is **true**, the Bessel function is expanded.



```
(%i1) besselexpand: false$
(%i2) bessel_j (3/2, z);

(%o2)
          3
      bessel_j(-, z)
          2

(%i3) besselexpand: true$
(%i4) bessel_j (3/2, z);

(%o4)
          2 z   sin(z)   cos(z)
      sqrt(---) (----- - -----)
          %pi      2       z
          z
```

- j0** ( $x$ ) Function  
 The Bessel function of the first kind of order 0.  
 This function is deprecated. Write `bessel_j (0, x)` instead.
- j1** ( $x$ ) Function  
 The Bessel function of the first kind of order 1.  
 This function is deprecated. Write `bessel_j (1, x)` instead.
- jn** ( $x, n$ ) Function  
 The Bessel function of the first kind of order  $n$ .  
 This function is deprecated. Write `bessel_j (n, x)` instead.
- i0** ( $x$ ) Function  
 The modified Bessel function of the first kind of order 0.  
 This function is deprecated. Write `bessel_i (0, x)` instead.
- i1** ( $x$ ) Function  
 The modified Bessel function of the first kind of order 1.  
 This function is deprecated. Write `bessel_i (1, x)` instead.
- beta** ( $x, y$ ) Function  
 The beta function, defined as  $\text{gamma}(x) \text{gamma}(y) / \text{gamma}(x + y)$ .
- gamma** ( $x$ ) Function  
 The gamma function.  
 See also `makegamma`.  
 The variable `gammalim` controls simplification of the gamma function.  
 The Euler-Mascheroni constant is `%gamma`.
- gammalim** Variable  
 Default value: 1000000  
`gammalim` controls simplification of the gamma function for integral and rational number arguments. If the absolute value of the argument is not greater than `gammalim`, then simplification will occur. Note that the `factlim` switch controls simplification of the result of `gamma` of an integer argument as well.

- intopois** (*a*) Function  
 Converts *a* into a Poisson encoding.
- makefact** (*expr*) Function  
 Transforms instances of binomial, gamma, and beta functions in *expr* into factorials.  
 See also **makegamma**.
- makegamma** (*expr*) Function  
 Transforms instances of binomial, factorial, and beta functions in *expr* into gamma functions.  
 See also **makefact**.
- numfactor** (*expr*) Function  
 Returns the numerical factor multiplying the expression *expr*, which should be a single term.  
**content** returns the greatest common divisor (gcd) of all terms in a sum.
- ```
(%i1) gamma (7/2);
(%o1)
          15 sqrt(%pi)
          -----
             8
(%i2) numfactor (%);
(%o2)
          15
          --
             8
```
- outofpois** (*a*) Function  
 Converts *a* from Poisson encoding to general representation. If *a* is not in Poisson form, **outofpois** carries out the conversion, i.e., the return value is **outofpois** (**intopois** (*a*)). This function is thus a canonical simplifier for sums of powers of sine and cosine terms of a particular type.
- poisdiff** (*a*, *b*) Function  
 Differentiates *a* with respect to *b*. *b* must occur only in the trig arguments or only in the coefficients.
- poisexpt** (*a*, *b*) Function  
 Functionally identical to **intopois** ( $a^b$ ). *b* must be a positive integer.
- poisint** (*a*, *b*) Function  
 Integrates in a similarly restricted sense (to **poisdiff**). Non-periodic terms in *b* are dropped if *b* is in the trig arguments.
- poislim** Variable  
 Default value: 5  
**poislim** determines the domain of the coefficients in the arguments of the trig functions. The initial value of 5 corresponds to the interval  $[-2^{(5-1)+1}, 2^{(5-1)}]$ , or  $[-15, 16]$ , but it can be set to  $[-2^{(n-1)+1}, 2^{(n-1)}]$ .

- poismap** (*series, sinfn, cosfn*) Function  
 will map the functions *sinfn* on the sine terms and *cosfn* on the cosine terms of the Poisson series given. *sinfn* and *cosfn* are functions of two arguments which are a coefficient and a trigonometric part of a term in series respectively.
- poisplus** (*a, b*) Function  
 Is functionally identical to `intopois (a + b)`.
- poissimp** (*a*) Function  
 Converts *a* into a Poisson series for *a* in general representation.
- poisson** special symbol  
 The symbol /P/ follows the line label of Poisson series expressions.
- poissubst** (*a, b, c*) Function  
 Substitutes *a* for *b* in *c*. *c* is a Poisson series.  
 (1) Where *B* is a variable *u, v, w, x, y, or z*, then *a* must be an expression linear in those variables (e.g., `6*u + 4*v`).  
 (2) Where *b* is other than those variables, then *a* must also be free of those variables, and furthermore, free of sines or cosines.  
**poissubst** (*a, b, c, d, n*) is a special type of substitution which operates on *a* and *b* as in type (1) above, but where *d* is a Poisson series, expands `cos(d)` and `sin(d)` to order *n* so as to provide the result of substituting *a + d* for *b* in *c*. The idea is that *d* is an expansion in terms of a small parameter. For example, `poissubst (u, v, cos(v), %e, 3)` yields `cos(u)*(1 - %e^2/2) - sin(u)*(%e - %e^3/6)`.
- poistimes** (*a, b*) Function  
 Is functionally identical to `intopois (a*b)`.
- poistrim** () Function  
 is a reserved function name which (if the user has defined it) gets applied during Poisson multiplication. It is a predicate function of 6 arguments which are the coefficients of the *u, v, ..., z* in a term. Terms for which `poistrim` is `true` (for the coefficients of that term) are eliminated during multiplication.
- printpois** (*a*) Function  
 Prints a Poisson series in a readable format. In common with `outofpois`, it will convert *a* into a Poisson encoding first, if necessary.
- psi** (*x*) Function  
**psi** [*n*](*x*) Function  
 The derivative of `log (gamma (x))`.  
 Maxima does not know how to compute a numerical value of `psi`. However, the function `bfpsi` in the `bfac` package can compute numerical values.

# 17 Orthogonal Polynomials

## 17.1 Introduction to Orthogonal Polynomials

The `specfun` package contains Maxima code for the evaluation of all orthogonal polynomials listed in Chapter 22 of Abramowitz and Stegun. These include Chebyshev, Laguerre, Hermite, Jacobi, Legendre, and ultraspherical (Gegenbauer) polynomials. Additionally, `specfun` contains code for spherical Bessel, spherical Hankel, and spherical harmonic functions. The `specfun` package is not part of Maxima proper; it is loaded at request of the user via `load` or automatically via the `autoload` system.

The following table lists each function in `specfun`, its Maxima name, restrictions on its arguments, and a reference to the algorithm `specfun` uses to evaluate it. With few exceptions, `specfun` follows the conventions of Abramowitz and Stegun. In all cases,  $m$  and  $n$  must be integers.

A&S refers to Abramowitz and Stegun, *Handbook of Mathematical Functions* (10th printing, December 1972), G&R to Gradshteyn and Ryzhik, *Table of Integrals, Series, and Products* (1980 corrected and enlarged edition), and Merzbacher to *Quantum Mechanics* (second edition, 1970).

<i>Function</i>	<i>Maxima Name</i>	<i>Restrictions</i>	<i>Reference(s)</i>
Chebyshev T	<code>chebyshev_t(n, x)</code>	$n > -1$	A&S 22.5.31
Chebyshev U	<code>chebyshev_u(n, x)</code>	$n > -1$	A&S 22.5.32
generalized Laguerre	<code>gen_laguerre(n,a,x)</code>	$n > -1$	A&S page 789
Laguerre	<code>laguerre(n,x)</code>	$n > -1$	A&S 22.5.67
Hermite	<code>hermite(n,x)</code>	$n > -1$	A&S 22.4.40, 22.5.41
Jacobi	<code>jacobi_p(n,a,b,x)</code>	$n > -1, a, b > -1$	A&S page 789
associated Legendre P	<code>assoc_legendre_p(n,m,x)</code>	$n > -1$	A&S 22.5.37, 8.6.6, 8.2.5
associated Legendre Q	<code>assoc_legendre_q(n,m,x)</code>	$n > -1, m > -1$	G & R 8.706
Legendre P	<code>legendre_p(n,m,x)</code>	$n > -1$	A&S 22.5.35
Legendre Q	<code>legendre_q(n,m,x)</code>	$n > -1$	A&S 8.6.19
spherical Hankel 1st	<code>spherical_hankel1(n, x)</code>	$n > -1$	A&S 10.1.36
spherical Hankel 2nd	<code>spherical_hankel2(n, x)</code>	$n > -1$	A&S 10.1.17
spherical Bessel J	<code>spherical_bessel_j(n,x)</code>	$n > -1$	A&S 10.1.8, 10.1.15
spherical Bessel Y	<code>spherical_bessel_y(n,x)</code>	$n > -1$	A&S 10.1.9, 10.1.15
spherical harmonic	<code>spherical_harmonic(n,m,x,y)</code>	$ m  \leq n$	Merzbacher 9.64
ultraspherical (Gegenbauer)	<code>ultraspherical(n,a,x)</code>	$n > -1$	A&S 22.5.27

The `specfun` package is primarily intended for symbolic computation. It is hoped that it gives accurate floating point results as well; however, no claims are made that the algorithms are well suited for numerical evaluation. Some effort, however, has been made to provide good numerical performance. When all arguments, except for the order, are floats (but

not bigfloats), many functions in `specfun` call a float modeddeclared version of the Jacobi function. This greatly speeds floating point evaluation of the orthogonal polynomials.

`specfun` handles most domain errors by returning an unevaluated function. No simplification rules (based on recursion relations) are defined for unevaluated functions. It is possible for an expression involving sums of unevaluated special functions to vanish, yet Maxima is unable to reduce it to zero.

`load ("specfun")` loads the `specfun` package. Alternatively, `setup_autoload` causes the package to be loaded when one of the `specfun` functions appears in an expression. `setup_autoload` may appear at the command line or in the `maxima-init.mac` file. See `setup_autoload`.

An example use of `specfun` is

```
(%i1) load ("specfun")$
(%i2) [hermite (0, x), hermite (1, x), hermite (2, x)];
(%o2)
          2
      [1, 2 x, - 2 (1 - 2 x )]
(%i3) diff (hermite (n, x), x);
(%o3)
          2 n hermite(n - 1, x)
```

Generally, compiled code runs faster than translated code; however, translated code may be better for program development.

Some functions (namely `jacobi_p`, `ultraspherical`, `chebyshev_t`, `chebyshev_u`, and `legendre_p`), return a series representation when the order is a symbolic integer. The series representation is not used by `specfun` for any computations, but it may be simplified by Maxima automatically, or it may be possible to use the series to evaluate the function through further manipulations. For example:

```
(%i1) load ("specfun")$
(%i2) legendre_p (n, x);
(%o2)
          legendre_p(n, x)
(%i3) ultraspherical (n, 3/2, 2);
(%o3)
          genfact(3, n, - 1) jacobi_p(n, 1, 1, 2)
          -----
          genfact(2, n, - 1)

(%i4) declare (n, integer)$
(%i5) legendre_p (n, x);
          n - 1
          ====
          \
(%o5) ( > binomial(n, i%) binomial(n, n - i%) (x - 1)
          /
          ====
          i% = 1

          i%          n          n n
          (x + 1)  + (x + 1)  + (x - 1) )/2
(%i6) ultraspherical (n, 3/2, 2);
          n - 1
          ====
          \          i%
```

```
(%o6) genfact(3, n, - 1) ( > 3 binomial(n + 1, i%)
/
====
i% = 1

binomial(n + 1, n - i%) + (n + 1) n 3 + n + 1)

/n(genfact(2, n, - 1) 2 )
```

The first and last terms of the sum are added outside the summation. Removing these two terms avoids Maxima bugs associated with  $0^0$  terms in a sum that should evaluate to 1, but evaluate to 0 in a Maxima summation. Because the sum index runs from 1 to  $n - 1$ , the lower sum index will exceed the upper sum index when  $n = 0$ ; setting `sumhack` to true provides a fix. For example:

```
(%i1) load ("specfun")$
(%i2) declare (n, integer)$
(%i3) e: legendre_p(n,x)$
(%i4) ev (e, sum, n=0);
Lower bound to sum: 1
is greater than the upper bound: - 1
-- an error. Quitting. To debug this try debugmode(true);
(%i5) ev (e, sum, n=0, sumhack=true);
(%o5) 1
```

Most functions in `specfun` have a `gradef` property; derivatives with respect to the order or other function parameters are undefined, and an attempt to compute such a derivative yields an error message.

The `specfun` package and its documentation were written by Barton Willis of the University of Nebraska at Kearney. It is released under the terms of the General Public License (GPL). Send bug reports and comments on this package to `willisb@unk.edu`. In your report, please include the Maxima version, as reported by `build_info()`, and the `specfun` version, as reported by `get ('specfun, 'version)`.

## 17.2 Definitions for Orthogonal Polynomials

**assoc\_legendre\_p** ( $n, m, x$ ) Function

Returns the associated Legendre function of the first kind for integers  $n > -1$  and  $m > -1$ . When  $|m| > n$  and  $n \geq 0$ , we have  $assoc\_legendre_p(n, m, x) = 0$ . Reference: A&S 22.5.37 page 779, A&S 8.6.6 (second equation) page 334, and A&S 8.2.5 page 333.

`load ("specfun")` loads this function.

See [\[assoc\\_legendre\\_q\]](#), page 141, [\[legendre\\_p\]](#), page 143, and [\[legendre\\_q\]](#), page 143.

**assoc\_legendre\_q** ( $n, m, x$ ) Function

Returns the associated Legendre function of the second kind for integers  $n > -1$  and  $m > -1$ .

Reference: Gradshteyn and Ryzhik 8.706 page 1000.

`load ("specfun")` loads this function.

See also [\[assoc\\_legendre\\_p\]](#), page 141, [\[legendre\\_p\]](#), page 143, and [\[legendre\\_q\]](#), page 143.

### **chebyshev\_t** ( $n, x$ )

Function

Returns the Chebyshev function of the first kind for integers  $n > -1$ .

Reference: A&S 22.5.31 page 778 and A&S 6.1.22 page 256.

`load ("specfun")` loads this function.

See also [\[chebyshev\\_u\]](#), page 142.

### **chebyshev\_u** ( $n, x$ )

Function

Returns the Chebyshev function of the second kind for integers  $n > -1$ .

Reference: A&S, 22.8.3 page 783 and A&S 6.1.22 page 256.

`load ("specfun")` loads this function.

See also [\[chebyshev\\_t\]](#), page 142.

### **gen\_laguerre** ( $n, a, x$ )

Function

Returns the generalized Laguerre polynomial for integers  $n > -1$ .

`load ("specfun")` loads this function.

Reference: table on page 789 in A&S.

### **hermite** ( $n, x$ )

Function

Returns the Hermite polynomial for integers  $n > -1$ .

`load ("specfun")` loads this function.

Reference: A&S 22.5.40 and 22.5.41, page 779.

### **jacobi\_p** ( $n, a, b, x$ )

Function

Returns the Jacobi polynomial for integers  $n > -1$  and  $a$  and  $b$  symbolic or  $a > -1$  and  $b > -1$ . (The Jacobi polynomials are actually defined for all  $a$  and  $b$ ; however, the Jacobi polynomial weight  $(1-x)^a(1+x)^b$  isn't integrable for  $a \leq -1$  or  $b \leq -1$ .)

When  $a, b$ , and  $x$  are floats (but not bfloats) `specfun` calls a special modeddeclared version of `jacobi_p`. For numerical values, the modeddeclared version is much faster than the other version. Many functions in `specfun` are computed as a special case of the Jacobi polynomials; they also enjoy the speed boost from the modeddeclared version of `jacobi`.

If  $n$  has been declared to be an integer, `jacobi_p(n, a, b, x)` returns a summation representation for the Jacobi function. Because Maxima simplifies  $0^0$  to 0 in a sum, two terms of the sum are added outside the summation.

`load ("specfun")` loads this function.

Reference: table on page 789 in A&S.

- laguerre** ( $n, x$ ) Function  
Returns the Laguerre polynomial for integers  $n > -1$ .  
Reference: A&S 22.5.16, page 778 and A&S page 789.  
`load ("specfun")` loads this function.  
See also [\[gen.laguerre\]](#), page 142.
- legendre\_p** ( $n, x$ ) Function  
Returns the Legendre polynomial of the first kind for integers  $n > -1$ .  
Reference: A&S 22.5.35 page 779.  
`load ("specfun")` loads this function.  
See [\[legendre\\_q\]](#), page 143.
- legendre\_q** ( $n, x$ ) Function  
Returns the Legendre polynomial of the first kind for integers  $n > -1$ .  
Reference: A&S 8.6.19 page 334.  
`load ("specfun")` loads this function.  
See also [\[legendre\\_p\]](#), page 143.
- spherical\_bessel\_j** ( $n, x$ ) Function  
Returns the spherical Bessel function of the first kind for integers  $n > -1$ .  
Reference: A&S 10.1.8 page 437 and A&S 10.1.15 page 439.  
`load ("specfun")` loads this function.  
See also [\[spherical\\_hankel1\]](#), page 143, [\[spherical\\_hankel2\]](#), page 143, and [\[spherical\\_bessel\\_y\]](#), page 143.
- spherical\_bessel\_y** ( $n, x$ ) Function  
Returns the spherical Bessel function of the second kind for integers  $n > -1$ .  
Reference: A&S 10.1.9 page 437 and 10.1.15 page 439.  
`load ("specfun")` loads this function.  
See also [\[spherical\\_hankel1\]](#), page 143, [\[spherical\\_hankel2\]](#), page 143, and [\[spherical\\_bessel\\_y\]](#), page 143.
- spherical\_hankel1** ( $n, x$ ) Function  
Returns the spherical hankel function of the first kind for integers  $n > -1$ .  
Reference: A&S 10.1.36 page 439.  
`load ("specfun")` loads this function.  
See also [\[spherical\\_hankel2\]](#), page 143, [\[spherical\\_bessel\\_j\]](#), page 143, and [\[spherical\\_bessel\\_y\]](#), page 143.



- spherical\_hankel2** ( $n, x$ ) Function  
Returns the spherical hankel function of the second kind for integers  $n > -1$ .  
Reference: A&S 10.1.17 page 439.  
`load ("specfun")` loads this function.  
See also [\[spherical\\_hankel1\]](#), page 143, [\[spherical\\_bessel\\_j\]](#), page 143, and [\[spherical\\_bessel\\_y\]](#), page 143.
- spherical\_harmonic** ( $n, m, x, y$ ) Function  
Returns the spherical harmonic function for integers  $n > -1$  and  $|m| \leq n$ .  
Reference: Merzbacher 9.64.  
`load ("specfun")` loads this function.  
See also [\[assoc\\_legendre\\_p\]](#), page 141.
- ultraspherical** ( $n, a, x$ ) Function  
Returns the ultraspherical polynomials for integers  $n > -1$ . The ultraspherical polynomials are also known as Gegenbauer polynomials.  
Reference: A&S 22.5.27  
`load ("specfun")` loads this function.  
See also [\[jacobi\\_p\]](#), page 142.

## 18 Elliptic Functions

### 18.1 Introduction to Elliptic Functions and Integrals

Maxima includes support for Jacobian elliptic functions and for complete and incomplete elliptic integrals. This includes symbolic manipulation of these functions and numerical evaluation as well. Definitions of these functions and many of their properties can be found in Abramowitz and Stegun, Chapter 16–17. As much as possible, we use the definitions and relationships given there.

In particular, all elliptic functions and integrals use the parameter  $m$  instead of the modulus  $k$  or the modular angle  $\alpha$ . This is one area where we differ from Abramowitz and Stegun who use the modular angle for the elliptic functions. The following relationships are true:

$$m = k^2$$

and

$$k = \sin \alpha$$

The elliptic functions and integrals are primarily intended to support symbolic computation. Therefore, most of derivatives of the functions and integrals are known. However, if floating-point values are given, a floating-point result is returned.

Support for most of the other properties of elliptic functions and integrals other than derivatives has not yet been written.

Some examples of elliptic functions:

```
(%i1) jacobi_sn (u, m);
(%o1) jacobi_sn(u, m)
(%i2) jacobi_sn (u, 1);
(%o2) tanh(u)
(%i3) jacobi_sn (u, 0);
(%o3) sin(u)
(%i4) diff (jacobi_sn (u, m), u);
(%o4) jacobi_cn(u, m) jacobi_dn(u, m)
(%i5) diff (jacobi_sn (u, m), m);
(%o5) jacobi_cn(u, m) jacobi_dn(u, m)

      elliptic_e(asin(jacobi_sn(u, m)), m)
(u - -----)/(2 m)
      1 - m

      2
      jacobi_cn (u, m) jacobi_sn(u, m)
+ -----
      2 (1 - m)
```

Some examples of elliptic integrals:

```
(%i1) elliptic_f (phi, m);
(%o1) elliptic_f(phi, m)
```

```

(%i2) elliptic_f (phi, 0);
(%o2) phi
(%i3) elliptic_f (phi, 1);
(%o3) log(tan(--- + ---))
          2      4
          phi      %pi
(%i4) elliptic_e (phi, 1);
(%o4) sin(phi)
(%i5) elliptic_e (phi, 0);
(%o5) phi
(%i6) elliptic_kc (1/2);
(%o6) elliptic_kc(-)
          1
          2
(%i7) makegamma (%);
(%o7) gamma (-)
          2 1
          4
          -----
          4 sqrt(%pi)
(%i8) diff (elliptic_f (phi, m), phi);
(%o8) -----
          2
          sqrt(1 - m sin (phi))
(%i9) diff (elliptic_f (phi, m), m);
(%o9) (-----)
          m
          elliptic_e(phi, m) - (1 - m) elliptic_f(phi, m)
          -----
          2
          cos(phi) sin(phi)
          - -----)/(2 (1 - m))
          2
          sqrt(1 - m sin (phi))

```

Support for elliptic functions and integrals was written by Raymond Toy. It is placed under the terms of the General Public License (GPL) that governs the distribution of Maxima.

## 18.2 Definitions for Elliptic Functions

<b>jacobi_sn</b> ( $u, m$ )	Function
The Jacobian elliptic function $sn(u, m)$ .	
<b>jacobi_cn</b> ( $u, m$ )	Function
The Jacobian elliptic function $cn(u, m)$ .	
<b>jacobi_dn</b> ( $u, m$ )	Function
The Jacobian elliptic function $dn(u, m)$ .	

<b>jacobi_ns</b> ( $u, m$ )	Function
The Jacobian elliptic function $ns(u, m) = 1/sn(u, m)$ .	
<b>jacobi_sc</b> ( $u, m$ )	Function
The Jacobian elliptic function $sc(u, m) = sn(u, m)/cn(u, m)$ .	
<b>jacobi_sd</b> ( $u, m$ )	Function
The Jacobian elliptic function $sd(u, m) = sn(u, m)/dn(u, m)$ .	
<b>jacobi_nc</b> ( $u, m$ )	Function
The Jacobian elliptic function $nc(u, m) = 1/cn(u, m)$ .	
<b>jacobi_cs</b> ( $u, m$ )	Function
The Jacobian elliptic function $cs(u, m) = cn(u, m)/sn(u, m)$ .	
<b>jacobi_cd</b> ( $u, m$ )	Function
The Jacobian elliptic function $cd(u, m) = cn(u, m)/dn(u, m)$ .	
<b>jacobi_nd</b> ( $u, m$ )	Function
The Jacobian elliptic function $nc(u, m) = 1/cn(u, m)$ .	
<b>jacobi_ds</b> ( $u, m$ )	Function
The Jacobian elliptic function $ds(u, m) = dn(u, m)/sn(u, m)$ .	
<b>jacobi_dc</b> ( $u, m$ )	Function
The Jacobian elliptic function $dc(u, m) = dn(u, m)/cn(u, m)$ .	
<b>inverse_jacobi_sn</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $sn(u, m)$ .	
<b>inverse_jacobi_cn</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $cn(u, m)$ .	
<b>inverse_jacobi_dn</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $dn(u, m)$ .	
<b>inverse_jacobi_ns</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $ns(u, m)$ .	
<b>inverse_jacobi_sc</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $sc(u, m)$ .	
<b>inverse_jacobi_sd</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $sd(u, m)$ .	

<b>inverse_jacobi_nc</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $nc(u, m)$ .	
<b>inverse_jacobi_cs</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $cs(u, m)$ .	
<b>inverse_jacobi_cd</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $cd(u, m)$ .	
<b>inverse_jacobi_nd</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $nc(u, m)$ .	
<b>inverse_jacobi_ds</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $ds(u, m)$ .	
<b>inverse_jacobi_dc</b> ( $u, m$ )	Function
The inverse of the Jacobian elliptic function $dc(u, m)$ .	

### 18.3 Definitions for Elliptic Integrals

<b>elliptic_f</b> ( $\phi, m$ )	Function
The incomplete elliptic integral of the first kind, defined as	

$$\int_0^\phi \frac{d\theta}{\sqrt{1 - m \sin^2 \theta}}$$

See also [\[elliptic-e\]](#), page 148 and [\[elliptic-ke\]](#), page 149.

<b>elliptic_e</b> ( $\phi, m$ )	Function
The incomplete elliptic integral of the second kind, defined as See also <a href="#">[elliptic-e]</a> , page 148 and <a href="#">[elliptic-ec]</a> , page 149.	

<b>elliptic_eu</b> ( $u, m$ )	Function
The incomplete elliptic integral of the second kind, defined as	

$$\int_0^u \operatorname{dn}(v, m) dv = \int_0^\tau \sqrt{\frac{1 - mt^2}{1 - t^2}} dt$$

where  $\tau = \operatorname{sn}(u, m)$

This is related to *elliptic<sub>e</sub>* by

$$E(u, m) = E(\phi, m)$$

where  $\phi = \sin^{-1} \operatorname{sn}(u, m)$ ,  $m$ ) See also [\[elliptic-e\]](#), page 148.

**elliptic\_pi** (*n*, *phi*, *m*)

Function

The incomplete elliptic integral of the third kind, defined as

$$\int_0^\phi \frac{d\theta}{(1 - n \sin^2 \theta) \sqrt{1 - m \sin^2 \theta}}$$

Only the derivative with respect to *phi* is known by Maxima.

**elliptic\_kc** (*m*)

Function

The complete elliptic integral of the first kind, defined as

$$\int_0^{\frac{\pi}{2}} \frac{d\theta}{\sqrt{1 - m \sin^2 \theta}}$$

For certain values of *m*, the value of the integral is known in terms of *Gamma* functions. Use `makegamma` to evaluate them.

**elliptic\_ec** (*m*)

Function

The complete elliptic integral of the second kind, defined as

$$\int_0^{\frac{\pi}{2}} \sqrt{1 - m \sin^2 \theta} d\theta$$

For certain values of *m*, the value of the integral is known in terms of *Gamma* functions. Use `makegamma` to evaluate them.



## 19 Limits

### 19.1 Definitions for Limits

**lhospitallim** option variable

Default: 4

**lhospitallim** is the maximum number of times L'Hospital's rule is used in **limit**. This prevents infinite looping in cases like **limit** (**cot(x)/csc(x)**, **x**, 0).

**limit** (*expr*, *x*, *val*, *dir*) Function

**limit** (*expr*, *x*, *val*) Function

**limit** (*expr*) Function

Computes the limit of *expr* as the real variable *x* approaches the value *val* from the direction *dir*. *dir* may have the value **plus** for a limit from above, **minus** for a limit from below, or may be omitted (implying a two-sided limit is to be computed).

**limit** uses the following special symbols: **inf** (positive infinity) and **minf** (negative infinity). On output it may also use **und** (undefined), **ind** (indefinite but bounded) and **infinity** (complex infinity).

**lhospitallim** is the maximum number of times L'Hospital's rule is used in **limit**. This prevents infinite looping in cases like **limit** (**cot(x)/csc(x)**, **x**, 0).

**tlimswitch** when true will cause the limit package to use Taylor series when possible.

**limsubst** prevents **limit** from attempting substitutions on unknown forms. This is to avoid bugs like **limit** (**f(n)/f(n+1)**, **n**, **inf**) giving 1. Setting **limsubst** to **true** will allow such substitutions.

**limit** with one argument is often called upon to simplify constant expressions, for example, **limit** (**inf-1**).

**example** (**limit**) displays some examples.

For the method see Wang, P., "Evaluation of Definite Integrals by Symbolic Manipulation", Ph.D. thesis, MAC TR-92, October 1971.

**limsubst** option variable

default value: **false** - prevents **limit** from attempting substitutions on unknown forms. This is to avoid bugs like **limit** (**f(n)/f(n+1)**, **n**, **inf**) giving 1. Setting **limsubst** to **true** will allow such substitutions.

**tlimit** (*expr*, *x*, *val*, *dir*) Function

**tlimit** (*expr*, *x*, *val*) Function

**tlimit** (*expr*) Function

Returns **limit** with **tlimswitch** set to **true**.

**tlimswitch** option variable

Default value: **false**

When **tlimswitch** is **true**, it causes the limit package to use Taylor series when possible.





## 20 Differentiation

### 20.1 Definitions for Differentiation

**antid** (*expr*, *x*, *u(x)*) Function

Returns a two-element list, such that an antiderivative of *expr* with respect to *x* can be constructed from the list. The expression *expr* may contain an unknown function *u* and its derivatives.

Let *L*, a list of two elements, be the return value of **antid**. Then *L*[1] + 'integrate (*L*[2], *x*) is an antiderivative of *expr* with respect to *x*.

When **antid** succeeds entirely, the second element of the return value is zero. Otherwise, the second element is nonzero, and the first element is nonzero or zero. If **antid** cannot make any progress, the first element is zero and the second nonzero.

`load ("antid")` loads this function. The **antid** package also defines the functions **nonzeroandfreeof** and **linear**.

**antid** is related to **antidiff** as follows. Let *L*, a list of two elements, be the return value of **antid**. Then the return value of **antidiff** is equal to *L*[1] + 'integrate (*L*[2], *x*) where *x* is the variable of integration.

Examples:

```
(%i1) load ("antid")$
(%i2) expr: exp (z(x)) * diff (z(x), x) * y(x);
(%o2)          y(x) %ez(x) (--- (z(x)))
                    dx
(%i3) a1: antid (expr, x, z(x));
(%o3)          [y(x) %ez(x), - %ez(x) (--- (y(x)))]
                    dx
(%i4) a2: antidiff (expr, x, z(x));
(%o4)          y(x) %ez(x) - I %ez(x) (--- (y(x))) dx
                    /
(%i5) a2 - (first (a1) + 'integrate (second (a1), x));
(%o5)          0
(%i6) antid (expr, x, y(x));
(%o6)          [0, y(x) %ez(x) (--- (z(x)))]
                    dx
(%i7) antidiff (expr, x, y(x));
(%o7)          [
                    I y(x) %ez(x) (--- (z(x))) dx
                    ]
                    /
```

**antidiff** (*expr*, *x*, *u(x)*) Function

Returns an antiderivative of *expr* with respect to *x*. The expression *expr* may contain an unknown function *u* and its derivatives.

When **antidiff** succeeds entirely, the resulting expression is free of integral signs (that is, free of the **integrate** noun). Otherwise, **antidiff** returns an expression which is partly or entirely within an integral sign. If **antidiff** cannot make any progress, the return value is entirely within an integral sign.

`load ("antid")` loads this function. The **antid** package also defines the functions **nonzeroandfreeof** and **linear**.

**antidiff** is related to **antid** as follows. Let *L*, a list of two elements, be the return value of **antid**. Then the return value of **antidiff** is equal to *L*[1] + 'integrate (*L*[2], *x*) where *x* is the variable of integration.

Examples:

```
(%i1) load ("antid")$
(%i2) expr: exp (z(x)) * diff (z(x), x) * y(x);
(%o2)          y(x) %e      z(x) d
              (--- (z(x)))
              dx
(%i3) a1: antid (expr, x, z(x));
(%o3)          [y(x) %e      z(x) d
              , - %e      z(x) d
              (--- (y(x)))
              dx
(%i4) a2: antidiff (expr, x, z(x));
(%o4)          /
              z(x) [ z(x) d
              y(x) %e - I %e (--- (y(x))) dx
              ]
              dx
              /
(%i5) a2 - (first (a1) + 'integrate (second (a1), x));
(%o5)          0
(%i6) antid (expr, x, y(x));
(%o6)          [0, y(x) %e      z(x) d
              (--- (z(x)))
              dx
(%i7) antidiff (expr, x, y(x));
(%o7)          /
              [
              I y(x) %e      z(x) d
              (--- (z(x))) dx
              ]
              dx
              /
```

**atomgrad** property  
**atomgrad** is the atomic gradient property of an expression. This property is assigned by **gradef**.

**atvalue** (*expr*, [*x\_1* = *a\_1*, ..., *x\_m* = *a\_m*], *c*)

Function

**atvalue** (*expr*, *x\_1* = *a\_1*, *c*)

Function

Assigns the value *c* to *expr* at the point  $x = a$ . Typically boundary values are established by this mechanism.

*expr* is a function evaluation,  $f(x_1, \dots, x_m)$ , or a derivative,  $\text{diff}(f(x_1, \dots, x_m), x_1, n_1, \dots, x_n, n_n)$  in which the function arguments explicitly appear.  $n_i$  is the order of differentiation with respect to  $x_i$ .

The point at which the *atvalue* is established is given by the list of equations [*x\_1* = *a\_1*, ..., *x\_m* = *a\_m*]. If there is a single variable *x\_1*, the sole equation may be given without enclosing it in a list.

**printprops** (*[f\_1, f\_2, ...]*, *atvalue*) displays the *atvalues* of the functions *f\_1*, *f\_2*, ... as specified by calls to *atvalue*. **printprops** (*f*, *atvalue*) displays the *atvalues* of one function *f*. **printprops** (*all*, *atvalue*) displays the *atvalues* of all functions for which *atvalues* are defined.

The symbols @1, @2, ... represent the variables *x\_1*, *x\_2*, ... when *atvalues* are displayed.

*atvalue* evaluates its arguments. *atvalue* returns *c*, the *atvalue*.

Examples:

```
(%i1) atvalue (f(x,y), [x = 0, y = 1], a^2);
                                2
(%o1)                                a
(%i2) atvalue ('diff (f(x,y), x), x = 0, 1 + y);
(%o2)                                @2 + 1
(%i3) printprops (all, atvalue);
                                !
                                d
                                --- (f(@1, @2))!      = @2 + 1
                                d@1                    !
                                !@1 = 0
                                2
                                f(0, 1) = a

(%o3)                                done
(%i4) diff (4*f(x,y)^2 - u(x,y)^2, x);
                                d
                                d
(%o4)  8 f(x, y) (--- (f(x, y))) - 2 u(x, y) (--- (u(x, y)))
                                dx                    dx
(%i5) at (% , [x = 0, y = 1]);
                                !
                                2
                                d
(%o5)  16 a - 2 u(0, 1) (--- (u(x, y)))!
                                dx                    !
                                !x = 0, y = 1
```

**cartan** - Function

The exterior calculus of differential forms is a basic tool of differential geometry developed by Elie Cartan and has important applications in the theory of partial differential equations. The **cartan** package implements the functions **ext\_diff** and **lie\_diff**, along with the operators  $\sim$  (wedge product) and  $|$  (contraction of a form with a vector.) Type **demo (tensor)** to see a brief description of these commands along with examples.

**cartan** was implemented by F.B. Estabrook and H.D. Wahlquist.

**del** (*x*) Function

**del** (*x*) represents the differential of the variable *x*.

**diff** returns an expression containing **del** if an independent variable is not specified. In this case, the return value is the so-called "total differential".

Examples:

```
(%i1) diff (log (x));
(%o1)
      del(x)
      -----
      x

(%i2) diff (exp (x*y));
(%o2)
      x y      x y
      x %e del(y) + y %e del(x)

(%i3) diff (x*y*z);
(%o3)
      x y del(z) + x z del(y) + y z del(x)
```

**delta** (*t*) Function

The Dirac Delta function.

Currently only **laplace** knows about the **delta** function.

Example:

```
(%i1) laplace (delta (t - a) * sin(b*t), t, s);
Is a positive, negative, or zero?

p;
(%o1)
      - a s
      sin(a b) %e
```

**dependencies** Variable

Default value: []

**dependencies** is the list of atoms which have functional dependencies, assigned by **depends** or **gradef**. The **dependencies** list is cumulative: each call to **depends** or **gradef** appends additional items.

See **depends** and **gradef**.

**depends** (*f<sub>1</sub>*, *x<sub>1</sub>*, ..., *f<sub>n</sub>*, *x<sub>n</sub>*) Function

Declares functional dependencies among variables for the purpose of computing derivatives. In the absence of declared dependence, **diff** (*f*, *x*) yields zero. If

`depends (f, x)` is declared, `diff (f, x)` yields a symbolic derivative (that is, a `diff` noun).

Each argument  $f_i$ ,  $x_i$ , etc., can be the name of a variable or array, or a list of names. Every element of  $f_i$  (perhaps just a single element) is declared to depend on every element of  $x_i$  (perhaps just a single element). If some  $f_i$  is the name of an array or contains the name of an array, all elements of the array depend on  $x_i$ .

`diff` recognizes indirect dependencies established by `depends` and applies the chain rule in these cases.

`remove (f, dependency)` removes all dependencies declared for  $f$ .

`depends` returns a list of the dependencies established. The dependencies are appended to the global variable `dependencies`. `depends` evaluates its arguments.

`diff` is the only Maxima command which recognizes dependencies established by `depends`. Other functions (`integrate`, `laplace`, etc.) only recognize dependencies explicitly represented by their arguments. For example, `integrate` does not recognize the dependence of  $f$  on  $x$  unless explicitly represented as `integrate (f(x), x)`.

```
(%i1) depends ([f, g], x);
(%o1) [f(x), g(x)]
(%i2) depends ([r, s], [u, v, w]);
(%o2) [r(u, v, w), s(u, v, w)]
(%i3) depends (u, t);
(%o3) [u(t)]
(%i4) dependencies;
(%o4) [f(x), g(x), r(u, v, w), s(u, v, w), u(t)]
(%i5) diff (r.s, u);
(%o5)
      dr      ds
      -- . s + r . --
      du      du

(%i6) diff (r.s, t);
(%o6)
      dr du      ds du
      -- -- . s + r . -- --
      du dt      du dt

(%i7) remove (r, dependency);
(%o7) done
(%i8) diff (r.s, t);
(%o8)
      ds du
      r . -- --
      du dt
```

### **derivabbrev**

Variable

Default value: `false`

When `derivabbrev` is `true`, symbolic derivatives (that is, `diff` nouns) are displayed as subscripts. Otherwise, derivatives are displayed in the Leibniz notation  $dy/dx$ .

### **derivdegree** ( $expr$ , $y$ , $x$ )

Function

Returns the highest degree of the derivative of the dependent variable  $y$  with respect to the independent variable  $x$  occurring in  $expr$ .

Example:

```
(%i1) 'diff (y, x, 2) + 'diff (y, z, 3) + 'diff (y, x) * x^2;
(%o1)

$$\frac{d^2 y}{dx^2} + \frac{d^3 y}{dz^3} + x \frac{dy}{dx}$$

(%i2) derivdegree (%o1, y, x);
(%o2) 2
```

**derivlist** (*var\_1*, ..., *var\_k*) Function  
 Causes only differentiations with respect to the indicated variables, within the `ev` command.

**derivsubst** Variable  
 Default value: `false`  
 When `derivsubst` is `true`, a non-syntactic substitution such as `subst (x, 'diff (y, t), 'diff (y, t, 2))` yields `'diff (x, t)`.

**diff** (*expr*, *x\_1*, *n\_1*, ..., *x\_m*, *n\_m*) Function  
**diff** (*expr*, *x*, *n*) Function  
**diff** (*expr*, *x*) Function  
**diff** (*expr*) Function

Returns the derivative or differential of *expr* with respect to some or all variables in *expr*.

`diff (expr, x, n)` returns the *n*'th derivative of *expr* with respect to *x*.

`diff (expr, x_1, n_1, ..., x_m, n_m)` returns the mixed partial derivative of *expr* with respect to *x\_1*, ..., *x\_m*. It is equivalent to `diff (... (diff (expr, x_m, n_m) ...), x_1, n_1)`.

`diff (expr, x)` returns the first derivative of *expr* with respect to the variable *x*.

`diff (expr)` returns the total differential of *expr*, that is, the sum of the derivatives of *expr* with respect to each its variables times the differential `del` of each variable. No further simplification of `del` is offered.

The noun form of `diff` is required in some contexts, such as stating a differential equation. In these cases, `diff` may be quoted (as `'diff`) to yield the noun form instead of carrying out the differentiation.

When `derivabbrev` is `true`, derivatives are displayed as subscripts. Otherwise, derivatives are displayed in the Leibniz notation, `dy/dx`.

Examples:

```
(%i1) diff (exp (f(x)), x, 2);
(%o1)

$$e^{f(x)} \frac{d^2 (f(x))}{dx^2} + e^{f(x)} \frac{d^2 (f(x))}{dx^2}$$

```

```

(%i2) derivabbrev: true$
(%i3) 'integrate (f(x, y), y, g(x), h(x));
      h(x)
      /
      [
(%o3) I      f(x, y) dy
      ]
      /
      g(x)

(%i4) diff (% , x);
      h(x)
      /
      [
(%o4) I      f(x, y) dy + f(x, h(x)) h(x) - f(x, g(x)) g(x)
      ]      x          x          x
      /
      g(x)

```

For the tensor package, the following modifications have been incorporated:

(1) The derivatives of any indexed objects in *expr* will have the variables  $x_i$  appended as additional arguments. Then all the derivative indices will be sorted.

(2) The  $x_i$  may be integers from 1 up to the value of the variable **dimension** [default value: 4]. This will cause the differentiation to be carried out with respect to the  $x_i$ 'th member of the list **coordinates** which should be set to a list of the names of the coordinates, e.g., [**x**, **y**, **z**, **t**]. If **coordinates** is bound to an atomic variable, then that variable subscripted by  $x_i$  will be used for the variable of differentiation. This permits an array of coordinate names or subscripted names like **X**[1], **X**[2], ... to be used. If **coordinates** has not been assigned a value, then the variables will be treated as in (1) above.

## **diff**

special symbol

When **diff** is present as an **evflag** in call to **ev**, all differentiations indicated in **expr** are carried out.

## **dscalar** (*f*)

Function

Applies the scalar d'Alembertian to the scalar function *f*.

**load** ("ctensor") loads this function.

## **express** (*expr*)

Function

Expands differential operator nouns into expressions in terms of partial derivatives. **express** recognizes the operators **grad**, **div**, **curl**, **laplacian**. **express** also expands the cross product  $\sim$ .

Symbolic derivatives (that is, **diff** nouns) in the return value of **express** may be evaluated by including **diff** in the **ev** function call or command line. In this context, **diff** acts as an **evfun**.

**load** ("vect") loads this function.

Examples:



```

(%i1) load ("vect")$
(%i2) grad (x^2 + y^2 + z^2);
(%o2)
      2      2      2
      grad (z  + y  + x )
(%i3) express (%);
      d      2      2      2      d      2      2      2      d      2      2      2
(%o3) [--- (z  + y  + x ), --- (z  + y  + x ), --- (z  + y  + x )]
      dx      dy      dz
(%i4) ev (% , diff);
(%o4)
      [2 x, 2 y, 2 z]
(%i5) div ([x^2, y^2, z^2]);
      2      2      2
(%o5)
      div [x , y , z ]
(%i6) express (%);
      d      2      d      2      d      2
(%o6)
      --- (z ) + --- (y ) + --- (x )
      dz      dy      dx
(%i7) ev (% , diff);
(%o7)
      2 z + 2 y + 2 x
(%i8) curl ([x^2, y^2, z^2]);
      2      2      2
(%o8)
      curl [x , y , z ]
(%i9) express (%);
      d      2      d      2      d      2      d      2      d      2      d      2
(%o9) [--- (z ) - --- (y ), --- (x ) - --- (z ), --- (y ) - --- (x )]
      dy      dz      dz      dx      dx      dy
(%i10) ev (% , diff);
(%o10)
      [0, 0, 0]
(%i11) laplacian (x^2 * y^2 * z^2);
      2      2      2
(%o11)
      laplacian (x y z )
(%i12) express (%);
      2      2      2      2      2      2
      d      2      2      2      d      2      2      2      d      2      2      2
(%o12) --- (x y z ) + --- (x y z ) + --- (x y z )
      dz      dy      dx
(%i13) ev (% , diff);
      2      2      2      2      2      2
(%o13)
      2 y z + 2 x z + 2 x y
(%i14) [a, b, c] ~ [x, y, z];
(%o14)
      [a, b, c] ~ [x, y, z]
(%i15) express (%);
(%o15)
      [b z - c y, c x - a z, a y - b x]

```

**gradef** ( $f(x_1, \dots, x_n), g_1, \dots, g_m$ )

Function

**gradef** ( $a, x, \text{expr}$ )

Function

Defines the partial derivatives (i.e., the components of the gradient) of the function  $f$  or variable  $a$ .

`gradef (f(x1, ..., xn), g1, ..., gm)` defines  $df/dx_i$  as  $g_i$ , where  $g_i$  is an expression;  $g_i$  may be a function call, but not the name of a function. The number of partial derivatives  $m$  may be less than the number of arguments  $n$ , in which case derivatives are defined with respect to  $x_1$  through  $x_m$  only.

`gradef (a, x, expr)` defines the derivative of variable  $a$  with respect to  $x$  as  $expr$ . This also establishes the dependence of  $a$  on  $x$  (via `depends (a, x)`).

The first argument  $f(x_1, \dots, x_n)$  or  $a$  is quoted, but the remaining arguments  $g_1, \dots, g_m$  are evaluated. `gradef` returns the function or variable for which the partial derivatives are defined.

`gradef` can redefine the derivatives of Maxima's built-in functions. For example, `gradef (sin(x), sqrt(1 - sin(x)^2))` redefines the derivative of `sin`.

`gradef` cannot define partial derivatives for a subscripted function.

`printprops ([f1, ..., fn], gradef)` displays the partial derivatives of the functions  $f_1, \dots, f_n$ , as defined by `gradef`.

`printprops ([an, ..., an], atomgrad)` displays the partial derivatives of the variables  $a_n, \dots, a_n$ , as defined by `gradef`.

`gradefs` is the list of the functions for which partial derivatives have been defined by `gradef`. `gradefs` does not include any variables for which partial derivatives have been defined by `gradef`.

Gradients are needed when, for example, a function is not known explicitly but its first derivatives are and it is desired to obtain higher order derivatives.

## **gradefs**

Variable

Default value: []

`gradefs` is the list of the functions for which partial derivatives have been defined by `gradef`. `gradefs` does not include any variables for which partial derivatives have been defined by `gradef`.

## **laplace (expr, t, s)**

Function

Attempts to compute the Laplace transform of  $expr$  with respect to the variable  $t$  and transform parameter  $s$ . If `laplace` cannot find a solution, a noun 'laplace is returned.

`laplace` recognizes in  $expr$  the functions `delta`, `exp`, `log`, `sin`, `cos`, `sinh`, `cosh`, and `erf`, as well as `derivative`, `integrate`, `sum`, and `ilt`. If some other functions are present, `laplace` may not be able to compute the transform.

$expr$  may also be a linear, constant coefficient differential equation in which case `atvalue` of the dependent variable is used. The required `atvalue` may be supplied either before or after the transform is computed. Since the initial conditions must be specified at zero, if one has boundary conditions imposed elsewhere he can impose these on the general solution and eliminate the constants by solving the general solution for them and substituting their values back.

`laplace` recognizes convolution integrals of the form `integrate (f(x) * g(t - x), x, 0, t)`; other kinds of convolutions are not recognized.

Functional relations must be explicitly represented in *expr*; implicit relations, established by *depends*, are not recognized. That is, if  $f$  depends on  $x$  and  $y$ ,  $f(x, y)$  must appear in *expr*.

See also *ilt*, the inverse Laplace transform.

Examples:

```
(%i1) laplace (exp (2*t + a) * sin(t) * t, t, s);
```

```
(%o1)
          a
          %e (2 s - 4)
-----
          2          2
          (s  - 4 s + 5)
```

```
(%i2) laplace ('diff (f (x), x), x, s);
```

```
(%o2) s laplace(f(x), x, s) - f(0)
```

```
(%i3) diff (diff (delta (t), t), t);
```

```
(%o3)
          2
          d
          --- (delta(t))
          2
          dt
```

```
(%i4) laplace (%o3, t, s);
```

```
(%o4)
          d          !          2
          -- (delta(t))!          + s  - delta(0) s
          dt          !
          !t = 0
```

## 21 Integration

### 21.1 Introduction to Integration

Maxima has several routines for handling integration. The `integrate` function makes use of most of them. There is also the `antid` package, which handles an unspecified function (and its derivatives, of course). For numerical uses, there is the `romberg` function; an adaptive integrator which uses the Newton-Cotes 8 panel quadrature rule, called `quanc8`; and a set of adaptive integrators from Quadpack, named `quad_qag`, `quad_qags`, etc. Hypergeometric functions are being worked on, see `specint` for details. Generally speaking, Maxima only handles integrals which are integrable in terms of the "elementary functions" (rational functions, trigonometrics, logs, exponentials, radicals, etc.) and a few extensions (error function, dilogarithm). It does not handle integrals in terms of unknown functions such as  $g(x)$  and  $h(x)$ .

### 21.2 Definitions for Integration

**changevar** (*expr*, *f(x,y)*, *y*, *x*)

Function

Makes the change of variable given by  $f(x,y) = 0$  in all integrals occurring in *expr* with integration with respect to *x*. The new variable is *y*.

```
(%i1) assume(a > 0)$
(%i2) 'integrate (%e**sqrt(a*y), y, 0, 4);
      4
      /
      [  sqrt(a) sqrt(y)
(%o2)  I  %e          dy
      ]
      /
      0
(%i3) changevar (%, y-z^2/a, z, y);
      0
      /
      [                abs(z)
      2 I                z %e          dz
      ]
      /
      - 2 sqrt(a)
(%o3)  -----
              a
```

An expression containing a noun form, such as the instances of `'integrate` above, may be evaluated by `ev` with the `nouns` flag. For example, the expression returned by `changevar` above may be evaluated by `ev (%o3, nouns)`.

`changevar` may also be used to changes in the indices of a sum or product. However, it must be realized that when a change is made in a sum or product, this change must be a shift, i.e.,  $i = j + \dots$ , not a higher degree function. E.g.,

```
(%i4) sum (a[i]*x^(i-2), i, 0, inf);
      inf
      ====
      \      i - 2
      >    a  x
      /      i
      ====
      i = 0
(%i5) changevar (%, i-2-n, n, i);
      inf
      ====
      \      n
      >    a  x
      /      n + 2
      ====
      n = - 2
```

**dblint** (*f, r, s, a, b*)

Function

A double-integral routine which was written in top-level Maxima and then translated and compiled to machine code. Use `load (dblint)` to access this package. It uses the Simpson's rule method in both the *x* and *y* directions to calculate

$$\int_a^b \int_{r(x)}^{s(x)} f(x,y) \, dy \, dx$$

The function *f* must be a translated or compiled function of two variables, and *r* and *s* must each be a translated or compiled function of one variable, while *a* and *b* must be floating point numbers. The routine has two global variables which determine the number of divisions of the *x* and *y* intervals: `dblint_x` and `dblint_y`, both of which are initially 10, and can be changed independently to other integer values (there are  $2*\text{dblint\_x}+1$  points computed in the *x* direction, and  $2*\text{dblint\_y}+1$  in the *y* direction). The routine subdivides the *X* axis and then for each value of *X* it first computes *r(x)* and *s(x)*; then the *Y* axis between *r(x)* and *s(x)* is subdivided and the integral along the *Y* axis is performed using Simpson's rule; then the integral along the *X* axis is done using Simpson's rule with the function values being the *Y*-integrals. This procedure may be numerically unstable for a great variety of reasons, but is reasonably fast: avoid using it on highly oscillatory functions and functions with singularities (poles or branch points in the region). The *Y* integrals depend on how far apart *r(x)* and *s(x)* are, so if the distance  $s(x) - r(x)$  varies rapidly with *X*, there may be substantial errors arising from truncation with different step-sizes in the various *Y* integrals. One can increase `dblint_x` and `dblint_y` in an effort to improve the coverage of the region, at the expense of computation time. The function values are not saved, so if the function is very time-consuming, you will have to wait for re-computation if you change anything (sorry). It is required that the functions *f*, *r*, and *s* be either translated or compiled prior to calling `dblint`. This will result in orders of magnitude speed improvement over interpreted code in many cases!

`demo (dblint)` executes a demonstration of `dblint` applied to an example problem.

**defint** (*expr*, *x*, *a*, *b*) Function

Attempts to compute a definite integral. `defint` is called by `integrate` when limits of integration are specified, i.e., when `integrate` is called as `integrate (expr, x, a, b)`. Thus from the user's point of view, it is sufficient to call `integrate`.

`defint` returns a symbolic expression, either the computed integral or the noun form of the integral. See `quad_qag` and related functions for numerical approximation of definite integrals.

**erf** (*x*) Function

Represents the error function, whose derivative is:  $2*\exp(-x^2)/\sqrt{\pi}$ .

**erfflag** Variable

Default value: `true`

When `erfflag` is `false`, prevents `risch` from introducing the `erf` function in the answer if there were none in the integrand to begin with.

**ilt** (*expr*, *t*, *s*) Function

Computes the inverse Laplace transform of *expr* with respect to *t* and parameter *s*. *expr* must be a ratio of polynomials whose denominator has only linear and quadratic factors. By using the functions `laplace` and `ilt` together with the `solve` or `linsolve` functions the user can solve a single differential or convolution integral equation or a set of them.

```
(%i1) 'integrate (sinh(a*x)*f(t-x), x, 0, t) + b*f(t) = t**2;
```

```

      t
      /
      [
(%o1)  I  f(t - x) sinh(a x) dx + b f(t) = t
      ]
      /
      0

```

```
(%i2) laplace (% , t, s);
```

```
(%o2)  b laplace(f(t), t, s) +  $\frac{a \text{laplace}(f(t), t, s) - 2}{s^2 - a^2} = \frac{2}{s^3}$ 
```

```
(%i3) linsolve ([%], ['laplace(f(t), t, s)]);
```

```
(%o3)  [laplace(f(t), t, s) =  $\frac{2 s^2 - 2 a^2}{5 b s^5 + (a - a^2 b) s^3}$ ]
```

```
(%i4) ilt (rhs (first (%)), s, t);
```

```
Is a b (a b - 1) positive, negative, or zero?
```

```
pos;
```

$$\begin{aligned}
 & \frac{\sqrt{a b (a b - 1)} t}{2 \cosh\left(\frac{b}{a t}\right)} \\
 (\%o4) & - \frac{a^3 b^2 - 2 a^2 b + a}{a^3 b^2 - 2 a^2 b + a} + \frac{a t}{a b - 1} \\
 & + \frac{2}{a^3 b^2 - 2 a^2 b + a}
 \end{aligned}$$

**integrate** (*expr*, *x*)

Function

**integrate** (*expr*, *x*, *a*, *b*)

Function

Attempts to symbolically compute the integral of *expr* with respect to *x*. **integrate** (*expr*, *x*) is an indefinite integral, while **integrate** (*expr*, *x*, *a*, *b*) is a definite integral, with limits of integration *a* and *b*. The limits should not contain *x*, although **integrate** does not enforce this restriction. *a* need not be less than *b*. If *b* is equal to *a*, **integrate** returns zero.

See **quad\_qag** and related functions for numerical approximation of definite integrals. See **residue** for computation of residues (complex integration). See **antid** for an alternative means of computing indefinite integrals.

The integral (an expression free of **integrate**) is returned if **integrate** succeeds. Otherwise the return value is the noun form of the integral (the quoted operator '**integrate**') or an expression containing one or more noun forms. The noun form of **integrate** is displayed with an integral sign.

In some circumstances it is useful to construct a noun form by hand, by quoting **integrate** with a single quote, e.g., '**integrate** (*expr*, *x*)'. For example, the integral may depend on some parameters which are not yet computed. The noun may be applied to its arguments by **ev** (*i*, **nouns**) where *i* is the noun form of interest.

**integrate** handles definite integrals separately from indefinite, and employs a range of heuristics to handle each case. Special cases of definite integrals include limits of integration equal to zero or infinity (**inf** or **minf**), trigonometric functions with limits of integration equal to zero and **%pi** or **2 %pi**, rational functions, integrals related to the definitions of the **beta** and **psi** functions, and some logarithmic and trigonometric integrals. Processing rational functions may include computation of residues. If an applicable special case is not found, an attempt will be made to compute the indefinite integral and evaluate it at the limits of integration. This may include taking a limit as a limit of integration goes to infinity or negative infinity; see also **ldefint**.

Special cases of indefinite integrals include trigonometric functions, exponential and logarithmic functions, and rational functions. **integrate** may also make use of a short table of elementary integrals.

**integrate** may carry out a change of variable if the integrand has the form **f(g(x)) \* diff(g(x), x)**. **integrate** attempts to find a subexpression **g(x)** such that the derivative of **g(x)** divides the integrand. This search may make use of derivatives defined by the **gradef** function. See also **changevar** and **antid**.

If none of the preceding heuristics find the indefinite integral, the Risch algorithm is executed. The flag `risch` may be set as an `evflag`, in a call to `ev` or on the command line, e.g., `ev (integrate (expr, x), risch)` or `integrate (expr, x), risch`. If `risch` is present, `integrate` calls the `risch` function without attempting heuristics first. See also `risch`.

`integrate` works only with functional relations represented explicitly with the `f(x)` notation. `integrate` does not respect implicit dependencies established by the `depends` function.

`integrate` may need to know some property of a parameter in the integrand. `integrate` will first consult the `assume` database, and, if the variable of interest is not there, `integrate` will ask the user. Depending on the question, suitable responses are `yes;` or `no;`, or `pos;`, `zero;`, or `neg;`.

`integrate` is not, by default, declared to be linear. See `declare` and `linear`.

`integrate` attempts integration by parts only in a few special cases.

Examples:

- Elementary indefinite and definite integrals.

```
(%i1) integrate (sin(x)^3, x);
```

```
(%o1)          3
              cos (x)
          ----- - cos(x)
              3
```

```
(%i2) integrate (x/ sqrt (b^2 - x^2), x);
```

```
(%o2)          2      2
              - sqrt(b  - x )
```

```
(%i3) integrate (cos(x)^2 * exp(x), x, 0, %pi);
```

```
(%o3)          %pi
              3 %e      3
          ----- - -
              5      5
```

```
(%i4) integrate (x^2 * exp(-x^2), x, minf, inf);
```

```
(%o4)          sqrt(%pi)
          -----
              2
```

- Use of `assume` and interactive query.

```
(%i1) assume (a > 1)$
```

```
(%i2) integrate (x**a/(x+1)**(5/2), x, 0, inf);
```

```
          2 a + 2
Is ----- an integer?
          5
```

```
no;
```

```
Is 2 a - 3 positive, negative, or zero?
```

```
neg;
```

```
(%o2)          3
          beta(a + 1, - - a)
              2
```



- Change of variable. There are two changes of variable in this example: one using a derivative established by `gradef`, and one using the derivation `diff(r(x))` of an unspecified function `r(x)`.

```
(%i3) gradef (q(x), sin(x**2));
(%o3)
      q(x)
(%i4) diff (log (q (r (x))), x);
      d
      2
      (-- (r(x))) sin(r (x))
      dx
(%o4) -----
      q(r(x))
(%i5) integrate (% , x);
(%o5) log(q(r(x)))
```

- Return value contains the 'integrate' noun form. In this example, Maxima can extract one factor of the denominator of a rational function, but cannot factor the remainder or otherwise find its integral. `grind` shows the noun form 'integrate' in the result. See also `integrate_use_rootsof` for more on integrals of rational functions.

```
(%i1) expand ((x-4) * (x^3+2*x+1));
      4      3      2
(%o1) x - 4 x + 2 x - 7 x - 4
(%i2) integrate (1/%, x);
      / 2
      [ x + 4 x + 18
      I ----- dx
      ] 3
      log(x - 4) / x + 2 x + 1
(%o2) ----- - -----
      73          73
(%i3) grind (%);
log(x-4)/73-(integrate((x^2+4*x+18)/(x^3+2*x+1),x))/73$
```

- Defining a function in terms of an integral. The body of a function is not evaluated when the function is defined. Thus the body of `f_1` in this example contains the noun form of `integrate`. The double-single-quotes operator `''` causes the integral to be evaluated, and the result becomes the body of `f_2`.

```
(%i1) f_1 (a) := integrate (x^3, x, 1, a);
      3
(%o1) f_1(a) := integrate(x , x, 1, a)
(%i2) ev (f_1 (7), nouns);
(%o2) 600
(%i3) /* Note parentheses around integrate(...) here */
      f_2 (a) := ''(integrate (x^3, x, 1, a));
      4
      a 1
(%o3) f_2(a) := -- - -
      4 4
(%i4) f_2 (7);
(%o4) 600
```

**integration\_constant\_counter**

Variable

Default value: 0

`integration_constant_counter` is a counter which is updated each time a constant of integration (named by Maxima, e.g., `integrationconstant1`) is introduced into an expression by indefinite integration of an equation.

**integrate\_use\_rootsof**

Variable

Default value: `false`

When `integrate_use_rootsof` is `true` and the denominator of a rational function cannot be factored, `integrate` returns the integral in a form which is a sum over the roots (not yet known) of the denominator.

For example, with `integrate_use_rootsof` set to `false`, `integrate` returns an unsolved integral of a rational function in noun form:

```
(%i1) integrate_use_rootsof: false$
(%i2) integrate (1/(1+x+x^5), x);
      / 2
      [ x  - 4 x + 5
      I ----- dx
      ] 3    2
      / x  - x  + 1
(%o2) ----- - ----- + -----
              7              14              7 sqrt(3)
```

Now we set the flag to be `true` and the unsolved part of the integral will be expressed as a summation over the roots of the denominator of the rational function:

```
(%i3) integrate_use_rootsof: true$
(%i4) integrate (1/(1+x+x^5), x);
====
\      2
\      (%r4  - 4 %r4 + 5) log(x - %r4)
> -----
/
====
      2
      3 %r4  - 2 %r4
      3    2
      %r4 in rootsof(x  - x  + 1)
(%o4) -----
              7
      2 x + 1
      2      5 atan(-----)
      log(x  + x + 1)  sqrt(3)
      ----- + -----
              14              7 sqrt(3)
```

Alternatively the user may compute the roots of the denominator separately, and then express the integrand in terms of these roots, e.g.,  $1/((x - a)*(x - b)*(x - c))$  or  $1/((x^2 - (a+b)*x + a*b)*(x - c))$  if the denominator is a cubic polynomial. Sometimes this will help Maxima obtain a more useful result.

**ldefint** (*expr*, *x*, *a*, *b*) Function

Attempts to compute the definite integral of *expr* by using `limit` to evaluate the indefinite integral of *expr* with respect to *x* at the upper limit *b* and at the lower limit *a*. If it fails to compute the definite integral, `ldefint` returns an expression containing limits as noun forms.

`ldefint` is not called from `integrate`, so executing `ldefint (expr, x, a, b)` may yield a different result than `integrate (expr, x, a, b)`. `ldefint` always uses the same method to evaluate the definite integral, while `integrate` may employ various heuristics and may recognize some special cases.

**potential** (*givengradient*) Function

The calculation makes use of the global variable `potentialzeroloc[0]` which must be `nonlist` or of the form

`[indeterminatej=expressionj, indeterminatek=expressionk, ...]`

the former being equivalent to the `nonlist` expression for all right-hand sides in the latter. The indicated right-hand sides are used as the lower limit of integration. The success of the integrations may depend upon their values and order. `potentialzeroloc` is initially set to 0.

**qq** Function

The package `qq` (which may be loaded with `load ("qq")`) contains a function `quanc8` which can take either 3 or 4 arguments. The 3 arg version computes the integral of the function specified as the first argument over the interval from `lo` to `hi` as in `quanc8 ('function, lo, hi)`. The function name should be quoted. The 4 arg version will compute the integral of the function or expression (first arg) with respect to the variable (second arg) over the interval from `lo` to `hi` as in `quanc8(<f(x) or expression in x>, x, lo, hi)`. The method used is the Newton-Cotes 8th order polynomial quadrature, and the routine is adaptive. It will thus spend time dividing the interval only when necessary to achieve the error conditions specified by the global variables `quanc8_relerr` (default value=1.0e-4) and `quanc8_abserr` (default value=1.0e-8) which give the relative error test:

`|integral(function) - computed value| < quanc8_relerr*|integral(function)|`

and the absolute error test:

`|integral(function) - computed value| < quanc8_abserr`

`printfile ("qq.usg")` yields additional information.

**quanc8** (*expr*, *a*, *b*) Function

An adaptive integrator. Demonstration and usage files are provided. The method is to use Newton-Cotes 8-panel quadrature rule, hence the function name `quanc8`, available in 3 or 4 arg versions. Absolute and relative error checks are used. To use it do `load ("qq")`. See also `qq`.

**residue** (*expr*, *z*, *z\_0*) Function

Computes the residue in the complex plane of the expression *expr* when the variable *z* assumes the value *z\_0*. The residue is the coefficient of  $(z - z_0)^{-1}$  in the Laurent series for *expr*.

```
(%i1) residue (s/(s**2+a**2), s, a*%i);
(%o1)
1
-
2
(%i2) residue (sin(a*x)/x**4, x, 0);
(%o2)
3
a
- --
6
```

**risch** (*expr*, *x*)

Function

Integrates *expr* with respect to *x* using the transcendental case of the Risch algorithm. (The algebraic case of the Risch algorithm has not been implemented.) This currently handles the cases of nested exponentials and logarithms which the main part of `integrate` can't do. `integrate` will automatically apply `risch` if given these cases.

`erfflag`, if `false`, prevents `risch` from introducing the `erf` function in the answer if there were none in the integrand to begin with.

```
(%i1) risch (x^2*erf(x), x);
(%o1)
3 2
%pi x erf(x) + (sqrt(%pi) x + sqrt(%pi)) %e - x
-----
3 %pi
(%i2) diff(%, x), ratsimp;
(%o2)
2
x erf(x)
```

**romberg** (*expr*, *x*, *a*, *b*)

Function

**romberg** (*expr*, *a*, *b*)

Function

Romberg integration. There are two ways to use this function. The first is an inefficient way like the definite integral version of `integrate`: `romberg (<integrand>, <variable of integration>, <lower limit>, <upper limit>)`.

Examples:

```
(%i1) showtime: true$
(%i2) romberg (sin(y), y, 0, %pi);
Evaluation took 0.00 seconds (0.01 elapsed) using 25.293 KB.
(%o2)
2.000000016288042
(%i3) 1/((x-1)^2+1/100) + 1/((x-2)^2+1/1000) + 1/((x-3)^2+1/200)$
(%i4) f(x) := ''%$
(%i5) rombergtol: 1e-6$
(%i6) rombergit: 15$
(%i7) romberg (f(x), x, -5, 5);
Evaluation took 11.97 seconds (12.21 elapsed) using 12.423 MB.
(%o7)
173.6730736617464
```

The second is an efficient way that is used as follows:

```
romberg (<function name>, <lower limit>, <upper limit>);
```

Continuing the above example, we have:

```
(%i8) f(x) := (mode_declare ([function(f), x], float), ''(%th(5)))$
(%i9) translate(f);
(%o9) [f]
(%i10) romberg (f, -5, 5);
Evaluation took 3.51 seconds (3.86 elapsed) using 6.641 MB.
(%o10) 173.6730736617464
```

The first argument must be a translated or compiled function. (If it is compiled it must be declared to return a `flonum`.) If the first argument is not already translated, `romberg` will not attempt to translate it but will give an error.

The accuracy of the integration is governed by the global variables `rombertol` (default value 1.E-4) and `rombergit` (default value 11). `romberg` will return a result if the relative difference in successive approximations is less than `rombertol`. It will try halving the stepsize `rombergit` times before it gives up. The number of iterations and function evaluations which `romberg` will do is governed by `rombergabs` and `rombergmin`.

`romberg` may be called recursively and thus can do double and triple integrals.

Example:

```
(%i1) assume (x > 0)$
(%i2) integrate (integrate (x*y/(x+y), y, 0, x/2), x, 1, 3)$
(%i3) radcan (%);
(%o3)
          26 log(3) - 26 log(2) - 13
          - -----
                   3
(%i4) %,numer;
(%o4) .8193023963959073
(%i5) define_variable (x, 0.0, float, "Global variable in function F")$
(%i6) f(y) := (mode_declare (y, float), x*y/(x+y))$
(%i7) g(x) := romberg ('f, 0, x/2)$
(%i8) romberg (g, 1, 3);
(%o8) .8193022864324522
```

The advantage with this way is that the function `f` can be used for other purposes, like plotting. The disadvantage is that you have to think up a name for both the function `f` and its free variable `x`. Or, without the global:

```
(%i1) g_1(x) := (mode_declare (x, float), romberg (x*y/(x+y), y, 0, x/2))$
(%i2) romberg (g_1, 1, 3);
(%o2) .8193022864324522
```

The advantage here is shortness.

```
(%i3) q (a, b) := romberg (romberg (x*y/(x+y), y, 0, x/2), x, a, b)$
(%i4) q (1, 3);
(%o4) .8193022864324522
```

It is even shorter this way, and the variables do not need to be declared because they are in the context of `romberg`. Use of `romberg` for multiple integrals can have great disadvantages, though. The amount of extra calculation needed because of

the geometric information thrown away by expressing multiple integrals this way can be incredible. The user should be sure to understand and use the `rombergtol` and `rombergit` switches.

### **rombergabs**

Variable

Default value: 0.0

Assuming that successive estimates produced by `romberg` are `y[0]`, `y[1]`, `y[2]`, etc., then `romberg` will return after `n` iterations if (roughly speaking)

```
(abs(y[n]-y[n-1]) <= rombergabs or
abs(y[n]-y[n-1])/(if y[n]=0.0 then 1.0 else y[n]) <= rombergtol)
```

is true. (The condition on the number of iterations given by `rombergmin` must also be satisfied.) Thus if `rombergabs` is 0.0 (the default) you just get the relative error test. The usefulness of the additional variable comes when you want to perform an integral, where the dominant contribution comes from a small region. Then you can do the integral over the small dominant region first, using the relative accuracy check, followed by the integral over the rest of the region using the absolute accuracy check.

Example: Suppose you want to compute

```
'integrate (exp(-x), x, 0, 50)
```

(numerically) with a relative accuracy of 1 part in 10000000. Define the function. `n` is a counter, so we can see how many function evaluations were needed. First of all try doing the whole integral at once.

```
(%i1) f(x) := (mode_declare (n, integer, x, float), n:n+1, exp(-x))$
(%i2) translate(f)$
Warning-> n is an undefined global variable.
(%i3) block ([rombergtol: 1.e-6, romberabs: 0.0], n:0, romberg (f, 0, 50));
(%o3) 1.000000000488271
(%i4) n;
(%o4) 257
```

That approach required 257 function evaluations. Now do the integral intelligently, by first doing `'integrate (exp(-x), x, 0, 10)` and then setting `rombergabs` to 1.E-6 times (this partial integral). This approach takes only 130 function evaluations.

```
(%i5) block ([rombergtol: 1.e-6, rombergabs:0.0, sum:0.0],
n: 0, sum: romberg (f, 0, 10), rombergabs: sum*rombergtol, rombergtol:0.0,
sum + romberg (f, 10, 50));
(%o5) 1.0000000001234793
(%i6) n;
(%o6) 130
```

So if `f(x)` were a function that took a long time to compute, the second method would be about 2 times quicker.

### **rombergit**

Variable

Default value: 11

The accuracy of the `romberg` integration command is governed by the global variables `rombergtol` and `rombergit`. `romberg` will return a result if the relative difference in successive approximations is less than `rombergtol`. It will try halving the stepsize `rombergit` times before it gives up.

**rombergmin**

Variable

Default value: 0

`rombergmin` governs the minimum number of function evaluations that `romberg` will make. `romberg` will evaluate its first arg. at least  $2^{(\text{rombergmin}+2)+1}$  times. This is useful for integrating oscillatory functions, when the normal converge test might sometimes wrongly pass.

**rombergtol**

Variable

Default value: 1e-4

The accuracy of the `romberg` integration command is governed by the global variables `rombergtol` and `rombergit`. `romberg` will return a result if the relative difference in successive approximations is less than `rombergtol`. It will try halving the stepsize `rombergit` times before it gives up.

**tldfint** (*expr*, *x*, *a*, *b*)

Function

Equivalent to `ldfint` with `tlimswitch` set to `true`.**quad\_qag** (*f(x)*, *x*, *a*, *b*, *key*, *epsrel*, *limit*)

Function

Numerically evaluate the integral

$$\int_a^b f(x)dx$$

using a simple adaptive integrator.

The function to be integrated is  $f(x)$ , with dependent variable  $x$ , and the function is to be integrated between the limits  $a$  and  $b$ . *key* is the integrator to be used and should be an integer between 1 and 6, inclusive. The value of *key* selects the order of the Gauss-Kronrod integration rule.

The numerical integration is done adaptively by subdividing the integration region into sub-intervals until the desired accuracy is achieved.

The optional arguments *epsrel* and *limit* are the desired relative error and the maximum number of subintervals, respectively. *epsrel* defaults to 1e-8 and *limit* is 200.

`quad_qag` returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0           if no problems were encountered;
- 1           if too many sub-intervals were done;
- 2           if excessive roundoff error is detected;
- 3           if extremely bad integrand behavior occurs;

6 if the input is invalid.

Examples:

```
(%i1) quad_qag (x^(1/2)*log(1/x), x, 0, 1, 3);
(%o1)  [.44444444444492108, 3.1700968502883E-9, 961, 0]
(%i2) integrate (x^(1/2)*log(1/x), x, 0, 1);
(%o2)  4
      -
      9
```

**quad\_qags** ( $f(x)$ ,  $x$ ,  $a$ ,  $b$ ,  $epsrel$ ,  $limit$ ) Function

Numerically integrate the given function using adaptive quadrature with extrapolation. The function to be integrated is  $f(x)$ , with dependent variable  $x$ , and the function is to be integrated between the limits  $a$  and  $b$ .

The optional arguments  $epsrel$  and  $limit$  are the desired relative error and the maximum number of subintervals, respectively.  $epsrel$  defaults to  $1e-8$  and  $limit$  is 200.

**quad\_qags** returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0 no problems were encountered;
- 1 too many sub-intervals were done;
- 2 excessive roundoff error is detected;
- 3 extremely bad integrand behavior occurs;
- 4 failed to converge
- 5 integral is probably divergent or slowly convergent
- 6 if the input is invalid.

Examples:

```
(%i1) quad_qags (x^(1/2)*log(1/x), x, 0, 1);
(%o1)  [.44444444444444448, 1.11022302462516E-15, 315, 0]
```

Note that **quad\_qags** is more accurate and efficient than **quad\_qag** for this integrand.

**quad\_qagi** ( $f(x)$ ,  $x$ ,  $a$ ,  $inftype$ ,  $epsrel$ ,  $limit$ ) Function

Numerically evaluate one of the following integrals

$$\int_a^{\infty} f(x)dx$$

$$\int_{-\infty}^a f(x)dx$$



$$\int_{-\infty}^{\infty} f(x)dx$$

using the Quadpack QAGI routine. The function to be integrated is  $f(x)$ , with dependent variable  $x$ , and the function is to be integrated over an infinite range.

The parameter *inf*type determines the integration interval as follows:

- inf**           The interval is from  $a$  to positive infinity.
- minf**         The interval is from negative infinity to  $a$ .
- both**         The interval is the entire real line.

The optional arguments *epsrel* and *limit* are the desired relative error and the maximum number of subintervals, respectively. *epsrel* defaults to 1e-8 and *limit* is 200.

**quad\_qagi** returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0           no problems were encountered;
- 1           too many sub-intervals were done;
- 2           excessive roundoff error is detected;
- 3           extremely bad integrand behavior occurs;
- 4           failed to converge
- 5           integral is probably divergent or slowly convergent
- 6           if the input is invalid.

Examples:

```
(%i1) quad_qagi (x^2*exp(-4*x), x, 0, inf);
(%o1) [0.03125, 2.95916102995002E-11, 105, 0]
(%i2) integrate (x^2*exp(-4*x), x, 0, inf);
(%o2) 1
      --
      32
```

**quad\_qawc** ( $f(x)$ ,  $x$ ,  $c$ ,  $a$ ,  $b$ , *epsrel*, *limit*)

Function

Numerically compute the Cauchy principal value of

$$\int_a^b \frac{f(x)}{x-c} dx$$

using the Quadpack QAWC routine. The function to be integrated is  $f(x)/(x-c)$ , with dependent variable  $x$ , and the function is to be integrated over the interval  $a$  to  $b$ .

The optional arguments *epsrel* and *limit* are the desired relative error and the maximum number of subintervals, respectively. *epsrel* defaults to 1e-8 and *limit* is 200.

`quad_qawc` returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0 no problems were encountered;
- 1 too many sub-intervals were done;
- 2 excessive roundoff error is detected;
- 3 extremely bad integrand behavior occurs;
- 6 if the input is invalid.

Examples:

```
(%i1) quad_qawc (2^(-5)*((x-1)^2+4^(-5))^( -1), x, 2, 0, 5);
(%o1) [- 3.130120337415925, 1.306830140249558E-8, 495, 0]
(%i2) integrate (2^(-alpha)*((x-1)^2 + 4^(-alpha))*(x-2))^( -1), x, 0, 5);
Principal Value
          alpha
          9 4          9
alpha      log(----- + -----)
4          64 4      + 4      64 4      + 4
(%o2) (-----)
          alpha
          2 4      + 2

          3 alpha          3 alpha
          -----          -----
          2          alpha/2          2          alpha/2
2 4      atan(4 4      ) 2 4      atan(4 4      ) alpha
- ----- - -----)/2
          alpha          alpha
          2 4      + 2          2 4      + 2
(%i3) ev (%, alpha=5, numer);
(%o3) - 3.130120337415917
```

**quad\_qawf** (*f(x)*, *x*, *a*, *omega*, *trig*, *epsabs*, *limit*, *maxp1*, *limlst*) Function  
 Numerically compute the a Fourier-type integral using the Quadpack QAWF routine.  
 The integral is

$$\int_a^\infty f(x)w(x)dx$$

The weight function *w* is selected by *trig*:

`cos`       $w(x) = \cos(\omega x)$

`cos`       $w(x) = \sin(\omega x)$

The optional arguments are:

`epsabs`    Desired absolute error of approximation. Default is 1d-10.

`limit`     Size of internal work array.  $(\text{limit} - \text{limlst})/2$  is the maximum number of subintervals to use. Default is 200.

`maxp1`    Maximum number of Chebyshev moments. Must be greater than 0. Default is 100.

`limlst`    Upper bound on the number of cycles. Must be greater than or equal to 3. Default is 10.

`epsabs` and `limit` are the desired relative error and the maximum number of subintervals, respectively. `epsrel` defaults to 1e-8 and `limit` is 200.

`quad_qawf` returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0            no problems were encountered;
- 1            too many sub-intervals were done;
- 2            excessive roundoff error is detected;
- 3            extremely bad integrand behavior occurs;
- 6            if the input is invalid.

Examples:

```
(%i1) quad_qawf (exp(-x^2), x, 0, 1, 'cos);
(%o1)  [.6901942235215714, 2.84846300257552E-11, 215, 0]
(%i2) integrate (exp(-x^2)*cos(x), x, 0, inf);
          - 1/4
          %e      sqrt(%pi)
(%o2)  -----
                2
(%i3) ev (% , numer);
(%o3)  .6901942235215714
```

**quad\_qawo** ( $f(x)$ ,  $x$ ,  $a$ ,  $b$ ,  $\omega$ ,  $\text{trig}$ ,  $\text{epsabs}$ ,  $\text{limit}$ ,  $\text{maxp1}$ ,  $\text{limlst}$ )

Function

Numerically compute the integral using the Quadpack QAWO routine:

$$\int_a^b f(x)w(x)dx$$

The weight function  $w$  is selected by `trig`:

`cos`       $w(x) = \cos(\omega x)$   
`sin`       $w(x) = \sin(\omega x)$

The optional arguments are:

`epsabs`    Desired absolute error of approximation. Default is 1d-10.  
`limit`     Size of internal work array.  $(limit - limlst)/2$  is the maximum number of subintervals to use. Default is 200.  
`maxpl`    Maximum number of Chebyshev moments. Must be greater than 0. Default is 100.  
`limlst`    Upper bound on the number of cycles. Must be greater than or equal to 3. Default is 10.

`epsabs` and `limit` are the desired relative error and the maximum number of subintervals, respectively. `epsrel` defaults to 1e-8 and `limit` is 200.

`quad_qawo` returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0            no problems were encountered;
- 1            too many sub-intervals were done;
- 2            excessive roundoff error is detected;
- 3            extremely bad integrand behavior occurs;
- 6            if the input is invalid.

Examples:

```
(%i1) quad_qawo (x^(-1/2)*exp(-2^(-2)*x), x, 1d-8, 20*2^2, 1, cos);
(%o1) [1.376043389877692, 4.72710759424899E-11, 765, 0]
(%i2) rectform (integrate (x^(-1/2)*exp(-2^(-alpha)*x) * cos(x), x, 0, inf));
          alpha/2 - 1/2                2 alpha
sqrt(%pi) 2                sqrt(sqrt(2      + 1) + 1)
(%o2) -----
              2 alpha
              sqrt(2      + 1)
(%i3) ev (% , alpha=2, numer);
(%o3) 1.376043390090716
```

**quad\_qaws** ( $f(x)$ ,  $x$ ,  $a$ ,  $b$ ,  $\alpha$ ,  $\beta$ ,  $wfun$ ,  $epsabs$ ,  $limit$ ) Function

Numerically compute the integral using the Quadpack QAWS routine:

$$\int_a^b f(x)w(x)dx$$

The weight function  $w$  is selected by `wfun`:

```

1      w(x) = (x - a)^alpha (b - x)^beta
2      w(x) = (x - a)^alpha (b - x)^beta log(x - a)
3      w(x) = (x - a)^alpha (b - x)^beta log(b - x)
2      w(x) = (x - a)^alpha (b - x)^beta log(x - a) log(b - x)

```

The optional arguments are:

*epsabs* Desired absolute error of approximation. Default is 1d-10.

*limit* Size of internal work array.  $(limit - limlst)/2$  is the maximum number of subintervals to use. Default is 200.

*epsabs* and *limit* are the desired relative error and the maximum number of subintervals, respectively. *epsrel* defaults to 1e-8 and *limit* is 200.

`quad_qaws` returns a list of four elements:

- an approximation to the integral,
- the estimated absolute error of the approximation,
- the number integrand evaluations,
- an error code.

The error code (fourth element of the return value) can have the values:

- 0 no problems were encountered;
- 1 too many sub-intervals were done;
- 2 excessive roundoff error is detected;
- 3 extremely bad integrand behavior occurs;
- 6 if the input is invalid.

Examples:

```

(%i1) quad_qaws (1/(x+1+2^(-4)), x, -1, 1, -0.5, -0.5, 1);
(%o1) [8.750097361672832, 1.24321522715422E-10, 170, 0]
(%i2) integrate ((1-x*x)^(-1/2)/(x+1+2^(-alpha)), x, -1, 1);
      alpha
Is 4 2      - 1 positive, negative, or zero?

```

pos;

```

      alpha      alpha
      2 %pi 2      sqrt(2 2      + 1)
(%o2) -----
      alpha
      4 2      + 2
(%i3) ev (% , alpha=4, numer);
(%o3) 8.750097361672829

```

## 22 Equations

### 22.1 Definitions for Equations

**%rnum\_list** Variable

Default value: []

**%rnum\_list** is the list of variables introduced in solutions by **algsys**. **%r** variables are added to **%rnum\_list** in the order they are created. This is convenient for doing substitutions into the solution later on. It's recommended to use this list rather than doing **concat ('%r, j)**.

**algexact** Variable

Default value: **false**

**algexact** affects the behavior of **algsys** as follows:

If **algexact** is **true**, **algsys** always calls **solve** and then uses **realroots** on **solve**'s failures.

If **algexact** is **false**, **solve** is called only if the eliminant was not univariate, or if it was a quadratic or biquadratic.

Thus **algexact: true** doesn't guarantee only exact solutions, just that **algsys** will first try as hard as it can to give exact solutions, and only yield approximations when all else fails.

**algsys** (*[expr\_1, ..., expr\_m], [x\_1, ..., x\_n]*) Function

**algsys** (*[eqn\_1, ..., eqn\_m], [x\_1, ..., x\_n]*) Function

Solves the simultaneous polynomials *expr\_1, ..., expr\_m* or polynomial equations *eqn\_1, ..., eqn\_m* for the variables *x\_1, ..., x\_n*. An expression *expr* is equivalent to an equation *expr = 0*. There may be more equations than variables or vice versa.

**algsys** returns a list of solutions, with each solution given as a list of equations stating values of the variables *x\_1, ..., x\_n* which satisfy the system of equations. If **algsys** cannot find a solution, an empty list [] is returned.

The symbols **%r1, %r2, ...**, are introduced as needed to represent arbitrary parameters in the solution; these variables are also appended to the list **%rnum\_list**.

The method is as follows:

- (1) First the equations are factored and split into subsystems.
- (2) For each subsystem *S<sub>i</sub>*, an equation *E* and a variable *x* are selected. The variable is chosen to have lowest nonzero degree. Then the resultant of *E* and *E<sub>j</sub>* with respect to *x* is computed for each of the remaining equations *E<sub>j</sub>* in the subsystem *S<sub>i</sub>*. This yields a new subsystem *S<sub>i</sub>'* in one fewer variables, as *x* has been eliminated. The process now returns to (1).
- (3) Eventually, a subsystem consisting of a single equation is obtained. If the equation is multivariate and no approximations in the form of floating point numbers have been introduced, then **solve** is called to find an exact solution.

In some cases, `solve` is not be able to find a solution, or if it does the solution may be a very large expression.

If the equation is univariate and is either linear, quadratic, or biquadratic, then again `solve` is called if no approximations have been introduced. If approximations have been introduced or the equation is not univariate and neither linear, quadratic, or biquadratic, then if the switch `realonly` is `true`, the function `realroots` is called to find the real-valued solutions. If `realonly` is `false`, then `allroots` is called which looks for real and complex-valued solutions.

If `algsys` produces a solution which has fewer significant digits than required, the user can change the value of `algepsilon` to a higher value.

If `algexact` is set to `true`, `solve` will always be called.

(4) Finally, the solutions obtained in step (3) are substituted into previous levels and the solution process returns to (1).

When `algsys` encounters a multivariate equation which contains floating point approximations (usually due to its failing to find exact solutions at an earlier stage), then it does not attempt to apply exact methods to such equations and instead prints the message: "algsys cannot solve - system too complicated."

Interactions with `radcan` can produce large or complicated expressions. In that case, it may be possible to isolate parts of the result with `pickapart` or `reveal`.

Occasionally, `radcan` may introduce an imaginary unit `%i` into a solution which is actually real-valued.

Examples:

```
(%i1) e1: 2*x*(1 - a1) - 2*(x - 1)*a2;
(%o1)          2 (1 - a1) x - 2 a2 (x - 1)
(%i2) e2: a2 - a1;
(%o2)          a2 - a1
(%i3) e3: a1*(-y - x^2 + 1);
(%o3)          a1 (- y - x^2 + 1)
(%i4) e4: a2*(y - (x - 1)^2);
(%o4)          a2 (y - (x - 1)^2)
(%i5) algsys ([e1, e2, e3, e4], [x, y, a1, a2]);
(%o5) [[x = 0, y = %r1, a1 = 0, a2 = 0],
[x = 1, y = 0, a1 = 1, a2 = 1]]
(%i6) e1: x^2 - y^2;
(%o6)          x^2 - y^2
(%i7) e2: -1 - y + 2*y^2 - x + x^2;
(%o7)          2 y^2 - y + x^2 - x - 1
(%i8) algsys ([e1, e2], [x, y]);
(%o8) [[x = - ----, y = ----],
          sqrt(3)      sqrt(3)]
```

$$\left[ x = \frac{1}{\sqrt{3}}, y = -\frac{1}{\sqrt{3}} \right], \left[ x = -\frac{1}{3}, y = -\frac{1}{3} \right], [x = 1, y = 1]$$

**allroots** (*expr*) Function  
**allroots** (*eqn*) Function

Computes numerical approximations of the real and complex roots of the polynomial *expr* or polynomial equation *eqn* of one variable.

The flag **polyfactor** when **true** causes **allroots** to factor the polynomial over the real numbers if the polynomial is real, or over the complex numbers, if the polynomial is complex.

**allroots** may give inaccurate results in case of multiple roots. If the polynomial is real, **allroots** (*%i\*p*) may yield more accurate approximations than **allroots** (*p*), as **allroots** invokes a different algorithm in that case.

**allroots** rejects non-polynomials. It requires that the numerator after **rat**'ing should be a polynomial, and it requires that the denominator be at most a complex number. As a result of this **allroots** will always return an equivalent (but factored) expression, if **polyfactor** is **true**.

For complex polynomials an algorithm by Jenkins and Traub is used (Algorithm 419, *Comm. ACM*, vol. 15, (1972), p. 97). For real polynomials the algorithm used is due to Jenkins (Algorithm 493, *ACM TOMS*, vol. 1, (1975), p.178).

Examples:

```
(%i1) eqn: (1 + 2*x)^3 = 13.5*(1 + x^5);
              3           5
(%o1)          (2 x + 1) = 13.5 (x + 1)
(%i2) soln: allroots (eqn);
(%o2) [x = .8296749902129361, x = - 1.015755543828121,
x = .9659625152196369 %i - .4069597231924075,
x = - .9659625152196369 %i - .4069597231924075, x = 1.0]
(%i3) for e in soln
      do (e2: subst (e, eqn), disp (expand (lhs(e2) - rhs(e2))));
          - 3.5527136788005E-15
          - 5.32907051820075E-15
          4.44089209850063E-15 %i - 4.88498130835069E-15
          - 4.44089209850063E-15 %i - 4.88498130835069E-15
          3.5527136788005E-15
(%o3)          done
(%i4) polyfactor: true$
(%i5) allroots (eqn);
```



$$\begin{aligned}
 (& \%o5) - 13.5 (x - 1.0) (x - .8296749902129361) \\
 & (x + 1.015755543828121)^2 (x + .8139194463848151 x \\
 & + 1.098699797110288)
 \end{aligned}$$

**backsubst**

Variable

Default value: true

When `backsubst` is false, prevents back substitution after the equations have been triangularized. This may be helpful in very big problems where back substitution would cause the generation of extremely large expressions.

**breakup**

Variable

Default value: true

When `breakup` is true, `solve` expresses solutions of cubic and quartic equations in terms of common subexpressions, which are assigned to intermediate expression labels (`%t1`, `%t2`, etc.). Otherwise, common subexpressions are not identified.

`breakup: true` has an effect only when `programmode` is false.

Examples:

```

(%i1) programmode: false$
(%i2) breakup: true$
(%i3) solve (x^3 + x^2 - 1);

```

```

(%t3)           sqrt(23)   25 1/3
             (----- + --)
              6 sqrt(3)   54

```

Solution:

```

(%t4)           sqrt(3) %i   1
              2           2   1
          x = (- ----- - -) %t3 + ----- - -
                2           2     9 %t3   3

```

```

(%t5)           sqrt(3) %i   1
              2           2   1
          x = (----- - -) %t3 + ----- - -
                2           2     9 %t3   3

```

```

(%t6)           1   1
              9 %t3  3
          x = %t3 + ----- - -

```

```

(%o6)           [%t4, %t5, %t6]

```

```

(%i6) breakup: false$
(%i7) solve (x^3 + x^2 - 1);

```

Solution:

$$\begin{aligned}
 (\%t7) \quad x &= \frac{\frac{\sqrt{3} i}{2} - \frac{1}{2}}{\frac{\sqrt{23}}{6\sqrt{3}} + \frac{25}{54}} + \left( \frac{\sqrt{23}}{6\sqrt{3}} + \frac{25}{54} \right)^{-1/3} \\
 (\%t8) \quad x &= \left( \frac{\sqrt{23}}{6\sqrt{3}} + \frac{25}{54} \right)^{-1/3} \left( \frac{\sqrt{3} i}{2} - \frac{1}{2} \right) \\
 (\%t9) \quad x &= \left( \frac{\sqrt{23}}{6\sqrt{3}} + \frac{25}{54} \right)^{-1/3} + \frac{1}{3 \left( \frac{\sqrt{23}}{6\sqrt{3}} + \frac{25}{54} \right)^{1/3}} \\
 (\%o9) \quad &[\%t7, \%t8, \%t9]
 \end{aligned}$$

**dimension** (*eqn*)

Function

**dimension** (*eqn\_1, ..., eqn\_n*)

Function

*dimen* is a package for dimensional analysis. `load("dimen")` loads this package. `demo("dimen")` displays a short demonstration.

**dispflag**

Variable

Default value: `true`

If set to `false` within a `block` will inhibit the display of output generated by the solve functions called from within the `block`. Termination of the `block` with a dollar sign, `$`, sets `dispflag` to `false`.

**funcsolve** (*eqn, g(t)*)

Function

Returns  $[g(t) = \dots]$  or  $[]$ , depending on whether or not there exists a rational function  $g(t)$  satisfying *eqn*, which must be a first order, linear polynomial in (for this case)  $g(t)$  and  $g(t+1)$

```
(%i1) eqn: (n + 1)*f(n) - (n + 3)*f(n + 1)/(n + 1) = (n - 1)/(n + 2);
(%o1)      (n + 1) f(n) - ----- = -----
              n + 1          n + 2
(%i2) funcsolve (eqn, f(n));

Dependent equations eliminated: (4 3)
(%o2)      f(n) = -----
              n
              (n + 1) (n + 2)
```

Warning: this is a very rudimentary implementation – many safety checks and obvious generalizations are missing.

## globalsolve

Variable

Default value: false

When `globalsolve` is true, solved-for variables are assigned the solution values found by `solve`.

Examples:

```
(%i1) globalsolve: true$
(%i2) solve ([x + 3*y = 2, 2*x - y = 5], [x, y]);
Solution

(%t2)      x : --
              7

(%t3)      y : - -
              7
(%o3)      [[%t2, %t3]]
(%i3) x;

(%o3)      17
              --
              7

(%i4) y;

(%o4)      1
              - -
              7

(%i5) globalsolve: false$
(%i6) kill (x, y)$
(%i7) solve ([x + 3*y = 2, 2*x - y = 5], [x, y]);
Solution

(%t7)      x = --
              7
```

```

(%t8)          y = - -
                7
(%o8)          [[%t7, %t8]]
(%i8) x;
(%o8)          x
(%i9) y;
(%o9)          y

```

**ieqn** (*ie, unk, tech, n, guess*) Function

`ieqn` is a package for solving integral equations. `load ("ieqn")` loads this package.

*ie* is the integral equation; *unk* is the unknown function; *tech* is the technique to be tried from those given above (*tech* = `first` means: try the first technique which finds a solution; *tech* = `all` means: try all applicable techniques); *n* is the maximum number of terms to take for `taylor`, `neumann`, `firstkindseries`, or `fredseries` (it is also the maximum depth of recursion for the differentiation method); *guess* is the initial guess for `neumann` or `firstkindseries`.

Default values for the 2nd thru 5th parameters are:

*unk*:  $p(x)$ , where  $p$  is the first function encountered in an integrand which is unknown to Maxima and  $x$  is the variable which occurs as an argument to the first occurrence of  $p$  found outside of an integral in the case of `secondkind` equations, or is the only other variable besides the variable of integration in `firstkind` equations. If the attempt to search for  $x$  fails, the user will be asked to supply the independent variable.

*tech*: `first`

*n*: 1

*guess*: `none` which will cause `neumann` and `firstkindseries` to use  $f(x)$  as an initial guess.

**ieqnprint** Variable

Default value: `true`

`ieqnprint` governs the behavior of the result returned by the `ieqn` command. When `ieqnprint` is `false`, the lists returned by the `ieqn` function are of the form

`[solution, technique used, nterms, flag]`

where *flag* is absent if the solution is exact.

Otherwise, it is the word `approximate` or `incomplete` corresponding to an inexact or non-closed form solution, respectively. If a series method was used, *nterms* gives the number of terms taken (which could be less than the *n* given to `ieqn` if an error prevented generation of further terms).

**lhs** (*eqn*) Function

Returns the left side of the equation *eqn*.

If the argument is not an equation, `lhs` returns the argument.

See also `rhs`.

Example:

```
(%i1) e: x^2 + y^2 = z^2;
(%o1)          2    2    2
          y  + x  = z
(%i2) lhs (e);
(%o2)          2    2
          y  + x
(%i3) rhs (e);
(%o3)          2
          z
```

**linsolve** (*[expr\_1, ..., expr\_m], [x\_1, ..., x\_n]*) Function

Solves the list of simultaneous linear equations for the list of variables. The expressions must each be polynomials in the variables and may be equations.

When `globalsolve` is `true` then variables which are solved for will be set to the solution of the set of simultaneous equations.

When `backsubst` is `false`, `linsolve` does not carry out back substitution after the equations have been triangularized. This may be necessary in very big problems where back substitution would cause the generation of extremely large expressions.

When `linsolve_params` is `true`, `linsolve` also generates the `%r` symbols used to represent arbitrary parameters described in the manual under `algsys`. Otherwise, `linsolve` solves an under-determined system of equations with some variables expressed in terms of others.

```
(%i1) e1: x + z = y$
(%i2) e2: 2*a*x - y = 2*a^2$
(%i3) e3: y - 2*z = 2$
(%i4) linsolve ([e1, e2, e3], [x, y, z]);
(%o4)          [x = a + 1, y = 2 a, z = a - 1]
```

**linsolvewarn** Variable

Default value: `true`

When `linsolvewarn` is `true`, `linsolve` prints a message "Dependent equations eliminated".

**linsolve\_params** Variable

Default value: `true`

When `linsolve_params` is `true`, `linsolve` also generates the `%r` symbols used to represent arbitrary parameters described in the manual under `algsys`. Otherwise, `linsolve` solves an under-determined system of equations with some variables expressed in terms of others.

**multiplicities** Variable

Default value: `not_set_yet`

`multiplicities` is set to a list of the multiplicities of the individual solutions returned by `solve` or `realroots`.

**nroots** (*p*, *low*, *high*) Function

Returns the number of real roots of the real univariate polynomial *p* in the half-open interval (*low*, *high*]. The endpoints of the interval may be `minf` or `inf`. infinity and plus infinity.

`nroots` uses the method of Sturm sequences.

```
(%i1) p: x^10 - 2*x^4 + 1/2$
(%i2) nroots (p, -6, 9.1);
(%o2)                                     4
```

**nthroot** (*p*, *n*) Function

where *p* is a polynomial with integer coefficients and *n* is a positive integer returns *q*, a polynomial over the integers, such that  $q^n = p$  or prints an error message indicating that *p* is not a perfect *n*th power. This routine is much faster than `factor` or even `sqfr`.

**programmode** Variable

Default value: `true`

When `programmode` is `true`, `solve`, `realroots`, `allroots`, and `linsolve` return solutions as elements in a list. (Except when `backsubst` is set to `false`, in which case `programmode: false` is assumed.)

When `programmode` is `false`, `solve`, etc. create intermediate expression labels `%t1`, `t2`, etc., and assign the solutions to them.

**realonly** Variable

Default value: `false`

When `realonly` is `true`, `algsys` returns only those solutions which are free of `%i`.

**realroots** (*poly*, *bound*) Function

Finds all of the real roots of the real univariate polynomial *poly* within a tolerance of *bound* which, if less than 1, causes all integral roots to be found exactly. The parameter *bound* may be arbitrarily small in order to achieve any desired accuracy. The first argument may also be an equation. `realroots` sets `multiplicities`, useful in case of multiple roots. `realroots (p)` is equivalent to `realroots (p, rootsepsilon)`. `rootsepsilon` is a real number used to establish the confidence interval for the roots. Do `example (realroots)` for an example.

**rhs** (*eqn*) Function

Returns the right side of the equation *eqn*.

If the argument is not an equation, `rhs` returns 0.

See also `lhs`.

Example:

```
(%i1) e: x^2 + y^2 = z^2;
(%o1)          2      2      2
              y  + x  = z
(%i2) lhs (e);
```

```

(%o2)          2    2
              y  + x
(%i3) rhs (e);
(%o3)          2
              z

```

**rootsconmode**

Variable

Default value: true

`rootsconmode` governs the behavior of the `rootscontract` command. See `rootscontract` for details.

**rootscontract** (*expr*)

Function

Converts products of roots into roots of products. For example, `rootscontract (sqrt(x)*y^(3/2))` yields `sqrt(x*y^3)`.

When `radexpand` is true and `domain` is real, `rootscontract` converts `abs` into `sqrt`, e.g., `rootscontract (abs(x)*sqrt(y))` yields `sqrt(x^2*y)`.

There is an option `rootsconmode` affecting `rootscontract` as follows:

Problem	Value of <code>rootsconmode</code>	Result of applying <code>rootscontract</code>
$x^{1/2}y^{3/2}$	false	$(x*y^3)^{1/2}$
$x^{1/2}y^{1/4}$	false	$x^{1/2}y^{1/4}$
$x^{1/2}y^{1/4}$	true	$(x*y^{1/2})^{1/2}$
$x^{1/2}y^{1/3}$	true	$x^{1/2}y^{1/3}$
$x^{1/2}y^{1/4}$	all	$(x^2*y)^{1/4}$
$x^{1/2}y^{1/3}$	all	$(x^3*y^2)^{1/6}$

When `rootsconmode` is false, `rootscontract` contracts only with respect to rational number exponents whose denominators are the same. The key to the `rootsconmode: true` examples is simply that 2 divides into 4 but not into 3. `rootsconmode: all` involves taking the least common multiple of the denominators of the exponents.

`rootscontract` uses `ratsimp` in a manner similar to `logcontract`.

Examples:

```

(%i1) rootsconmode: false$
(%i2) rootscontract (x^(1/2)*y^(3/2));
(%o2)          3
              sqrt(x y )
(%i3) rootscontract (x^(1/2)*y^(1/4));
(%o3)          1/4
              sqrt(x) y
(%i4) rootsconmode: true$
(%i5) rootscontract (x^(1/2)*y^(1/4));
(%o5)          sqrt(x sqrt(y))
(%i6) rootscontract (x^(1/2)*y^(1/3));
(%o6)          1/3
              sqrt(x) y
(%i7) rootsconmode: all$

```

```
(%i8) rootscontract (x^(1/2)*y^(1/4));
              2  1/4
(%o8)          (x y)
(%i9) rootscontract (x^(1/2)*y^(1/3));
              3  2 1/6
(%o9)          (x y )
(%i10) rootsconmode: false$
(%i11) rootscontract (sqrt(sqrt(x) + sqrt(1 + x))
              *sqrt(sqrt(1 + x) - sqrt(x)));
(%o11)          1
(%i12) rootsconmode: true$
(%i13) rootscontract (sqrt(5 + sqrt(5)) - 5^(1/4)*sqrt(1 + sqrt(5)));
(%o13)          0
```

**rootsepsilon**

Variable

Default value: 1.0e-7

**rootsepsilon** is the tolerance which establishes the confidence interval for the roots found by the **realroots** function.

**solve** (*expr*, *x*)

Function

**solve** (*expr*)

Function

**solve** ([*eqn\_1*, ..., *eqn\_n*], [*x\_1*, ..., *x\_n*])

Function

Solves the algebraic equation *expr* for the variable *x* and returns a list of solution equations in *x*. If *expr* is not an equation, the equation  $expr = 0$  is assumed in its place. *x* may be a function (e.g.  $f(x)$ ), or other non-atomic expression except a sum or product. *x* may be omitted if *expr* contains only one variable. *expr* may be a rational expression, and may contain trigonometric functions, exponentials, etc.

The following method is used:

Let *E* be the expression and *X* be the variable. If *E* is linear in *X* then it is trivially solved for *X*. Otherwise if *E* is of the form  $A * X^N + B$  then the result is  $(-B/A)^{1/N}$  times the *N*'th roots of unity.

If *E* is not linear in *X* then the gcd of the exponents of *X* in *E* (say *N*) is divided into the exponents and the multiplicity of the roots is multiplied by *N*. Then **solve** is called again on the result. If *E* factors then **solve** is called on each of the factors. Finally **solve** will use the quadratic, cubic, or quartic formulas where necessary.

In the case where *E* is a polynomial in some function of the variable to be solved for, say  $F(X)$ , then it is first solved for  $F(X)$  (call the result *C*), then the equation  $F(X)=C$  can be solved for *X* provided the inverse of the function *F* is known.

**breakup** if **false** will cause **solve** to express the solutions of cubic or quartic equations as single expressions rather than as made up of several common subexpressions which is the default.

**multiplicities** - will be set to a list of the multiplicities of the individual solutions returned by **solve**, **realroots**, or **allroots**. Try **apropos (solve)** for the switches which affect **solve**. **describe** may then be used on the individual switch names if their purpose is not clear.



`solve ([eqn_1, ..., eqn_n], [x_1, ..., x_n])` solves a system of simultaneous (linear or non-linear) polynomial equations by calling `linsolve` or `algsys` and returns a list of the solution lists in the variables. In the case of `linsolve` this list would contain a single list of solutions. It takes two lists as arguments. The first list represents the equations to be solved; the second list is a list of the unknowns to be determined. If the total number of variables in the equations is equal to the number of equations, the second argument-list may be omitted. For linear systems if the given equations are not compatible, the message `inconsistent` will be displayed (see the `solve_inconsistent_error` switch); if no unique solution exists, then `singular` will be displayed.

Examples:

```
(%i1) solve (asin (cos (3*x))*(f(x) - 1), x);
```

SOLVE is using arc-trig functions to get a solution.  
Some solutions will be lost.

```
(%o1) [x = ---, f(x) = 1]
          %pi
          6
```

```
(%i2) ev (solve (5^f(x) = 125, f(x)), solveradcan);
          log(125)
```

```
(%o2) [f(x) = -----]
          log(5)
```

```
(%i3) [4*x^2 - y^2 = 12, x*y - x = 2];
          2      2
```

```
(%o3) [4 x  - y  = 12, x y - x = 2]
```

```
(%i4) solve (%, [x, y]);
```

```
(%o4) [[x = 2, y = 2], [x = .5202594388652008 %i
```

```
- .1331240357358706, y = .0767837852378778
```

```
- 3.608003221870287 %i], [x = - .5202594388652008 %i
```

```
- .1331240357358706, y = 3.608003221870287 %i
```

```
+ .0767837852378778], [x = - 1.733751846381093,
```

```
y = - .1535675710019696]]
```

```
(%i5) solve (1 + a*x + x^3, x);
```

```
(%o5) [x = (- ----- - -) (----- - -)
          sqrt(3) %i  1  sqrt(4 a  + 27)  1 1/3
          2          2          6 sqrt(3)  2
```

```
          sqrt(3) %i  1
          (----- - -) a
          2          2
```

```
- -----, x =
          3
```

$$3 \left( \frac{\sqrt{4a+27}}{6\sqrt{3}} - \frac{1}{2} \right)^{1/3}$$

$$\left( \frac{\sqrt{3}i}{2} - \frac{1}{2} \right) \left( \frac{\sqrt{4a+27}}{6\sqrt{3}} - \frac{1}{2} \right)^{1/3}$$

$$- \frac{\left( -\frac{\sqrt{3}i}{2} - \frac{1}{2} \right) a}{3}, x =$$

$$3 \left( \frac{\sqrt{4a+27}}{6\sqrt{3}} - \frac{1}{2} \right)^{1/3}$$

$$\left( \frac{\sqrt{4a+27}}{6\sqrt{3}} - \frac{1}{2} \right)^{1/3} - \frac{a}{3 \left( \frac{\sqrt{4a+27}}{6\sqrt{3}} - \frac{1}{2} \right)^{1/3}}$$

```
(%i6) solve (x^3 - 1);
```

$$[x = \frac{\sqrt{3}i - 1}{2}, x = \frac{\sqrt{3}i + 1}{2}, x = 1]$$

```
(%i7) solve (x^6 - 1);
```

$$[x = \frac{\sqrt{3}i + 1}{2}, x = \frac{\sqrt{3}i - 1}{2}, x = -1,$$

$$x = -\frac{\sqrt{3}i + 1}{2}, x = -\frac{\sqrt{3}i - 1}{2}, x = 1]$$

```
(%i8) ev (x^6 - 1, %[1]);
```

$$\frac{(\sqrt{3}i + 1)^6}{64} - 1$$

```
(%i9) expand (%);
```

$$0$$

```
(%i10) x^2 - 1;
```

$$x^2 - 1$$

```
(%i11) solve (%, x);
```

$$[x = -1, x = 1]$$

```
(%i12) ev (%th(2), %[1]);
(%o12)                                0
```

**solvedecomposes** Variable

Default value: true

When `solvedecomposes` is true, `solve` calls `polydecomp` if asked to solve polynomials.

**solveexplicit** Variable

Default value: false

When `solveexplicit` is true, inhibits `solve` from returning implicit solutions, that is, solutions of the form  $F(x) = 0$  where  $F$  is some function.

**solvefactors** Variable

Default value: true

When `solvefactors` is false, `solve` does not try to factor the expression. The false setting may be desired in some cases where factoring is not necessary.

**solvenullwarn** Variable

Default value: true

When `solvenullwarn` is true, `solve` prints a warning message if called with either a null equation list or a null variable list. For example, `solve ([], [])` would print two warning messages and return `[]`.

**solveradcan** Variable

Default value: false

When `solveradcan` is true, `solve` calls `radcan` which makes `solve` slower but will allow certain problems containing exponentials and logarithms to be solved.

**solvetricwarn** Variable

Default value: true

When `solvetricwarn` is true, `solve` may print a message saying that it is using inverse trigonometric functions to solve the equation, and thereby losing solutions.

**solve\_inconsistent\_error** Variable

Default value: true

When `solve_inconsistent_error` is true, `solve` and `linsolve` give an error if the equations to be solved are inconsistent.

If false, `solve` and `linsolve` return an empty list `[]` if the equations are inconsistent.

Example:

```
(%i1) solve_inconsistent_error: true$
(%i2) solve ([a + b = 1, a + b = 2], [a, b]);
Inconsistent equations: (2)
-- an error. Quitting. To debug this try debugmode(true);
(%i3) solve_inconsistent_error: false$
(%i4) solve ([a + b = 1, a + b = 2], [a, b]);
(%o4)                                []
```

## 23 Differential Equations

### 23.1 Definitions for Differential Equations

**bc2** (*solution*, *xval1*, *yval1*, *xval2*, *yval2*) Function

Solves boundary value problem for second order differential equation. Here: *solution* is a general solution to the equation, as found by *ode2*, *xval1* is an equation for the independent variable in the form  $x = x0$ , and *yval1* is an equation for the dependent variable in the form  $y = y0$ . The *xval2* and *yval2* are equations for these variables at another point. See *ode2* for example of usage.

**dsolve** (*eqn*, *x*) Function

**dsolve** (*[eqn\_1, ..., eqn\_n]*, *[x\_1, ..., x\_n]*) Function

The function *dsolve* solves systems of linear ordinary differential equations using Laplace transform. Here the *eqn*'s are differential equations in the dependent variables  $x_1, \dots, x_n$ . The functional relationships must be explicitly indicated in both the equations and the variables. For example

```
'diff(f,x,2)=sin(x)+'diff(g,x);
'diff(f,x)+x^2-f=2*'diff(g,x,2);
```

is not the proper format. The correct way is:

```
'diff(f(x),x,2)=sin(x)+'diff(g(x),x);
'diff(f(x),x)+x^2-f=2*'diff(g(x),x,2);
```

The call is then `dsolve(['%o3,%o4],[f(x),g(x)]); .`

If initial conditions at 0 are known, they should be supplied before calling *dsolve* by using *atvalue*.

```
(%i1) 'diff(f(x),x)='diff(g(x),x)+sin(x);
      d          d
```

```
(%o1)      -- (f(x)) = -- (g(x)) + sin(x)
      dx          dx
```

```
(%i2) 'diff(g(x),x,2)='diff(f(x),x)-cos(x);
      2
```

```
(%o2)      --- (g(x)) = -- (f(x)) - cos(x)
      2          dx
```

```
(%i3) atvalue('diff(g(x),x),x=0,a);
```

```
(%o3)      a
```

```
(%i4) atvalue(f(x),x=0,1);
```

```
(%o4)      1
```

```
(%i5) dsolve(['%o1,%o2],[f(x),g(x)]);
```

```
(%o5) [f(x) = a %ex - a + 1, g(x) =
```

```
cos(x) + a %ex - a + g(0) - 1]
```

```
(%i6) [%o1,%o2],%o5,diff;
(%o6) [a %ex = a %ex , a %ex - cos(x) = a %ex - cos(x)]
```

If `desolve` cannot obtain a solution, it returns `false`.

**ic1** (*solution, xval, yval*) Function

Solves initial value problem for first order differential equation. Here: *solution* is a general solution to the equation, as found by `ode2`, *xval* is an equation for the independent variable in the form  $x = x_0$ , and *yval* is an equation for the dependent variable in the form  $y = y_0$ . See `ode2` for example of usage.

**ic2** (*solution, xval, yval, dval*) Function

Solves initial value problem for second order differential equation. Here: *solution* is a general solution to the equation, as found by `ode2`, *xval* is an equation for the independent variable in the form  $x = x_0$ , *yval* is an equation for the dependent variable in the form  $y = y_0$ , and *dval* is an equation for the derivative of the dependent variable with respect to independent variable evaluated at the point *xval*. See `ode2` for example of usage.

**ode2** (*eqn, dvar, ivar*) Function

The function `ode2` solves ordinary differential equations of first or second order. It takes three arguments: an ODE *eqn*, the dependent variable *dvar*, and the independent variable *ivar*. When successful, it returns either an explicit or implicit solution for the dependent variable. `%c` is used to represent the constant in the case of first order equations, and `%k1` and `%k2` the constants for second order equations. If `ode2` cannot obtain a solution for whatever reason, it returns `false`, after perhaps printing out an error message. The methods implemented for first order equations in the order in which they are tested are: linear, separable, exact - perhaps requiring an integrating factor, homogeneous, Bernoulli's equation, and a generalized homogeneous method. For second order: constant coefficient, exact, linear homogeneous with non-constant coefficients which can be transformed to constant coefficient, the Euler or equidimensional equation, the method of variation of parameters, and equations which are free of either the independent or of the dependent variable so that they can be reduced to two first order linear equations to be solved sequentially. In the course of solving ODEs, several variables are set purely for informational purposes: `method` denotes the method of solution used e.g. `linear`, `intfactor` denotes any integrating factor used, `odeindex` denotes the index for Bernoulli's method or for the generalized homogeneous method, and `yp` denotes the particular solution for the variation of parameters technique.

In order to solve initial value problems (IVPs) and boundary value problems (BVPs), the routine `ic1` is available for first order equations, and `ic2` and `bc2` for second order IVPs and BVPs, respectively.

Example:

```
(%i1) x^2*'diff(y,x) + 3*y*x = sin(x)/x;
      2 dy          sin(x)
```

```
(%o1)          x  -- + 3 x y = -----
              dx          x
(%i2) ode2(%,y,x);
              %c - cos(x)
(%o2)          y = -----
              3
              x
(%i3) ic1(%o2,x=%pi,y=0);
              cos(x) + 1
(%o3)          y = - -----
              3
              x
(%i4) 'diff(y,x,2) + y*'diff(y,x)^3 = 0;
              2
              d y      dy 3
(%o4)          --- + y (---) = 0
              2        dx
              dx
(%i5) ode2(%,y,x);
              3
              y  + 6 %k1 y
(%o5)          ----- = x + %k2
              6
(%i6) ratsimp(ic2(%o5,x=0,y=0,'diff(y,x)=2));
              3
              2 y  - 3 y
(%o6)          - ----- = x
              6
(%i7) bc2(%o5,x=0,y=1,x=1,y=3);
              3
              y  - 10 y      3
(%o7)          ----- = x - -
              6                2
```



## 24 Numerical

### 24.1 Introduction to Numerical

### 24.2 Fourier packages

The `fft` package comprises functions for the numerical (not symbolic) computation of the fast Fourier transform. `load ("fft")` loads this package. See `fft`.

The `fourie` package comprises functions for the symbolic computation of Fourier series. `load ("fourie")` loads this package. There are functions in the `fourie` package to calculate Fourier integral coefficients and some functions for manipulation of expressions. See `Definitions for Fourier Series`.

### 24.3 Definitions for Numerical

**polartorect** (*magnitude\_array, phase\_array*) Function

Translates complex values of the form  $r e^{i t}$  to the form  $a + b i$ . `load ("fft")` loads this function into Maxima. See also `fft`.

The magnitude and phase,  $r$  and  $t$ , are taken from *magnitude\_array* and *phase\_array*, respectively. The original values of the input arrays are replaced by the real and imaginary parts,  $a$  and  $b$ , on return. The outputs are calculated as

$$\begin{aligned} a &: r \cos (t) \\ b &: r \sin (t) \end{aligned}$$

The input arrays must be the same size and 1-dimensional. The array size need not be a power of 2.

`polartorect` is the inverse function of `recttopolar`.

**recttopolar** (*real\_array, imaginary\_array*) Function

Translates complex values of the form  $a + b i$  to the form  $r e^{i t}$ . `load ("fft")` loads this function into Maxima. See also `fft`.

The real and imaginary parts,  $a$  and  $b$ , are taken from *real\_array* and *imaginary\_array*, respectively. The original values of the input arrays are replaced by the magnitude and angle,  $r$  and  $t$ , on return. The outputs are calculated as

$$\begin{aligned} r &: \sqrt{a^2 + b^2} \\ t &: \operatorname{atan2} (b, a) \end{aligned}$$

The computed angle is in the range  $-\pi$  to  $\pi$ .

The input arrays must be the same size and 1-dimensional. The array size need not be a power of 2.

`recttopolar` is the inverse function of `polartorect`.



**ift** (*real\_array*, *imaginary\_array*) Function  
 Fast inverse discrete Fourier transform. `load ("fft")` loads this function into Maxima.

`ift` carries out the inverse complex fast Fourier transform on 1-dimensional floating point arrays. The inverse transform is defined as

$$x[j]: \text{sum} (y[j] \exp (+2 \%i \%pi j k / n), k, 0, n-1)$$

See `fft` for more details.

**fft** (*real\_array*, *imaginary\_array*) Function

**ift** (*real\_array*, *imaginary\_array*) Function

**recttopolar** (*real\_array*, *imaginary\_array*) Function

**polartorect** (*magnitude\_array*, *phase\_array*) Function

Fast Fourier transform and related functions. `load ("fft")` loads these functions into Maxima.

`fft` and `ift` carry out the complex fast Fourier transform and inverse transform, respectively, on 1-dimensional floating point arrays. The size of *imaginary\_array* must equal the size of *real\_array*.

`fft` and `ift` operate in-place. That is, on return from `fft` or `ift`, the original content of the input arrays is replaced by the output. The `fillarray` function can make a copy of an array, should it be necessary.

The discrete Fourier transform and inverse transform are defined as follows. Let *x* be the original data, with

$$x[i]: \text{real\_array}[i] + \%i \text{imaginary\_array}[i]$$

Let *y* be the transformed data. The forward and inverse transforms are

$$y[k]: (1/n) \text{sum} (x[j] \exp (-2 \%i \%pi j k / n), j, 0, n-1)$$

$$x[j]: \text{sum} (y[k] \exp (+2 \%i \%pi j k / n), k, 0, n-1)$$

Suitable arrays can be allocated by the `array` function. For example:

```
array (my_array, float, n-1)$
```

declares a 1-dimensional array with *n* elements, indexed from 0 through *n*-1 inclusive. The number of elements *n* must be equal to  $2^m$  for some *m*.

`fft` can be applied to real data (imaginary array all zeros) to obtain sine and cosine coefficients. After calling `fft`, the sine and cosine coefficients, say *a* and *b*, can be calculated as

```
a[0]: real_array[0]
b[0]: 0
```

and

```
a[j]: real_array[j] + real_array[n-j]
b[j]: imaginary_array[j] - imaginary_array[n-j]
```

for *j* equal to 1 through *n*/2-1, and

```
a[n/2]: real_array[n/2]
b[n/2]: 0
```

`recttopolar` translates complex values of the form  $a + b \%i$  to the form  $r \%e^{(\%i t)}$ . See `recttopolar`.

`polartorect` translates complex values of the form  $r e^{i t}$  to the form  $a + b i$ . See `polartorect`.

`demo ("fft")` displays a demonstration of the `fft` package.

### **fortindent**

Variable

Default value: 0

`fortindent` controls the left margin indentation of expressions printed out by the `fortran` command. 0 gives normal printout (i.e., 6 spaces), and positive values will cause the expressions to be printed farther to the right.

### **fortran** (*expr*)

Function

Prints *expr* as a Fortran statement. The output line is indented with spaces. If the line is too long, `fortran` prints continuation lines. `fortran` prints the exponentiation operator  $\wedge$  as `**`, and prints a complex number  $a + b i$  in the form `(a,b)`.

*expr* may be an equation. If so, `fortran` prints an assignment statement, assigning the right-hand side of the equation to the left-hand side. In particular, if the right-hand side of *expr* is the name of a matrix, then `fortran` prints an assignment statement for each element of the matrix.

If *expr* is not something recognized by `fortran`, the expression is printed in `grind` format without complaint. `fortran` does not know about lists, arrays, or functions.

`fortindent` controls the left margin of the printed lines. 0 is the normal margin (i.e., indented 6 spaces). Increasing `fortindent` causes expressions to be printed further to the right.

When `fortspaces` is `true`, `fortran` fills out each printed line with spaces to 80 columns.

`fortran` evaluates its arguments; quoting an argument defeats evaluation. `fortran` always returns `done`.

Examples:

```
(%i1) expr: (a + b)^12$
(%i2) fortran (expr);
      (b+a)**12
(%o2)                                     done
(%i3) fortran ('x=expr);
      x = (b+a)**12
(%o3)                                     done
(%i4) fortran ('x=expand (expr));
      x = b**12+12*a*b**11+66*a**2*b**10+220*a**3*b**9+495*a**4*b**8+792
1      *a**5*b**7+924*a**6*b**6+792*a**7*b**5+495*a**8*b**4+220*a**9*b
2      **3+66*a**10*b**2+12*a**11*b+a**12
(%o4)                                     done
(%i5) fortran ('x=7+5%i);
      x = (7,5)
(%o5)                                     done
(%i6) fortran ('x=[1,2,3,4]);
      x = [1,2,3,4]
(%o6)                                     done
```

```
(%i7) f(x) := x^2$
(%i8) fortran (f);
      f
(%o8)                               done
```

**fortspaces**

Variable

Default value: `false`

When `fortspaces` is `true`, `fortran` fills out each printed line with spaces to 80 columns.

**horner** (*expr*, *x*)

Function

**horner** (*expr*)

Function

Returns a rearranged representation of *expr* as in Horner's rule, using *x* as the main variable if it is specified. *x* may be omitted in which case the main variable of the canonical rational expression form of *expr* is used.

`horner` sometimes improves stability if *expr* is to be numerically evaluated. It is also useful if Maxima is used to generate programs to be run in Fortran. See also `stringout`.

```
(%i1) expr: 1e-155*x^2 - 5.5*x + 5.2e155;
      2
(%o1)          1.0E-155 x  - 5.5 x + 5.2E+155
(%i2) expr2: horner (% , x), keepfloat: true;
(%o2)          (1.0E-155 x - 5.5) x + 5.2E+155
(%i3) ev (expr, x=1e155);
Maxima encountered a Lisp error:

floating point overflow

Automatically continuing.
To reenable the Lisp debugger set *debugger-hook* to nil.
(%i4) ev (expr2, x=1e155);
(%o4)          7.0E+154
```

**interpolate** (*f(x)*, *x*, *a*, *b*)

Function

**interpolate** (*f*, *a*, *b*)

Function

Finds the zero of function *f* as variable *x* varies over the range [*a*, *b*]. The function must have a different sign at each endpoint. If this condition is not met, the action of the of the function is governed by `intpolerror`. If `intpolerror` is `true` then an error occurs, otherwise the value of `intpolerror` is returned (thus for plotting `intpolerror` might be set to 0.0). Otherwise (given that Maxima can evaluate the first argument in the specified range, and that it is continuous) `interpolate` is guaranteed to come up with the zero (or one of them if there is more than one zero). The accuracy of `interpolate` is governed by `intpolabs` and `intpolrel` which must be non-negative floating point numbers. `interpolate` will stop when the first arg evaluates to something less than or equal to `intpolabs` or if successive approximants to the root differ by no more than `intpolrel * <one of the approximants>`. The default values of `intpolabs` and `intpolrel` are 0.0 so `interpolate` gets as good an

answer as is possible with the single precision arithmetic we have. The first arg may be an equation. The order of the last two args is irrelevant. Thus

```
interpolate (sin(x) = x/2, x, %pi, 0.1);
```

is equivalent to

```
interpolate (sin(x) = x/2, x, 0.1, %pi);
```

The method used is a binary search in the range specified by the last two args. When it thinks the function is close enough to being linear, it starts using linear interpolation.

```
(%i1) f(x) := sin(x) - x/2;
(%o1)          f(x) := sin(x) -  $\frac{x}{2}$ 
(%i2) interpolate (sin(x) - x/2, x, 0.1, %pi);
(%o2)          1.895494267033981
(%i3) interpolate (sin(x) = x/2, x, 0.1, %pi);
(%o3)          1.895494267033981
(%i4) interpolate (f(x), x, 0.1, %pi);
(%o4)          1.895494267033981
(%i5) interpolate (f, 0.1, %pi);
(%o5)          1.895494267033981
```

There is also a Newton method interpolation routine. See `newton`.

### **intpolabs**

Variable

Default value: 0.0

`intpolabs` is the accuracy of the `interpolate` command is governed by `intpolabs` and `intpolrel` which must be non-negative floating point numbers. `interpolate` will stop when the first arg evaluates to something less than or equal to `intpolabs` or if successive approximants to the root differ by no more than `intpolrel * <one of the approximants>`. The default values of `intpolabs` and `intpolrel` are 0.0 so `interpolate` gets as good an answer as is possible with the single precision arithmetic we have.

### **intpolerror**

Variable

Default value: `true`

`intpolerror` governs the behavior of `interpolate`. When `interpolate` is called, it determines whether or not the function to be interpolated satisfies the condition that the values of the function at the endpoints of the interpolation interval are opposite in sign. If they are of opposite sign, the interpolation proceeds. If they are of like sign, and `intpolerror` is `true`, then an error is signaled. If they are of like sign and `intpolerror` is not `true`, the value of `intpolerror` is returned. Thus for plotting, `intpolerror` might be set to 0.0.

### **intpolrel**

Variable

Default value: 0.0

`intpolrel` is the accuracy of the `interpolate` command is governed by `intpolabs` and `intpolrel` which must be non-negative floating point numbers. `interpolate`

will stop when the first arg evaluates to something less than or equal to `intpolabs` or if successive approximants to the root differ by no more than `intpolrel * <one of the approximants>`. The default values of `intpolabs` and `intpolrel` are 0.0 so `interpolate` gets as good an answer as is possible with the single precision arithmetic we have.

**newton** (*expr*, *x*, *x\_0*, *eps*) Function

Interpolation by Newton's method. `load ("newton1")` loads this function.

`newton` can handle some expressions that `interpolate` refuses to handle, since `interpolate` requires that everything evaluate to a floating point number. Thus `newton (x^2 - a^2, x, a/2, a^2/100)` complains that it can't tell if  $6.098490481853958E-4 a^2 < a^2/100$ . After `assume (a>0)`, the same function call succeeds, yielding a symbolic result,  $1.00030487804878 a$ .

On the other hand, `interpolate (x^2 - a^2, x, a/2, 2*a)` complains that  $0.5 a$  is not a floating point number.

An adaptive integrator which uses the Newton-Cotes 8 panel quadrature rule is available. See `qq`.

## 24.4 Definitions for Fourier Series

**equalp** (*x*, *y*) Function

Returns `true` if `equal (x, y)` otherwise `false` (doesn't give an error message like `equal (x, y)` would do in this case).

**remfun** (*f*, *expr*) Function

**remfun** (*f*, *expr*, *x*) Function

`remfun (f, expr)` replaces all occurrences of *f* (*arg*) by *arg* in *expr*.

`remfun (f, expr, x)` replaces all occurrences of *f* (*arg*) by *arg* in *expr* only if *arg* contains the variable *x*.

**funp** (*f*, *expr*) Function

**funp** (*f*, *expr*, *x*) Function

`funp (f, expr)` returns `true` if *expr* contains the function *f*.

`funp (f, expr, x)` returns `true` if *expr* contains the function *f* and the variable *x* is somewhere in the argument of one of the instances of *f*.

**absint** (*f*, *x*, *halfplane*) Function

**absint** (*f*, *x*) Function

**absint** (*f*, *x*, *a*, *b*) Function

`absint (f, x, halfplane)` returns the indefinite integral of *f* with respect to *x* in the given halfplane (`pos`, `neg`, or `both`). *f* may contain expressions of the form `abs (x)`, `abs (sin (x))`, `abs (a) * exp (-abs (b) * abs (x))`.

`absint (f, x)` is equivalent to `absint (f, x, pos)`.

`absint (f, x, a, b)` returns the definite integral of *f* with respect to *x* from *a* to *b*. *f* may include absolute values.

- fourier** ( $f, x, p$ ) Function  
Returns a list of the Fourier coefficients of  $f(x)$  defined on the interval  $[-\pi, \pi]$ .
- foursimp** ( $l$ ) Function  
Simplifies  $\sin(n\pi)$  to 0 if `sinnpiflag` is `true` and  $\cos(n\pi)$  to  $(-1)^n$  if `cosnpiflag` is `true`.
- sinnpiflag** Variable  
Default value: `true`  
See `foursimp`.
- cosnpiflag** Variable  
Default value: `true`  
See `foursimp`.
- fourexpend** ( $l, x, p, limit$ ) Function  
Constructs and returns the Fourier series from the list of Fourier coefficients  $l$  up through  $limit$  terms ( $limit$  may be `inf`).  $x$  and  $p$  have same meaning as in `fourier`.
- fourcos** ( $f, x, p$ ) Function  
Returns the Fourier cosine coefficients for  $f(x)$  defined on  $[0, \pi]$ .
- foursin** ( $f, x, p$ ) Function  
Returns the Fourier sine coefficients for  $f(x)$  defined on  $[0, \pi]$ .
- totalfourier** ( $f, x, p$ ) Function  
Returns `fourexpend(foursimp(fourier(f, x, p)), x, p, 'inf')`.
- fourint** ( $f, x$ ) Function  
Constructs and returns a list of the Fourier integral coefficients of  $f(x)$  defined on  $[\min, \max]$ .
- fourintcos** ( $f, x$ ) Function  
Returns the Fourier cosine integral coefficients for  $f(x)$  on  $[0, \max]$ .
- fourintsin** ( $f, x$ ) Function  
Returns the Fourier sine integral coefficients for  $f(x)$  on  $[0, \max]$ .



## 25 Statistics

### 25.1 Definitions for Statistics

**gauss** (*mean*, *sd*) Function  
Returns a random floating point number from a normal distribution with mean *mean* and standard deviation *sd*.





## 26 Arrays and Tables

### 26.1 Definitions for Arrays and Tables

**array** (*name*, *dim\_1*, ..., *dim\_n*) Function  
**array** (*name*, *type*, *dim\_1*, ..., *dim\_n*) Function  
**array** (*[name\_1, ..., name\_m]*, *dim\_1*, ..., *dim\_n*) Function

Creates an *n*-dimensional array. *n* may be less than or equal to 5. The subscripts for the *i*'th dimension are the integers running from 0 to *dim\_i*.

**array** (*name*, *dim\_1*, ..., *dim\_n*) creates a general array.

**array** (*name*, *type*, *dim\_1*, ..., *dim\_n*) creates an array, with elements of a specified type. *type* can be **fixnum** for integers of limited size or **flonum** for floating-point numbers.

**array** (*[name\_1, ..., name\_m]*, *dim\_1*, ..., *dim\_n*) creates *m* arrays, all of the same dimensions.

If the user assigns to a subscripted variable before declaring the corresponding array, an undeclared array is created. Undeclared arrays, otherwise known as hashed arrays (because hash coding is done on the subscripts), are more general than declared arrays. The user does not declare their maximum size, and they grow dynamically by hashing as more elements are assigned values. The subscripts of undeclared arrays need not even be numbers. However, unless an array is rather sparse, it is probably more efficient to declare it when possible than to leave it undeclared. The **array** function can be used to transform an undeclared array into a declared array.

**arrayapply** (*A*, [*i\_1*, ..., *i\_n*]) Function  
 Evaluates *A* [*i\_1*, ..., *i\_n*], where *A* is an array and *i\_1*, ..., *i\_n* are integers.

This is reminiscent of **apply**, except the first argument is an array instead of a function.

**arrayinfo** (*A*) Function  
 Returns a list of information about the array *A*. For hashed arrays it returns a list of **hashed**, the number of subscripts, and the subscripts of every element which has a value. For declared arrays it returns a list of **declared**, the number of subscripts, and the bounds that were given the the **array** function when it was called on *A*. Do **example(arrayinfo)**; for an example.

**arraymake** (*name*, [*i\_1*, ..., *i\_n*]) Function  
 Returns the expression *name* [*i\_1*, ..., *i\_n*].

This is reminiscent of **funmake**, except the return value is an unevaluated array reference instead of an unevaluated function call.

**arrays** system variable

Default value: []

`arrays` is a list of all the arrays that have been allocated, both declared and undeclared.

See also `array`, `arrayapply`, `arrayinfo`, `arraymake`, `fillarray`, `listarray`, and `rearray`.

**bashindices** (*expr*) Function

Transforms the expression *expr* by giving each summation and product a unique index. This gives `changevar` greater precision when it is working with summations or products. The form of the unique index is *jnumber*. The quantity *number* is determined by referring to `gensumnum`, which can be changed by the user. For example, `gensumnum:0$` resets it.

**fillarray** (*A*, *B*) Function

Fills array *A* from *B*, which is a list or an array.

If *A* is a floating-point (integer) array then *B* should be either a list of floating-point (integer) numbers or another floating-point (integer) array.

If the dimensions of the arrays are different *A* is filled in row-major order. If there are not enough elements in *B* the last element is used to fill out the rest of *A*. If there are too many the remaining ones are thrown away.

`fillarray` returns its first argument.

**getchar** (*a*, *i*) Function

Returns the *i*'th character of the quoted string or atomic name *a*. This function is useful in manipulating the `labels` list.

**listarray** (*A*) Function

Returns a list of the elements of a declared or hashed array *A*. The order is row-major. Elements which are not yet defined are represented by #####.

**make\_array** (*type*, *dim\_1*, ..., *dim\_n*) Function

Creates and returns a Lisp array. *type* may be `any`, `flonum`, `fixnum`, `hashed` or `functional`. There are *n* indices, and the *i*'th index runs from 0 to *dim<sub>i</sub>* - 1.

The advantage of `make_array` over `array` is that the return value doesn't have a name, and once a pointer to it goes away, it will also go away. For example, if `y: make_array (...)` then `y` points to an object which takes up space, but after `y: false`, `y` no longer points to that object, so the object can be garbage collected.

`y: make_array ('functional, 'f, 'hashed, 1)` - the second argument to `make_array` in this case is the function to call to calculate array elements, and the rest of the arguments are passed recursively to `make_array` to generate the "memory" for the array function object.

**rearray** (*A, dim\_1, ..., dim\_n*) Function

Changes the dimensions of an array. The new array will be filled with the elements of the old one in row-major order. If the old array was too small, the remaining elements are filled with **false**, 0.0 or 0, depending on the type of the array. The type of the array cannot be changed.

**remarray** (*A\_1, ..., A\_n*) Function

**remarray** (*all*) Function

Removes arrays and array associated functions and frees the storage occupied.

**remarray** (**all**) removes all items in the global list **arrays**.

It may be necessary to use this function if it is desired to redefine the values in a hashed array.

**remarray** returns the list of arrays removed.

**use\_fast\_arrays** option variable

- if **true** then only two types of arrays are recognized.

1) The art-q array (t in Common Lisp) which may have several dimensions indexed by integers, and may hold any Lisp or Maxima object as an entry. To construct such an array, enter **a:make\_array(any,3,4)**; then **a** will have as value, an array with twelve slots, and the indexing is zero based.

2) The Hash\_table array which is the default type of array created if one does **b[x+1]:y^2** (and **b** is not already an array, a list, or a matrix – if it were one of these an error would be caused since **x+1** would not be a valid subscript for an art-q array, a list or a matrix). Its indices (also known as keys) may be any object. It only takes one key at a time (**b[x+1,u]:y** would ignore the **u**). Referencing is done by **b[x+1] ==> y^2**. Of course the key may be a list, e.g. **b[[x+1,u]]:y** would be valid. This is incompatible with the old Maxima hash arrays, but saves consing.

An advantage of storing the arrays as values of the symbol is that the usual conventions about local variables of a function apply to arrays as well. The Hash\_table type also uses less consing and is more efficient than the old type of Maxima hashar. To obtain consistent behaviour in translated and compiled code set **translate\_fast\_arrays** to be **true**.



## 27 Matrices and Linear Algebra

### 27.1 Introduction to Matrices and Linear Algebra

#### 27.1.1 Dot

The operator `.` represents noncommutative multiplication and scalar product. When the operands are 1-column or 1-row matrices `a` and `b`, the expression `a.b` is equivalent to `sum(a[i]*b[i], i, 1, length(a))`. If `a` and `b` are not complex, this is the scalar product, also called the inner product or dot product, of `a` and `b`. The scalar product is defined as `conjugate(a).b` when `a` and `b` are complex; `innerproduct` in the `eigen` package provides the complex scalar product.

When the operands are more general matrices, the product is the matrix product `a` and `b`. The number of rows of `b` must equal the number of columns of `a`, and the result has number of rows equal to the number of rows of `a` and number of columns equal to the number of columns of `b`.

To distinguish `.` as an arithmetic operator from the decimal point in a floating point number, it may be necessary to leave spaces on either side. For example, `5.e3` is 5000.0 but `5 . e3` is 5 times `e3`.

There are several flags which govern the simplification of expressions involving `.`, namely `dot`, `dot0nscsimp`, `dot0simp`, `dot1simp`, `dotassoc`, `dotconstrules`, `dotdistrib`, `dotexptsimp`, `dotident`, and `dotscrules`.

#### 27.1.2 Vectors

`vect` is a package of functions for vector analysis. `load("vect")` loads this package, and `demo("vect")` displays a demonstration.

The vector analysis package can combine and simplify symbolic expressions including dot products and cross products, together with the gradient, divergence, curl, and Laplacian operators. The distribution of these operators over sums or products is governed by several flags, as are various other expansions, including expansion into components in any specific orthogonal coordinate systems. There are also functions for deriving the scalar or vector potential of a field.

The `vect` package contains these functions: `vectorsimp`, `scalefactors`, `express`, `potential`, and `vectorpotential`.

Warning: the `vect` package declares the dot operator `.` to be a commutative operator.

#### 27.1.3 eigen

The package `eigen` contains several functions devoted to the symbolic computation of eigenvalues and eigenvectors. Maxima loads the package automatically if one of the functions `eigenvalues` or `eigenvectors` is invoked. The package may be loaded explicitly as `load("eigen")`.

`demo ("eigen")` displays a demonstration of the capabilities of this package. `batch ("eigen")` executes the same demonstration, but without the user prompt between successive computations.

The functions in the `eigen` package are `conjugate`, `innerproduct`, `unitvector`, `columnvector`, `gramschmidt`, `eigenvalues`, `eigenvectors`, `uniteigenvectors`, and `similaritytransform`.

## 27.2 Definitions for Matrices and Linear Algebra

**addcol** ( $M$ ,  $list\_1$ , ...,  $list\_n$ ) Function

Appends the column(s) given by the one or more lists (or matrices) onto the matrix  $M$ .

**addrow** ( $M$ ,  $list\_1$ , ...,  $list\_n$ ) Function

Appends the row(s) given by the one or more lists (or matrices) onto the matrix  $M$ .

**adjoint** ( $M$ ) Function

Returns the adjoint of the matrix  $M$ .

**augcoefmatrix** ( $[eqn\_1, \dots, eqn\_m]$ ,  $[x\_1, \dots, x\_n]$ ) Function

Returns the augmented coefficient matrix for the variables  $x_1, \dots, x_n$  of the system of linear equations  $eqn_1, \dots, eqn_m$ . This is the coefficient matrix with a column adjoined for the constant terms in each equation (i.e., those terms not dependent upon  $x_1, \dots, x_n$ ).

```
(%i1) m: [2*x - (a - 1)*y = 5*b, c + b*y + a*x = 0]$
(%i2) augcoefmatrix (m, [x, y]);
          [ 2  1 - a  - 5 b ]
(%o2)    [
          [ a   b   c   ]
```

**charpoly** ( $M$ ,  $x$ ) Function

Returns the characteristic polynomial for the matrix  $M$  with respect to variable  $x$ . That is, `determinant (M - diagsmatrix (length (M), x))`.

```
(%i1) a: matrix ([3, 1], [2, 4]);
          [ 3  1 ]
(%o1)    [
          [ 2  4 ]
(%i2) expand (charpoly (a, lambda));
          2
(%o2)    lambda  - 7 lambda + 10
(%i3) (programmode: true, solve (%));
(%o3)    [lambda = 5, lambda = 2]
(%i4) matrix ([x1], [x2]);
          [ x1 ]
(%o4)    [
          [ x2 ]
```

```
(%i5) ev (a . % - lambda*%, %th(2)[1]);
                                [ x2 - 2 x1 ]
(%o5)                                [          ]
                                [ 2 x1 - x2 ]
(%i6) %[1, 1] = 0;
(%o6)                                x2 - 2 x1 = 0
(%i7) x2^2 + x1^2 = 1;
                                2      2
(%o7)                                x2  + x1  = 1
(%i8) solve ([%th(2), %], [x1, x2]);
                                1          2
(%o8) [[x1 = - ----, x2 = - ----],
                                sqrt(5)    sqrt(5)
```

$$\left[ x1 = \frac{1}{\sqrt{5}}, x2 = \frac{2}{\sqrt{5}} \right]$$

**coefmatrix** ( $[eqn_1, \dots, eqn_m], [x_1, \dots, x_n]$ ) Function  
 Returns the coefficient matrix for the variables  $eqn_1, \dots, eqn_m$  of the system of linear equations  $x_1, \dots, x_n$ .

**col** ( $M, i$ ) Function  
 Returns the  $i$ 'th column of the matrix  $M$ . The return value is a matrix.

**columnvector** ( $L$ ) Function

**covect** ( $L$ ) Function

Returns a matrix of one column and `length` ( $L$ ) rows, containing the elements of the list  $L$ .

`covect` is a synonym for `columnvector`.

`load ("eigen")` loads this function.

This is useful if you want to use parts of the outputs of the functions in this package in matrix calculations.

Example:

```
(%i1) load ("eigen")$
Warning - you are redefining the Macsyma function eigenvalues
Warning - you are redefining the Macsyma function eigenvectors
(%i2) columnvector ([aa, bb, cc, dd]);
                                [ aa ]
                                [      ]
                                [ bb ]
(%o2)                                [      ]
                                [ cc ]
                                [      ]
                                [ dd ]
```



**conjugate** ( $x$ ) Function  
**conj** ( $x$ ) Function

Substitutes  $\overline{\%i}$  for  $\%i$  in  $x$ . Depending on  $x$ , the result may be the complex conjugate of  $x$ .

`conj` is a synonym for `conjugate`.

`load ("eigen")` loads this function.

**copymatrix** ( $M$ ) Function

Returns a copy of the matrix  $M$ . This is the only way to make a copy aside from copying  $M$  element by element.

Note that an assignment of one matrix to another, as in `m2: m1`, does not copy `m1`. An assignment `m2 [i,j]: x` or `setelmx (x, i, j, m2)` also modifies `m1 [i,j]`. Creating a copy with `copymatrix` and then using assignment creates a separate, modified copy.

**determinant** ( $M$ ) Function

Computes the determinant of  $M$  by a method similar to Gaussian elimination.

The form of the result depends upon the setting of the switch `ratmx`.

There is a special routine for computing sparse determinants which is called when the switches `ratmx` and `sparse` are both `true`.

**detout** Variable

Default value: `false`

When `detout` is `true`, the determinant of a matrix whose inverse is computed is factored out of the inverse.

For this switch to have an effect `DOALLMXOPS` and `DOSCMXOPS` should be `false` (see their descriptions). Alternatively this switch can be given to `EV` which causes the other two to be set correctly.

Example:

```
(%i1) m: matrix ([a, b], [c, d]);
                                [ a  b ]
(%o1)                                [   ]
                                [ c  d ]

(%i2) detout: true$
(%i3) doallmxops: false$
(%i4) doscmxops: false$
(%i5) invert (m);
                                [ d  - b ]
                                [   ]
                                [ - c  a ]
(%o5) -----
                                a d - b c
```

**diagmatrix** ( $n, x$ ) Function

Returns a diagonal matrix of size  $n$  by  $n$  with the diagonal elements all equal to  $x$ . `diagmatrix (n, 1)` returns an identity matrix (same as `ident (n)`).

$n$  must evaluate to an integer, otherwise `diagmatrix` complains with an error message.  $x$  can be any kind of expression, including another matrix. If  $x$  is a matrix, it is not copied; all diagonal elements refer to the same instance,  $x$ .

**doallmxops**

Variable

Default value: `true`

When `doallmxops` is `true`, all operations relating to matrices are carried out. When it is `false` then the setting of the individual dot switches govern which operations are performed.

**domxexpt**

Variable

Default value: `true`

When `domxexpt` is `true`, a matrix exponential, `exp(M)` where  $M$  is a matrix, is interpreted as a matrix with element  $[i,j]$  equal to `exp(m[i,j])`. Otherwise `exp(M)` evaluates to `exp(ev(M))`.

`domxexpt` affects all expressions of the form  $base^{power}$  where  $base$  is an expression assumed scalar or constant, and  $power$  is a list or matrix.

Example:

```
(%i1) m: matrix ([1, %i], [a+b, %pi]);
              [ 1   %i ]
(%o1)          [          ]
              [ b + a %pi ]

(%i2) domxexpt: false$
(%i3) (1 - c)^m;
              [ 1   %i ]
              [          ]
              [ b + a %pi ]

(%o3)          (1 - c)
(%i4) domxexpt: true$
(%i5) (1 - c)^m;
              [          %i ]
              [ 1 - c   (1 - c) ]
(%o5)          [          ]
              [          b + a   %pi ]
              [ (1 - c)   (1 - c) ]
```

**domxmxops**

Variable

Default value: `true`

When `domxmxops` is `true`, all matrix-matrix or matrix-list operations are carried out (but not scalar-matrix operations); if this switch is `false` such operations are not carried out.

**domxnctimes**

Variable

Default value: `false`

When `domxnctimes` is `true`, non-commutative products of matrices are carried out.

- dontfactor** Variable  
 Default value: []  
**dontfactor** may be set to a list of variables with respect to which factoring is not to occur. (The list is initially empty.) Factoring also will not take place with respect to any variables which are less important, according the variable ordering assumed for canonical rational expression (CRE) form, than those on the **dontfactor** list.
- doscmxops** Variable  
 Default value: `false`  
 When **doscmxops** is `true`, scalar-matrix operations are carried out.
- doscmxplus** Variable  
 Default value: `false`  
 When **doscmxplus** is `true`, scalar-matrix operations yield a matrix result. This switch is not subsumed under **doallmxops**.
- dot0nscsimp** Variable  
 Default value: `true`  
 When **dot0nscsimp** is `true`, a non-commutative product of zero and a nonscalar term is simplified to a commutative product.
- dot0simp** Variable  
 Default value: `true`  
 When **dot0simp** is `true`, a non-commutative product of zero and a scalar term is simplified to a commutative product.
- dot1simp** Variable  
 Default value: `true`  
 When **dot1simp** is `true`, a non-commutative product of one and another term is simplified to a commutative product.
- dotassoc** Variable  
 Default value: `true`  
 When **dotassoc** is `true`, an expression  $(A.B).C$  simplifies to  $A.(B.C)$ .
- dotconstrules** Variable  
 Default value: `true`  
 When **dotconstrules** is `true`, a non-commutative product of a constant and another term is simplified to a commutative product. Turning on this flag effectively turns on **dot0simp**, **dot0nscsimp**, and **dot1simp** as well.
- dotdistrib** Variable  
 Default value: `false`  
 When **dotdistrib** is `true`, an expression  $A.(B + C)$  simplifies to  $A.B + A.C$ .

**dotexptsimp**

Variable

Default value: `true`When `dotexptsimp` is `true`, an expression `A.A` simplifies to `A2`.**dotident**

Variable

Default value: `1``dotident` is the value returned by `X0`.**dotscrules**

Variable

Default value: `false`When `dotscrules` is `true`, an expression `A.SC` or `SC.A` simplifies to `SC*A` and `A.(SC*B)` simplifies to `SC*(A.B)`.**echelon** (*M*)

Function

Returns the echelon form of the matrix *M*. The echelon form is computed from *M* by elementary row operations such that the first non-zero element in each row in the resulting matrix is a one and the column elements under the first one in each row are all zero.

```
(%i1) m: matrix ([2, 1-a, -5*b], [a, b, c]);
          [ 2  1 - a  - 5 b ]
(%o1)      [                ]
          [ a    b    c    ]
(%i2) echelon (m);
          [      a - 1      5 b      ]
          [ 1 - ----- - ---- ]
          [      2          2      ]
(%o2)/R/  [                ]
          [      2 c + 5 a b ]
          [ 0    1  ----- ]
          [                2 ]
          [      2 b + a  - a ]
```

**eigenvalues** (*M*)

Function

**eivals** (*M*)

Function

Returns a list of two lists containing the eigenvalues of the matrix *M*. The first sublist of the return value is the list of eigenvalues of the matrix, and the second sublist is the list of the multiplicities of the eigenvalues in the corresponding order.`eivals` is a synonym for `eigenvalues`.`eigenvalues` calls the function `solve` to find the roots of the characteristic polynomial of the matrix. Sometimes `solve` may not be able to find the roots of the polynomial; in that case some other functions in this package (except `conjugate`, `innerproduct`, `unitvector`, `columnvector` and `gramschmidt`) will not work.In some cases the eigenvalues found by `solve` may be complicated expressions. (This may happen when `solve` returns a not-so-obviously real expression for an eigenvalue which is known to be real.) It may be possible to simplify the eigenvalues using some other functions.

The package `eigen.mac` is loaded automatically when `eigenvalues` or `eigenvectors` is referenced. If `eigen.mac` is not already loaded, `load("eigen")` loads it. After loading, all functions and variables in the package are available.

**eigenvectors** ( $M$ ) Function  
**eivects** ( $M$ ) Function

takes a matrix  $M$  as its argument and returns a list of lists the first sublist of which is the output of `eigenvalues` and the other sublists of which are the eigenvectors of the matrix corresponding to those eigenvalues respectively. The calculated eigenvectors and the unit eigenvectors of the matrix are the right eigenvectors and the right unit eigenvectors respectively.

`eivects` is a synonym for `eigenvectors`.

The package `eigen.mac` is loaded automatically when `eigenvalues` or `eigenvectors` is referenced. If `eigen.mac` is not already loaded, `load("eigen")` loads it. After loading, all functions and variables in the package are available.

The flags that affect this function are:

`nondiagonalizable` is set to `true` or `false` depending on whether the matrix is nondiagonalizable or diagonalizable after `eigenvectors` returns.

`hermitianmatrix` when `true`, causes the degenerate eigenvectors of the Hermitian matrix to be orthogonalized using the Gram-Schmidt algorithm.

`knowneigvals` when `true` causes the `eigen` package to assume the eigenvalues of the matrix are known to the user and stored under the global name `listeigvals`. `listeigvals` should be set to a list similar to the output `eigenvalues`.

The function `algsys` is used here to solve for the eigenvectors. Sometimes if the eigenvalues are messy, `algsys` may not be able to find a solution. In some cases, it may be possible to simplify the eigenvalues by first finding them using `eigenvalues` command and then using other functions to reduce them to something simpler. Following simplification, `eigenvectors` can be called again with the `knowneigvals` flag set to `true`.

**ematrix** ( $m, n, x, i, j$ ) Function  
 Returns an  $m$  by  $n$  matrix, all elements of which are zero except for the  $[i, j]$  element which is  $x$ .

**entermatrix** ( $m, n$ ) Function  
 Returns an  $m$  by  $n$  matrix, reading the elements interactively.

If  $n$  is equal to  $m$ , Maxima prompts for the type of the matrix (diagonal, symmetric, antisymmetric, or general) and for each element. Each response is terminated by a semicolon ; or dollar sign \$.

If  $n$  is not equal to  $m$ , Maxima prompts for each element.

The elements may be any expressions, which are evaluated. `entermatrix` evaluates its arguments.

```
(%i1) n: 3$
(%i2) m: entermatrix (n, n)$
```

Is the matrix 1. Diagonal 2. Symmetric 3. Antisymmetric 4. General

Answer 1, 2, 3 or 4 :

1\$

Row 1 Column 1:

$(a+b)^n$

Row 2 Column 2:

$(a+b)^{(n+1)}$

Row 3 Column 3:

$(a+b)^{(n+2)}$

Matrix entered.

(%i3) m;

```

      [          3          ]
      [ (b + a)      0      0      ]
      [          ]
(%o3) [          4          ]
      [  0      (b + a)      0      ]
      [          ]
      [          ]
      [          5          ]
      [  0      0      (b + a)      ]

```

**genmatrix** (*a*, *i\_2*, *j\_2*, *i\_1*, *j\_1*)

Function

**genmatrix** (*a*, *i\_2*, *j\_2*, *i\_1*)

Function

**genmatrix** (*a*, *i\_2*, *j\_2*)

Function

Returns a matrix generated from *a*, taking element  $a[i_1, j_1]$  as the upper-left element and  $a[i_2, j_2]$  as the lower-right element of the matrix. Here *a* is an array (created by `array` but not by `make_array`) or an array function. (An array function is created like other functions with `:=` or `define`, but arguments are enclosed in square brackets instead of parentheses.)

If *j\_1* is omitted, it is assumed equal to *i\_1*. If both *j\_1* and *i\_1* are omitted, both are assumed equal to 1.

If a selected element *i*, *j* of the array is undefined, the matrix will contain a symbolic element  $a[i, j]$ .

(%i1)  $h[i, j] := 1/(i+j-1)$

(%i2) `genmatrix (h, 3, 3);`

```

      [  1  1 ]
      [ 1 - - ]
      [  2  3 ]
      [          ]
(%o2) [ 1 1 1 ]
      [ - - - ]
      [ 2 3 4 ]
      [          ]
      [ 1 1 1 ]
      [ - - - ]
      [ 3 4 5 ]

```

(%i3) `array (a, fixnum, 2, 2)`

```
(%i4) a[1,1]: %e$
(%i5) a[2,2]: %pi$
(%i6) kill (a[1,2], a[2,1])$
(%i7) genmatrix (a, 2, 2);

(%o7)
      [ %e      a      ]
      [          1, 2 ]
      [          ]
      [ a        %pi ]
      [ 2, 1      ]
```

**gramschmidt** (*x*)

Function

**gschmit** (*x*)

Function

Carries out the Gram-Schmidt orthogonalization algorithm on *x*, which is either a matrix or a list of lists. *x* is not modified by **gramschmidt**.

If *x* is a matrix, the algorithm is applied to the rows of *x*. If *x* is a list of lists, the algorithm is applied to the sublists, which must have equal numbers of elements. In either case, the return value is a list of lists, the sublists of which are orthogonal and span the same space as *x*. If the dimension of the span of *x* is less than the number of rows or sublists, some sublists of the return value are zero.

**factor** is called at each stage of the algorithm to simplify intermediate results. As a consequence, the return value may contain factored integers.

**gschmit** (note spelling) is a synonym for **gramschmidt**.

**load** ("eigen") loads this function.

Example:

```
(%i1) load ("eigen")$
Warning - you are redefining the Macsyma function eigenvalues
Warning - you are redefining the Macsyma function eigenvectors
(%i2) x: matrix ([1, 2, 3], [9, 18, 30], [12, 48, 60]);

(%o2)
      [ 1  2  3 ]
      [          ]
      [ 9  18 30 ]
      [          ]
      [ 12 48 60 ]

(%i3) y: gschmidt (x);

(%o3) [[1, 2, 3], [- 2/7, - 3/7, 5/7], [- 4/5, 3/5, 0]]

(%i4) i: innerproduct$
(%i5) [i (y[1], y[2]), i (y[2], y[3]), i (y[3], y[1])];
(%o5) [0, 0, 0]
```

**hach** (*a*, *b*, *m*, *n*, *l*)

Function

**hach** is an implementation of Hacijan's linear programming algorithm.

**load** ("kach") loads this function. **demo** ("kach") executes a demonstration of this function.

**ident** (*n*) Function  
Returns an *n* by *n* identity matrix.

**innerproduct** (*x*, *y*) Function

**inprod** (*x*, *y*) Function

Returns the inner product (also called the scalar product or dot product) of *x* and *y*, which are lists of equal length, or both 1-column or 1-row matrices of equal length. The return value is `conjugate (x) . y`, where `.` is the noncommutative multiplication operator.

`load ("eigen")` loads this function.

`inprod` is a synonym for `innerproduct`.

**invert** (*M*) Function

Returns the inverse of the matrix *M*. The inverse is computed by the adjoint method.

This allows a user to compute the inverse of a matrix with `bfloat` entries or polynomials with floating `pt.` coefficients without converting to `cre`-form.

Cofactors are computed by the `determinant` function, so if `ratmx` is `false` the inverse is computed without changing the representation of the elements.

The current implementation is inefficient for matrices of high order.

When `detout` is `true`, the determinant is factored out of the inverse.

The elements of the inverse are not automatically expanded. If *M* has polynomial elements, better appearing output can be generated by `expand (invert (m))`, `detout`. If it is desirable to then divide through by the determinant this can be accomplished by `xthru (%)` or alternatively from scratch by

```
expand (adjoint (m)) / expand (determinant (m))
invert (m) := adjoint (m) / determinant (m)
```

See `^^` (noncommutative exponent) for another method of inverting a matrix.

**lmxchar** Variable

Default value: `[`

`lmxchar` is the character displayed as the left delimiter of a matrix. See also `rmxchar`.

Example:

```
(%i1) lmxchar: "|"$
(%i2) matrix ([a, b, c], [d, e, f], [g, h, i]);
          | a  b  c ]
          |          ]
(%o2)    | d  e  f ]
          |          ]
          | g  h  i ]
```

**matrix** (*row\_1*, ..., *row\_n*) Function

Returns a rectangular matrix which has the rows *row\_1*, ..., *row\_n*. Each row is a list of expressions. All rows must be the same length.



The operations + (addition), - (subtraction), \* (multiplication), and / (division), are carried out element by element when the operands are two matrices, a scalar and a matrix, or a matrix and a scalar. The operation ^ (exponentiation, equivalently \*\*) is carried out element by element if the operands are a scalar and a matrix or a matrix and a scalar, but not if the operands are two matrices. All operations are normally carried out in full, including . (noncommutative multiplication).

Matrix multiplication is represented by the noncommutative multiplication operator .. The corresponding noncommutative exponentiation operator is ^^. For a matrix A, A.A = A^^2 and A^^-1 is the inverse of A, if it exists.

There are switches for controlling simplification of expressions involving dot and matrix-list operations. These are doallmxops, domxexpt domxmxops, doscmxops, and doscmxplus.

There are additional options which are related to matrices. These are: lmxchar, rmxchar, ratmx, listarith, detout, scalarmatrix, and sparse.

There are a number of functions which take matrices as arguments or yield matrices as return values. See eigenvalues, eigenvectors, determinant, charpoly, genmatrix, addcol, addrow, copymatrix, transpose, echelon, and rank.

Examples:

- Construction of matrices from lists.

```
(%i1) x: matrix ([17, 3], [-8, 11]);
      [ 17  3 ]
(%o1)  [      ]
      [ - 8  11 ]
(%i2) y: matrix ([%pi, %e], [a, b]);
      [ %pi %e ]
(%o2)  [      ]
      [ a  b ]
```

- Addition, element by element.

```
(%i3) x + y;
      [ %pi + 17 %e + 3 ]
(%o3)  [      ]
      [ a - 8    b + 11 ]
```

- Subtraction, element by element.

```
(%i4) x - y;
      [ 17 - %pi  3 - %e ]
(%o4)  [      ]
      [ - a - 8   11 - b ]
```

- Multiplication, element by element.

```
(%i5) x * y;
      [ 17 %pi  3 %e ]
(%o5)  [      ]
      [ - 8 a   11 b ]
```

- Division, element by element.

```
(%i6) x / y;
      [ 17      - 1 ]
```

$$\begin{array}{l}
 (\%o6) \quad \begin{bmatrix} \text{---} & 3 \text{ \%e} & ] \\
 \text{\%pi} & & ] \\
 [ & & ] \\
 [ 8 & 11 & ] \\
 [ - & - & -- ] \\
 [ a & b & ]
 \end{array}
 \end{array}$$

- Matrix to a scalar exponent, element by element.

$$\begin{array}{l}
 (\%i7) \quad x \wedge 3; \\
 (\%o7) \quad \begin{bmatrix} 4913 & 27 ] \\
 [ & ] \\
 [ - 512 & 1331 ]
 \end{array}
 \end{array}$$

- Scalar base to a matrix exponent, element by element.

$$\begin{array}{l}
 (\%i8) \quad \exp(y); \\
 (\%o8) \quad \begin{bmatrix} \text{\%pi} & \text{\%e} ] \\
 [ \text{\%e} & \text{\%e} ] \\
 [ & ] \\
 [ a & b ] \\
 [ \text{\%e} & \text{\%e} ]
 \end{array}
 \end{array}$$

- Matrix base to a matrix exponent. This is not carried out element by element.

$$\begin{array}{l}
 (\%i9) \quad x \wedge y; \\
 (\%o9) \quad \begin{bmatrix} \text{\%pi} & \text{\%e} ] \\
 [ & ] \\
 [ a & b ] \\
 [ 17 & 3 ] \\
 [ & ] \\
 [ - 8 & 11 ]
 \end{array}
 \end{array}$$

- Noncommutative matrix multiplication.

$$\begin{array}{l}
 (\%i10) \quad x \cdot y; \\
 (\%o10) \quad \begin{bmatrix} 3 a + 17 \text{\%pi} & 3 b + 17 \text{\%e} ] \\
 [ & ] \\
 [ 11 a - 8 \text{\%pi} & 11 b - 8 \text{\%e} ]
 \end{array} \\
 (\%i11) \quad y \cdot x; \\
 (\%o11) \quad \begin{bmatrix} 17 \text{\%pi} - 8 \text{\%e} & 3 \text{\%pi} + 11 \text{\%e} ] \\
 [ & ] \\
 [ 17 a - 8 b & 11 b + 3 a ]
 \end{array}
 \end{array}$$

- Noncommutative matrix exponentiation. A scalar base  $b$  to a matrix power  $M$  is carried out element by element and so  $b^M$  is the same as  $b^m$ .

$$\begin{array}{l}
 (\%i12) \quad x \wedge \wedge 3; \\
 (\%o12) \quad \begin{bmatrix} 3833 & 1719 ] \\
 [ & ] \\
 [ - 4584 & 395 ]
 \end{array} \\
 (\%i13) \quad \text{\%e} \wedge \wedge y; \\
 (\%o13) \quad \begin{bmatrix} \text{\%pi} & \text{\%e} ] \\
 [ \text{\%e} & \text{\%e} ] \\
 [ & ] \\
 [ a & b ] \\
 [ \text{\%e} & \text{\%e} ]
 \end{array}
 \end{array}$$

- A matrix raised to a -1 exponent with noncommutative exponentiation is the matrix inverse, if it exists.

```
(%i14) x ^^ -1;
          [ 11      3 ]
          [ --- - --- ]
          [ 211     211 ]
(%o14)    [          ]
          [ 8      17 ]
          [ ---   --- ]
          [ 211     211 ]

(%i15) x . (x ^^ -1);
          [ 1  0 ]
(%o15)    [    ]
          [ 0  1 ]
```

**matrixmap** (*f*, *M*) Function

Returns a matrix with element *i*,*j* equal to *f*(*M*[*i*,*j*]).

See also `map`, `fullmap`, `fullmap1`, and `apply`.

**matrixp** (*expr*) Function

Returns `true` if *expr* is a matrix, otherwise `false`.

**matrix\_element\_add** Variable

Default value: +

`matrix_element_add` is the operation invoked in place of addition in a matrix multiplication. `matrix_element_add` can be assigned any n-ary operator (that is, a function which handles any number of arguments). The assigned value may be the name of an operator enclosed in quote marks, the name of a function, or a lambda expression.

See also `matrix_element_mult` and `matrix_element_transpose`.

Example:

```
(%i1) matrix_element_add: "*"
(%i2) matrix_element_mult: "^"
(%i3) aa: matrix ([a, b, c], [d, e, f]);
          [ a  b  c ]
(%o3)    [          ]
          [ d  e  f ]
(%i4) bb: matrix ([u, v, w], [x, y, z]);
          [ u  v  w ]
(%o4)    [          ]
          [ x  y  z ]
(%i5) aa . transpose (bb);
          [ u  v  w  x  y  z ]
          [ a  b  c  a  b  c ]
(%o5)    [          ]
          [ u  v  w  x  y  z ]
          [ d  e  f  d  e  f ]
```

**matrix\_element\_mult**

Variable

Default value: \*

`matrix_element_mult` is the operation invoked in place of multiplication in a matrix multiplication. `matrix_element_mult` can be assigned any binary operator. The assigned value may be the name of an operator enclosed in quote marks, the name of a function, or a lambda expression.

The dot operator `.` is a useful choice in some contexts.

See also `matrix_element_add` and `matrix_element_transpose`.

Example:

```
(%i1) matrix_element_add: lambda ([[x]], sqrt (apply ("+", x)))$
(%i2) matrix_element_mult: lambda ([x, y], (x - y)^2)$
(%i3) [a, b, c] . [x, y, z];

(%o3)          2          2          2
      sqrt((c - z)  + (b - y)  + (a - x) )
(%i4) aa: matrix ([a, b, c], [d, e, f]);
      [ a  b  c ]
(%o4)          [          ]
      [ d  e  f ]
(%i5) bb: matrix ([u, v, w], [x, y, z]);
      [ u  v  w ]
(%o5)          [          ]
      [ x  y  z ]
(%i6) aa . transpose (bb);
      [          2          2          2 ]
      [ sqrt((c - w)  + (b - v)  + (a - u) ) ]
(%o6) Col 1 = [          ]
      [          2          2          2 ]
      [ sqrt((f - w)  + (e - v)  + (d - u) ) ]

      [          2          2          2 ]
      [ sqrt((c - z)  + (b - y)  + (a - x) ) ]
Col 2 = [          ]
      [          2          2          2 ]
      [ sqrt((f - z)  + (e - y)  + (d - x) ) ]
```

**matrix\_element\_transpose**

Variable

Default value: false

`matrix_element_transpose` is the operation applied to each element of a matrix when it is transposed. `matrix_element_mult` can be assigned any unary operator. The assigned value may be the name of an operator enclosed in quote marks, the name of a function, or a lambda expression.

When `matrix_element_transpose` equals `transpose`, the `transpose` function is applied to every element. When `matrix_element_transpose` equals `nonscalars`, the `transpose` function is applied to every nonscalar element. If some element is an atom, the `nonscalars` option applies `transpose` only if the atom is declared nonscalar, while the `transpose` option always applies `transpose`.

The default value, `false`, means no operation is applied.

See also `matrix_element_add` and `matrix_element_mult`.

Examples:

```
(%i1) declare (a, nonscalar)$
(%i2) transpose ([a, b]);
(%o2)          [ transpose(a) ]
          [          ]
          [          b          ]
(%i3) matrix_element_transpose: nonscalars$
(%i4) transpose ([a, b]);
(%o4)          [ transpose(a) ]
          [          ]
          [          b          ]
(%i5) matrix_element_transpose: transpose$
(%i6) transpose ([a, b]);
(%o6)          [ transpose(a) ]
          [          ]
          [ transpose(b) ]
(%i7) matrix_element_transpose: lambda ([x], realpart(x) - %i*imagpart(x))$
(%i8) m: matrix ([1 + 5*%i, 3 - 2*%i], [7*%i, 11]);
(%o8)          [ 5 %i + 1  3 - 2 %i ]
          [          ]
          [ 7 %i          11          ]
(%i9) transpose (m);
(%o9)          [ 1 - 5 %i  - 7 %i ]
          [          ]
          [ 2 %i + 3    11          ]
```

**mattrace** ( $M$ ) Function

Returns the trace (that is, the sum of the elements on the main diagonal) of the square matrix  $M$ .

`mattrace` is called by `ncharpoly`, an alternative to Maxima's `charpoly`.

`load ("nchrpl")` loads this function.

**minor** ( $M, i, j$ ) Function

Returns the  $i, j$  minor of the matrix  $M$ . That is,  $M$  with row  $i$  and column  $j$  removed.

**ncexpt** ( $a, b$ ) Function

If a non-commutative exponential expression is too wide to be displayed as  $a^b$  it appears as `ncexpt (a, b)`.

`ncexpt` is not the name of a function or operator; the name only appears in output, and is not recognized in input.

**ncharpoly** ( $M, x$ ) Function

Returns the characteristic polynomial of the matrix  $M$  with respect to  $x$ . This is an alternative to Maxima's `charpoly`.

`ncharpoly` works by computing traces of powers of the given matrix, which are known to be equal to sums of powers of the roots of the characteristic polynomial. From these quantities the symmetric functions of the roots can be calculated, which are nothing more than the coefficients of the characteristic polynomial. `charpoly` works by forming the determinant of  $x * \text{ident } [n] - a$ . Thus `ncharpoly` wins, for example, in the case of large dense matrices filled with integers, since it avoids polynomial arithmetic altogether.

`load ("nchrpl")` loads this file.

- newdet** ( $M, n$ ) Function  
 Computes the determinant of the matrix or array  $M$  by the Johnson-Gentleman tree minor algorithm. The argument  $n$  is the order; it is optional if  $M$  is a matrix.
- nonscalar** declaration  
 Makes atoms behave as does a list or matrix with respect to the dot operator.
- nonscalarp** ( $expr$ ) Function  
 Returns `true` if  $expr$  is a non-scalar, i.e., it contains atoms declared as non-scalars, lists, or matrices.
- permanent** ( $M, n$ ) Function  
 Computes the permanent of the matrix  $M$ . A permanent is like a determinant but with no sign changes.
- rank** ( $M$ ) Function  
 Computes the rank of the matrix  $M$ . That is, the order of the largest non-singular subdeterminant of  $M$ .  
*rank* may return the wrong answer if it cannot determine that a matrix element that is equivalent to zero is indeed so.
- ratmx** Variable  
 Default value: `false`  
 When `ratmx` is `false`, determinant and matrix addition, subtraction, and multiplication are performed in the representation of the matrix elements and cause the result of matrix inversion to be left in general representation.  
 When `ratmx` is `true`, the 4 operations mentioned above are performed in CRE form and the result of matrix inverse is in CRE form. Note that this may cause the elements to be expanded (depending on the setting of `ratfac`) which might not always be desired.
- row** ( $M, i$ ) Function  
 Returns the  $i$ 'th row of the matrix  $M$ . The return value is a matrix.

**scalarmatrixp**

Variable

Default value: true

When `scalarmatrixp` is true, then whenever a 1 x 1 matrix is produced as a result of computing the dot product of matrices it is simplified to a scalar, namely the sole element of the matrix.

When `scalarmatrixp` is all, then all 1 x 1 matrices are simplified to scalars.

When `scalarmatrixp` is false, 1 x 1 matrices are not simplified to scalars.

**scalefactors** (*coordinatetransform*)

Function

Here `coordinatetransform` evaluates to the form `[[expression1, expression2, ...], indeterminate1, indeterminate2, ...]`, where `indeterminate1`, `indeterminate2`, etc. are the curvilinear coordinate variables and where a set of rectangular Cartesian components is given in terms of the curvilinear coordinates by `[expression1, expression2, ...]`. `coordinates` is set to the vector `[indeterminate1, indeterminate2, ...]`, and `dimension` is set to the length of this vector. `SF[1]`, `SF[2]`, ..., `SF[DIMENSION]` are set to the coordinate scale factors, and `sfprod` is set to the product of these scale factors. Initially, `coordinates` is `[X, Y, Z]`, `dimension` is 3, and `SF[1]=SF[2]=SF[3]=SFPROD=1`, corresponding to 3-dimensional rectangular Cartesian coordinates. To expand an expression into physical components in the current coordinate system, there is a function with usage of the form

**setelmx** (*x, i, j, M*)

Function

Assigns `x` to the  $(i, j)$ 'th element of the matrix `M`, and returns the altered matrix.

`M [i, j]`: `x` has the same effect, but returns `x` instead of `M`.

**similaritytransform** (*M*)

Function

**simtran** (*M*)

Function

`similaritytransform` computes a similarity transform of the matrix `M`. It returns a list which is the output of the `uniteigenvectors` command. In addition if the flag `nondiagonalizable` is false two global matrices `leftmatrix` and `rightmatrix` are computed. These matrices have the property that `leftmatrix . M . rightmatrix` is a diagonal matrix with the eigenvalues of `M` on the diagonal. If `nondiagonalizable` is true the left and right matrices are not computed.

If the flag `hermitianmatrix` is true then `leftmatrix` is the complex conjugate of the transpose of `rightmatrix`. Otherwise `leftmatrix` is the inverse of `rightmatrix`. `rightmatrix` is the matrix the columns of which are the unit eigenvectors of `M`. The other flags (see `eigenvalues` and `eigenvectors`) have the same effects since `similaritytransform` calls the other functions in the package in order to be able to form `rightmatrix`.

`load ("eigen")` loads this function.

`simtran` is a synonym for `similaritytransform`.

**sparse**

Variable

Default value: false

When `sparse` is true, and if `ratmx` is true, then `determinant` will use special routines for computing sparse determinants.

- submatrix** ( $i_1, \dots, i_m, M, j_1, \dots, j_n$ ) Function  
**submatrix** ( $i_1, \dots, i_m, M$ ) Function  
**submatrix** ( $M, j_1, \dots, j_n$ ) Function  
 Returns a new matrix composed of the matrix  $M$  with rows  $i_1, \dots, i_m$  deleted, and columns  $j_1, \dots, j_n$  deleted.
- transpose** ( $M$ ) Function  
 Returns the transpose of  $M$ .  
 If  $M$  is a matrix, the return value is another matrix  $N$  such that  $N[i, j] = M[j, i]$ .  
 Otherwise  $M$  is a list, and the return value is a matrix  $N$  of length ( $m$ ) rows and 1 column, such that  $N[i, 1] = M[i]$ .
- triangularize** ( $M$ ) Function  
 Returns the upper triangular form of the matrix  $M$ .  
 $M$  need not be square.
- uniteigenvectors** ( $M$ ) Function  
**ueivects** ( $M$ ) Function  
 Computes unit eigenvectors of the matrix  $M$ . The return value is a list of lists, the first sublist of which is the output of the **eigenvalues** command, and the other sublists of which are the unit eigenvectors of the matrix corresponding to those eigenvalues respectively.  
 The flags mentioned in the description of the **eigenvectors** command have the same effects in this one as well.  
 When **knoweigvects** is **true**, the **eigen** package assumes that the eigenvectors of the matrix are known to the user and are stored under the global name **listeigvects**. **listeigvects** should be set to a list similar to the output of the **eigenvectors** command.  
 If **knoweigvects** is set to **true** and the list of eigenvectors is given the setting of the flag **nondiagonalizable** may not be correct. If that is the case please set it to the correct value. The author assumes that the user knows what he is doing and will not try to diagonalize a matrix the eigenvectors of which do not span the vector space of the appropriate dimension.  
**load** ("eigen") loads this function.  
**ueivects** is a synonym for **uniteigenvectors**.
- unitvector** ( $x$ ) Function  
**uvect** ( $x$ ) Function  
 Returns  $x/\text{norm}(x)$ ; this is a unit vector in the same direction as  $x$ .  
**load** ("eigen") loads this function.  
**uvect** is a synonym for **unitvector**.



**vectorsimp** (*expr*) Function

Applies simplifications and expansions according to the following global flags:

`expandall`, `expanddot`, `expanddotplus`, `expandcross`, `expandcrossplus`,  
`expandcrosscross`, `expandgrad`, `expandgradplus`, `expandgradprod`, `expanddiv`,  
`expanddivplus`, `expanddivprod`, `expandcurl`, `expandcurlplus`, `expandcurlcurl`,  
`expandlaplacian`, `expandlaplacianplus`, and `expandlaplacianprod`.

All these flags have default value `false`. The `plus` suffix refers to employing additivity or distributivity. The `prod` suffix refers to the expansion for an operand that is any kind of product.

`expandcrosscross`

Simplifies  $p(qr)$  to  $(p.r) * q - (p.q) * r$ .

`expandcurlcurl`

Simplifies  $curlcurlp$  to  $graddivp + divgradp$ .

`expandlaplaciantodivgrad`

Simplifies  $laplacianp$  to  $divgradp$ .

`expandcross`

Enables `expandcrossplus` and `expandcrosscross`.

`expandplus`

Enables `expanddotplus`, `expandcrossplus`, `expandgradplus`,  
`expanddivplus`, `expandcurlplus`, and `expandlaplacianplus`.

`expandprod`

Enables `expandgradprod`, `expanddivprod`, and `expandlaplacianprod`.

These flags have all been declared `evflag`.

**vect\_cross** Variable

Default value: `false`

When `vect_cross` is `true`, it allows `DIFF(X~Y,T)` to work where `~` is defined in `SHARE;VECT` (where `VECT_CROSS` is set to `true`, anyway.)

**zeromatrix** (*m*, *n*) Function

Returns an  $m$  by  $n$  matrix, all elements of which are zero.

**"["** special symbol

[ and ] mark the beginning and end, respectively, of a list.

[ and ] also enclose the subscripts of a list, array, hash array, or array function.

Examples:

```
(%i1) x: [a, b, c];
(%o1) [a, b, c]
(%i2) x[3];
(%o2) c
(%i3) array (y, fixnum, 3);
(%o3) y
```





## 28 Affine

### 28.1 Definitions for Affine

**fast\_linsolve** ( $[expr\_1, \dots, expr\_m], [x\_1, \dots, x\_n]$ ) Function

Solves the simultaneous linear equations  $expr\_1, \dots, expr\_m$  for the variables  $x\_1, \dots, x\_n$ . Each  $expr\_i$  may be an equation or a general expression; if given as a general expression, it is treated as an equation of the form  $expr\_i = 0$ .

The return value is a list of equations of the form  $[x\_1 = a\_1, \dots, x\_n = a\_n]$  where  $a\_1, \dots, a\_n$  are all free of  $x\_1, \dots, x\_n$ .

`fast_linsolve` is faster than `linsolve` for system of equations which are sparse.

**groebner\_basis** ( $[expr\_1, \dots, expr\_m]$ ) Function

Returns a Groebner basis for the equations  $expr\_1, \dots, expr\_m$ . The function `polysimp` can then be used to simplify other functions relative to the equations.

```
groebner_basis ([3*x^2+1, y*x])$
```

```
polysimp (y^2*x + x^3*9 + 2) ==> -3*x + 2
```

`polysimp(f)` yields 0 if and only if  $f$  is in the ideal generated by  $expr\_1, \dots, expr\_m$ , that is, if and only if  $f$  is a polynomial combination of the elements of  $expr\_1, \dots, expr\_m$ .

**set\_up\_dot\_simplifications** ( $eqns, check\_through\_degree$ ) Function

**set\_up\_dot\_simplifications** ( $eqns$ ) Function

The  $eqns$  are polynomial equations in non commutative variables. The value of `current_variables` is the list of variables used for computing degrees. The equations must be homogeneous, in order for the procedure to terminate.

If you have checked overlapping simplifications in `dot_simplifications` above the degree of  $f$ , then the following is true: `dotsimp(f)` yields 0 if and only if  $f$  is in the ideal generated by the equations, i.e., if and only if  $f$  is a polynomial combination of the elements of the equations.

The degree is that returned by `nc_degree`. This in turn is influenced by the weights of individual variables.

**declare\_weight** ( $x\_1, w\_1, \dots, x\_n, w\_n$ ) Function

Assigns weights  $w\_1, \dots, w\_n$  to  $x\_1, \dots, x\_n$ , respectively. These are the weights used in computing `nc_degree`.

**nc\_degree** ( $p$ ) Function

Returns the degree of a noncommutative polynomial  $p$ . See `declare_weights`.

**dotsimp** ( $f$ ) Function

Returns 0 if and only if  $f$  is in the ideal generated by the equations, i.e., if and only if  $f$  is a polynomial combination of the elements of the equations.

**fast\_central\_elements** ( $[x_1, \dots, x_n], n$ ) Function

If `set_up_dot_simplifications` has been previously done, finds the central polynomials in the variables  $x_1, \dots, x_n$  in the given degree,  $n$ .

For example:

```
set_up_dot_simplifications ([y.x + x.y], 3);
fast_central_elements ([x, y], 2);
[y.y, x.x];
```

**check\_overlaps** ( $n, add\_to\_simps$ ) Function

Checks the overlaps thru degree  $n$ , making sure that you have sufficient simplification rules in each degree, for `dotsimp` to work correctly. This process can be speeded up if you know before hand what the dimension of the space of monomials is. If it is of finite global dimension, then `hilbert` should be used. If you don't know the monomial dimensions, do not specify a `rank_function`. An optional third argument `reset, false` says don't bother to query about resetting things.

**mono** ( $[x_1, \dots, x_n], n$ ) Function

Returns the list of independent monomials relative to the current dot simplifications of degree  $n$  in the variables  $x_1, \dots, x_n$ .

**monomial\_dimensions** ( $n$ ) Function

Compute the Hilbert series through degree  $n$  for the current algebra.

**extract\_linear\_equations** ( $[p_1, \dots, p_n], [m_1, \dots, m_n]$ ) Function

Makes a list of the coefficients of the noncommutative polynomials  $p_1, \dots, p_n$  of the noncommutative monomials  $m_1, \dots, m_n$ . The coefficients should be scalars. Use `list_nc_monomials` to build the list of monomials.

**list\_nc\_monomials** ( $[p_1, \dots, p_n]$ ) Function

**list\_nc\_monomials** ( $p$ ) Function

Returns a list of the non commutative monomials occurring in a polynomial  $p$  or a list of polynomials  $p_1, \dots, p_n$ .

**create\_list** ( $form, x_1, list_1, \dots, x_n, list_n$ ) Function

Create a list by evaluating `form` with  $x_1$  bound to each element of `list_1`, and for each such binding bind  $x_2$  to each element of `list_2`, .... The number of elements in the result will be the product of the number of elements in each list. Each variable  $x_i$  must actually be a symbol—it will not be evaluated. The list arguments will be evaluated once at the beginning of the iteration.

```
(%i82) create_list1(x^i,i,[1,3,7]);
(%o82) [x,x^3,x^7]
```

With a double iteration:

```
(%i79) create_list([i,j],i,[a,b],j,[e,f,h]);
(%o79) [[a,e],[a,f],[a,h],[b,e],[b,f],[b,h]]
```

Instead of `list_i` two args may be supplied each of which should evaluate to a number. These will be the inclusive lower and upper bounds for the iteration.

```
(%i81) create_list([i,j],i,[1,2,3],j,1,i);  
(%o81) [[1,1],[2,1],[2,2],[3,1],[3,2],[3,3]]
```

Note that the limits or list for the  $j$  variable can depend on the current value of  $i$ .

**all\_dotsimp\_denoms**

Variable

Default value: `false`

When `all_dotsimp_denoms` is a list, the denominators encountered by `dotsimp` are appended to the list. `all_dotsimp_denoms` may be initialized to an empty list `[]` before calling `dotsimp`.

By default, denominators are not collected by `dotsimp`.



## 29 itensor

### 29.1 Introduction to itensor

Maxima implements symbolic tensor manipulation of two distinct types: component tensor manipulation (`ctensor` package) and indicial tensor manipulation (`itensor` package).

Nota bene: Please see the note on 'new tensor notation' below.

Component tensor manipulation means that geometrical tensor objects are represented as arrays or matrices. Tensor operations such as contraction or covariant differentiation are carried out by actually summing over repeated (dummy) indices with `do` statements. That is, one explicitly performs operations on the appropriate tensor components stored in an array or matrix.

Indicial tensor manipulation is implemented by representing tensors as functions of their covariant, contravariant and derivative indices. Tensor operations such as contraction or covariant differentiation are performed by manipulating the indices themselves rather than the components to which they correspond.

These two approaches to the treatment of differential, algebraic and analytic processes in the context of Riemannian geometry have various advantages and disadvantages which reveal themselves only through the particular nature and difficulty of the user's problem. However, one should keep in mind the following characteristics of the two implementations:

The representation of tensors and tensor operations explicitly in terms of their components makes `ctensor` easy to use. Specification of the metric and the computation of the induced tensors and invariants is straightforward. Although all of Maxima's powerful simplification capacity is at hand, a complex metric with intricate functional and coordinate dependencies can easily lead to expressions whose size is excessive and whose structure is hidden. In addition, many calculations involve intermediate expressions which swell causing programs to terminate before completion. Through experience, a user can avoid many of these difficulties.

Because of the special way in which tensors and tensor operations are represented in terms of symbolic operations on their indices, expressions which in the component representation would be unmanageable can sometimes be greatly simplified by using the special routines for symmetrical objects in `itensor`. In this way the structure of a large expression may be more transparent. On the other hand, because of the the special indicial representation in `itensor`, in some cases the user may find difficulty with the specification of the metric, function definition, and the evaluation of differentiated "indexed" objects.

#### 29.1.1 New tensor notation

Until now, the `itensor` package in Maxima has used a notation that sometimes led to incorrect index ordering. Consider the following, for instance:

```
(%i2) imetric(g);
(%o2) done
(%i3) ishow(g([], [j,k])*g([], [i,l])*a([i,j], []))$
          i l j k
(%t3)    g   g   a
```



```
(%i4) ishow(contract(%))$
                                     i j
                                     k l
(%t4)                                a
```

This result is incorrect unless  $a$  happens to be a symmetric tensor. The reason why this happens is that although `itensor` correctly maintains the order within the set of covariant and contravariant indices, once an index is raised or lowered, its position relative to the other set of indices is lost.

To avoid this problem, a new notation has been developed that remains fully compatible with the existing notation and can be used interchangeably. In this notation, contravariant indices are inserted in the appropriate positions in the covariant index list, but with a minus sign prepended. Functions like `contract` and `ishow` are now aware of this new index notation and can process tensors appropriately.

In this new notation, the previous example yields a correct result:

```
(%i5) ishow(g([-j,-k],[ ])*g([-i,-l],[ ])*a([i,j],[ ]))$
                                     i l      j k
(%t5)                                g  a      g
                                     i j
(%i6) ishow(contract(%))$
                                     l k
(%t6)                                a
```

Presently, the only code that makes use of this notation is the `lc2kdt` function. Through this notation, it achieves consistent results as it applies the metric tensor to resolve Levi-Civita symbols without resorting to numeric indices.

Since this code is brand new, it probably contains bugs. While it has been tested to make sure that it doesn't break anything using the "old" tensor notation, there is a considerable chance that "new" tensors will fail to interoperate with certain functions or features. These bugs will be fixed as they are encountered... until then, caveat emptor!

### 29.1.2 Indicial tensor manipulation

The indicial tensor manipulation package may be loaded by `load(itensor)`. Demos are also available: try `demo(tensor)`.

In `itensor` a tensor is represented as an "indexed object". This is a function of 3 groups of indices which represent the covariant, contravariant and derivative indices. The covariant indices are specified by a list as the first argument to the indexed object, and the contravariant indices by a list as the second argument. If the indexed object lacks either of these groups of indices then the empty list `[]` is given as the corresponding argument. Thus, `g([a,b],[c])` represents an indexed object called  $g$  which has two covariant indices  $(a,b)$ , one contravariant index  $(c)$  and no derivative indices.

The derivative indices, if they are present, are appended as additional arguments to the symbolic function representing the tensor. They can be explicitly specified by the user or be created in the process of differentiation with respect to some coordinate variable. Since ordinary differentiation is commutative, the derivative indices are sorted alphanumerically, unless `iframe_flag` is set to `true`, indicating that a frame metric is being used.. This canonical ordering makes it possible for Maxima to recognize that, for example,  $t([a],[b],i,j)$

is the same as  $t([a], [b], j, i)$ . Differentiation of an indexed object with respect to some coordinate whose index does not appear as an argument to the indexed object would normally yield zero. This is because Maxima would not know that the tensor represented by the indexed object might depend implicitly on the corresponding coordinate. By modifying the existing Maxima function `diff` in `itensor`, Maxima now assumes that all indexed objects depend on any variable of differentiation unless otherwise stated. This makes it possible for the summation convention to be extended to derivative indices. It should be noted that `itensor` does not possess the capabilities of raising derivative indices, and so they are always treated as covariant.

The following functions are available in the tensor package for manipulating indexed objects. At present, with respect to the simplification routines, it is assumed that indexed objects do not by default possess symmetry properties. This can be overridden by setting the variable `allsym[false]` to `true`, which will result in treating all indexed objects completely symmetric in their lists of covariant indices and symmetric in their lists of contravariant indices.

The `itensor` package generally treats tensors as opaque objects. Tensorial equations are manipulated based on algebraic rules, specifically symmetry and contraction rules. In addition, the `itensor` package understands covariant differentiation, curvature, and torsion. Calculations can be performed relative to a metric of moving frame, depending on the setting of the `iframe_flag` variable.

A sample session below demonstrates how to load the `itensor` package, specify the name of the metric, and perform some simple calculations.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) imetric(g);
(%o2)
done
(%i3) components(g([i,j],[ ]),p([i,j],[ ])*e([ ],[ ]))$
(%i4) ishow(g([k,l],[ ]))$
(%t4)
      e p
      k l

(%i5) ishow(diff(v([i],[ ]),t))$
(%t5)
      0
(%i6) depends(v,t);
(%o6)
[v(t)]
(%i7) ishow(diff(v([i],[ ]),t))$
(%t7)
      d
      -- (v )
      dt   i

(%i8) ishow(idiff(v([i],[ ]),j))$
(%t8)
      v
      i,j

(%i9) ishow(extdiff(v([i],[ ]),j))$
(%t9)
      v   - v
      j,i   i,j
      -----
      2

(%i10) ishow(liediff(v,w([i],[ ])))$
```

```

(%t10)

$$v_{i, \%3} w_{i, \%3} + v_{i, \%3} w_{i, \%3}$$

(%i11) ishow(covdiff(v([i],[ ]),j))$
(%t11)

$$v_{i,j} - v_{i,j}^{ic2}$$

(%i12) ishow(ev(%,ic2))$
(%t12)

$$v_{i,j} - g_{i,j} v_{j, \%4} (e_{j, \%5, i} p_{i, \%5, j} + e_{i, \%5, j} p_{j, \%5, i} - e_{i, \%5, j} p_{i, \%5, j} - e_{j, \%5, i} p_{j, \%5, i} + e_{i, \%5, j} p_{j, \%5, i} + e_{j, \%5, i} p_{i, \%5, j})/2$$

(%i13) iframe_flag:true;
(%o13)
true
(%i14) ishow(covdiff(v([i],[ ]),j))$
(%t14)

$$v_{i,j} - v_{i,j}^{ic2}$$

(%i15) ishow(ev(%,ic2))$
(%t15)

$$v_{i,j} - v_{i,j}^{ic2}$$

(%i16) ishow(radcan(ev(%,ifc2,ifc1)))$
(%t16)

$$- (ifg_{i,j, \%6} v_{j, \%8} ifb_{i, \%6} + ifg_{i,j, \%6} v_{i, \%6} ifb_{j, \%8} - 2 v_{i,j} - ifg_{i,j, \%6} v_{j, \%8} ifb_{i, \%6} - ifg_{i,j, \%6} v_{i, \%6} ifb_{j, \%8})/2$$

(%i17) ishow(canform(s([i,j],[ ])-s([j,i])))$
(%t17)

$$s_{i,j} - s_{j,i}$$

(%i18) decsym(s,2,0,[sym(all)], [ ]);
(%o18)
done
(%i19) ishow(canform(s([i,j],[ ])-s([j,i])))$
(%t19)
0
(%i20) ishow(canform(a([i,j],[ ])+a([j,i])))$
(%t20)

$$a_{j,i} + a_{i,j}$$

(%i21) decsym(a,2,0,[anti(all)], [ ]);
(%o21)
done
(%i22) ishow(canform(a([i,j],[ ])+a([j,i])))$
(%t22)
0

```

## 29.2 Definitions for itensor

### 29.2.1 Managing indexed objects

**entertensor** (*name*) Function  
 is a function which, by prompting, allows one to create an indexed object called *name* with any number of tensorial and derivative indices. Either a single index or a list of indices (which may be null) is acceptable input (see the example under `covdiff`).

**changename** (*old, new, expr*) Function  
 will change the name of all indexed objects called *old* to *new* in *expr*. *old* may be either a symbol or a list of the form [*name, m, n*] in which case only those indexed objects called *name* with *m* covariant and *n* contravariant indices will be renamed to *new*.

**listoftens** Function  
 Lists all tensors in a tensorial expression, complete with their indices. E.g.,

```
(%i6) ishow(a([i,j],[k])*b([u],[v])+c([x,y],[])*d([],[])*e)$
(%t6)
          d e c      + a      b
                x y      i j u,v
(%i7) ishow(listoftens(%))$
(%t7)
          k
[a      , b      , c      , d]
  i j      u,v      x y
```

**ishow** (*expr*) Function  
 displays *expr* with the indexed objects in it shown having their covariant indices as subscripts and contravariant indices as superscripts. The derivative indices are displayed as subscripts, separated from the covariant indices by a comma (see the examples throughout this document).

**indices** (*expr*) Function  
 Returns a list of two elements. The first is a list of the free indices in *expr* (those that occur only once). The second is the list of the dummy indices in *expr* (those that occur exactly twice) as the following example demonstrates.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) ishow(a([i,j],[k,l],m,n)*b([k,o],[j,m,p],q,r))$
(%t2)
          k l      j m p
a      b
  i j,m n      k o,q r
(%i3) indices(%);
(%o3)      [[l, p, i, n, o, q, r], [k, j, m]]
```

A tensor product containing the same index more than twice is syntactically illegal. `indices` attempts to deal with these expressions in a reasonable manner; however, when it is called to operate upon such an illegal expression, its behavior should be considered undefined.

**rename** (*expr*) Function  
**rename** (*expr*, *count*) Function

Returns an expression equivalent to *expr* but with the dummy indices in each term chosen from the set [%1, %2, ...], if the optional second argument is omitted. Otherwise, the dummy indices are indexed beginning at the value of *count*. Each dummy index in a product will be different. For a sum, `rename` will operate upon each term in the sum resetting the counter with each term. In this way `rename` can serve as a tensorial simplifier. In addition, the indices will be sorted alphanumerically (if `allsym` is `true`) with respect to covariant or contravariant indices depending upon the value of `flipflag`. If `flipflag` is `false` then the indices will be renamed according to the order of the contravariant indices. If `flipflag` is `true` the renaming will occur according to the to the order of the contravariant indices. It often happens that the combined effect of the two renamings will reduce an expression more than either one by itself.

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) allsym:true;
(%o2) true
(%i3) g([], [%4,%5])*g([], [%6,%7])*ichr2([%1,%4], [%3])*
ichr2([%2,%3], [u])*ichr2([%5,%6], [%1])*ichr2([%7,r], [%2])-
g([], [%4,%5])*g([], [%6,%7])*ichr2([%1,%2], [u])*
ichr2([%3,%5], [%1])*ichr2([%4,%6], [%3])*ichr2([%7,r], [%2]),noeval$
(%i4) expr:ishow(%)$

      %4 %5 %6 %7 %3 u %1 %2
(%t4) g g ichr2 ichr2 ichr2 ichr2
      %1 %4 %2 %3 %5 %6 %7 r

      %4 %5 %6 %7 u %1 %3 %2
- g g ichr2 ichr2 ichr2 ichr2
      %1 %2 %3 %5 %4 %6 %7 r■

(%i5) flipflag:true;
(%o5) true
(%i6) ishow(rename(expr))$
      %2 %5 %6 %7 %4 u %1 %3
(%t6) g g ichr2 ichr2 ichr2 ichr2
      %1 %2 %3 %4 %5 %6 %7 r

      %4 %5 %6 %7 u %1 %3 %2
- g g ichr2 ichr2 ichr2 ichr2
      %1 %2 %3 %4 %5 %6 %7 r■

(%i7) flipflag:false;
```

```

(%o7)                                false
(%i8) rename(%th(2));
(%o8)                                0
(%i9) ishow(rename(expr))$
      %1 %2 %3 %4      %5      %6      %7      u
(%t9) g      g      ichr2      ichr2      ichr2      ichr2
      %1 %6      %2 %3      %4 r      %5 %7
      %1 %2 %3 %4      %6      %5      %7      u
- g      g      ichr2      ichr2      ichr2      ichr2
      %1 %3      %2 %6      %4 r      %5 %7

```

**flipflag**

Variable

Default: `false`. If `false` then the indices will be renamed according to the order of the contravariant indices, otherwise according to the order of the covariant indices. The function influences `rename` in the following way: If `flipflag` is `false` then `rename` forms a list of the contravariant indices as they are encountered from left to right (if `true` then of the covariant indices). The first dummy index in the list is renamed to `%1`, the next to `%2`, etc. Then sorting occurs after the `rename`-ing (see the example under `rename`).

**defcon** (*tensor\_1*)

Function

**defcon** (*tensor\_1*, *tensor\_2*, *tensor\_3*)

Function

gives *tensor\_1* the property that the contraction of a product of *tensor\_1* and *tensor\_2* results in *tensor\_3* with the appropriate indices. If only one argument, *tensor\_1*, is given, then the contraction of the product of *tensor\_1* with any indexed object having the appropriate indices (say `my_tensor`) will yield an indexed object with that name, i.e. `my_tensor`, and with a new set of indices reflecting the contractions performed. For example, if `imetric:g`, then `defcon(g)` will implement the raising and lowering of indices through contraction with the metric tensor. More than one `defcon` can be given for the same indexed object; the latest one given which applies in a particular contraction will be used. `contractions` is a list of those indexed objects which have been given contraction properties with `defcon`.

**remcon** (*tensor\_1*, ..., *tensor\_n*)

Function

**remcon** (*all*)

Function

removes all the contraction properties from the *tensor\_1*, ..., *tensor\_n*). `remcon(all)` removes all contraction properties from all indexed objects.

**contract** (*expr*)

Function

Carries out the tensorial contractions in *expr* which may be any combination of sums and products. This function uses the information given to the `defcon` function. For best results, *expr* should be fully expanded. `ratexpand` is the fastest way to expand products and powers of sums if there are no variables in the denominators of the terms. The `gcd` switch should be `false` if GCD cancellations are unnecessary.

**indexed\_tensor** (*tensor*) Function

Must be executed before assigning components to a *tensor* for which a built in value already exists as with `ichr1`, `ichr2`, `icurvature`. See the example under `icurvature`.

**components** (*tensor, expr*) Function

permits one to assign an indicial value to an expression *expr* giving the values of the components of *tensor*. These are automatically substituted for the tensor whenever it occurs with all of its indices. The tensor must be of the form  $\tau([\dots], [\dots])$  where either list may be empty. *expr* can be any indexed expression involving other objects with the same free indices as *tensor*. When used to assign values to the metric tensor wherein the components contain dummy indices one must be careful to define these indices to avoid the generation of multiple dummy indices. Removal of this assignment is given to the function `remcomps`.

It is important to keep in mind that `components` cares only about the valence of a tensor, not about any particular index ordering. Thus assigning components to, say,  $x([i, -j], [])$ ,  $x([-j, i], [])$ , or  $x([i], [j])$  all produce the same result, namely components being assigned to a tensor named *x* with valence (1,1).

Components can be assigned to an indexed expression in four ways, two of which involve the use of the `components` command:

1) As an indexed expression. For instance:

```
(%i2) components(g([], [i, j]), e([], [i])*p([], [j]))$
(%i3) ishow(g([], [i, j]))$
(%t3)
          i  j
        e  p
```

2) As a matrix:

```
(%i6) components(g([i, j], []), lg);
(%o6)
done
(%i7) ishow(g([i, j], []))$
(%t7)
          g
          i  j
(%i8) g([3,3], []);
(%o8)
          1
(%i9) g([4,4], []);
(%o9)
          - 1
```

3) As a function. You can use a Maxima function to specify the components of a tensor based on its indices. For instance, the following code assigns `kdelta` to `h` if `h` has the same number of covariant and contravariant indices and no derivative indices, and `g` otherwise:

```
(%i4) h(l1,l2,[l3]):=if length(l1)=length(l2) and length(l3)=0
      then kdelta(l1,l2) else apply(g,append([l1,l2], l3))$
(%i5) ishow(h([i], [j]))$
```

```

(%t5)
          j
      kdelta
          i

(%i6) ishow(h([i,j],[k],1))$

(%t6)
          k
          g
          i j,1

```

4) Using Maxima's pattern matching capabilities, specifically the `defrule` and `applyb1` commands:

```

(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) matchdeclare(l1,listp);
(%o2) done
(%i3) defrule(r1,m(l1,[ ]),(i1:idummy(),
      g([l1[1],l1[2]],[ ])*q([i1],[ ])*e([ ],[i1])))$

(%i4) defrule(r2,m([ ],l1),(i1:idummy(),
      w([ ],[l1[1],l1[2]])*e([i1],[ ])*q([ ],[i1])))$

(%i5) ishow(m([i,n],[ ])*m([ ],[i,m]))$
          i m
(%t5)          m m
          i n

(%i6) ishow(rename(applyb1(% ,r1,r2)))$
          %1 %2 %3 m
(%t6)          e q w q e g
          %1 %2 %3 n

```

### **remcomps** (*tensor*)

Function

Unbinds all values from *tensor* which were assigned with the `components` function.

### **showcomps**

Function

Shows component assignments of a tensor, as made using the `components` command. This function can be particularly useful when a matrix is assigned to an indicial tensor using `components`, as demonstrated by the following example:

```

(%i1) load(ctensor);
(%o1) /share/tensor/ctensor.mac
(%i2) load(itensor);
(%o2) /share/tensor/itensor.lisp
(%i3) lg:matrix([sqrt(r/(r-2*m)),0,0,0],[0,r,0,0],
      [0,0,sin(theta)*r,0],[0,0,0,sqrt((r-2*m)/r)]);
          [
          r
          [ sqrt(-----) 0 0 0 ]

```



```

      [      r - 2 m      ]
      [                  ]
      [      0      r      0      0      ]
(%o3)  [                  ]
      [      0      0 r sin(theta)      0      ]
      [                  ]
      [                  r - 2 m      ]
      [      0      0      0      sqrt(-----) ]
      [                  r      ]
(%i4) components(g([i,j],[ ]),lg);
(%o4) done
(%i5) showcomps(g([i,j],[ ]));
      [      r      ]
      [ sqrt(-----) 0      0      0      ]
      [      r - 2 m      ]
      [                  ]
(%t5)  g      = [      0      r      0      0      ]
      i j      [      0      0 r sin(theta)      0      ]
      [                  ]
      [                  r - 2 m      ]
      [      0      0      0      sqrt(-----) ]
      [                  r      ]
(%o5) false

```

The `showcomps` command can also display components of a tensor of rank higher than 2.

### **idummy** ()

Function

Increments `icounter` and returns as its value an index of the form `%n` where `n` is a positive integer. This guarantees that dummy indices which are needed in forming expressions will not conflict with indices already in use (see the example under `indices`).

### **idummyx**

Variable

Is the prefix for dummy indices (see the example under `indices`).

### **icounter**

Variable

default: [1] determines the numerical suffix to be used in generating the next dummy index in the tensor package. The prefix is determined by the option `idummy` (default: `%`).

### **kdelta** (*L1*, *L2*)

Function

is the generalized Kronecker delta function defined in the `itensor` package with *L1* the list of covariant indices and *L2* the list of contravariant indices. `kdelta([i],[j])` returns the ordinary Kronecker delta. The command `ev(expr,kdelta)` causes the evaluation of an expression containing `kdelta([],[ ])` to the dimension of the manifold.

In what amounts to an abuse of this notation, `itensor` also allows `kdelta` to have 2 covariant and no contravariant, or 2 contravariant and no covariant indices, in effect providing a co(ntra)variant "unit matrix" capability. This is strictly considered a programming aid and not meant to imply that `kdelta([i,j],[,])` is a valid tensorial object.

**kdels** (*L1, L2*)

Function

Symmetricized Kronecker delta, used in some calculations. For instance:

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) kdelta([1,2],[2,1]);
(%o2)
          - 1
(%i3) kdels([1,2],[2,1]);
(%o3)
          1
(%i4) ishow(kdelta([a,b],[c,d]))$
          c      d      d      c
(%t4)      kdelta kdelta - kdelta kdelta
          a      b      a      b
(%i4) ishow(kdels([a,b],[c,d]))$
          c      d      d      c
(%t4)      kdelta kdelta + kdelta kdelta
          a      b      a      b
```

**levi\_civita** (*L*)

Function

is the permutation (or Levi-Civita) tensor which yields 1 if the list *L* consists of an even permutation of integers, -1 if it consists of an odd permutation, and 0 if some indices in *L* are repeated.

**lc2kdt** (*expr*)

Function

Simplifies expressions containing the Levi-Civita symbol, converting these to Kronecker-delta expressions when possible. The main difference between this function and simply evaluating the Levi-Civita symbol is that direct evaluation often results in Kronecker expressions containing numerical indices. This is often undesirable as it prevents further simplification. The `lc2kdt` function avoids this problem, yielding expressions that are more easily simplified with `rename` or `contract`.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) expr:isshow('levi_civita([],[i,j])*levi_civita([k,l],[,])*a([j],[k]))$
          i j k
(%t2)      levi_civita      a      levi_civita
          j      k l
(%i3) ishow(ev(expr,levi_civita))$
          i j k      1 2
```

```
(%t3)
          kdelta    a    kdelta
          1 2 j      k l
(%i4) ishow(ev(%,kdelta))$
          i      j      j      i    k
(%t4) (kdelta  kdelta - kdelta  kdelta ) a
          1      2      1      2 j
          1      2      2      1
          (kdelta  kdelta - kdelta  kdelta )
          k      l      k      l
(%i5) ishow(lc2kdt(expr))$
          k      i      j      k      j      i
(%t5)      a  kdelta  kdelta - a  kdelta  kdelta
          j      k      l      j      k      l
(%i6) ishow(contract(expand(%)))$
          i      i
(%t6)      a - a kdelta
          1      1
```

The `lc2kdt` function sometimes makes use of the metric tensor. If the metric tensor was not defined previously with `imetric`, this results in an error.

```
(%i7) expr:ishow('levi_civita([], [i,j])*levi_civita([], [k,l])*a([j,k], []))$
          i j      k l
(%t7)      levi_civita      levi_civita      a
          j k

(%i8) ishow(lc2kdt(expr))$
Maxima encountered a Lisp error:

Error in $IMETRIC [or a callee]: $IMETRIC [or a callee] requires less than two
arguments.

Automatically continuing.
To reenale the Lisp debugger set *debugger-hook* to nil.
(%i9) imetric(g);
(%o9)      done
(%i10) ishow(lc2kdt(expr))$
          %3 i      k      %4 j      l      %3 i      l      %4 j      k
(%t10) (g      kdelta  g      kdelta - g      kdelta  g      kdelta ) a
          %3      %4      %3      %4      %3      %4      %4 j k
(%i11) ishow(contract(expand(%)))$
          l i      l i
(%t11)      a - a g
```

## lc\_l

## Function

Simplification rule used for expressions containing the unevaluated Levi-Civita symbol (`levi_civita`). Along with `lc_u`, it can be used to simplify many expressions more efficiently than the evaluation of `levi_civita`. For example:

```

(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2)  e11:ishow('levi_civita([i,j,k],[i,j,k]))$
          i j
          a a levi_civita
(%t2)
(%i3)  e12:ishow('levi_civita([i,j,k],[i,j,k]))$
          i j k
          levi_civita a a
(%t3)
(%i4)  ishow(canform(contract(expand(applyb1(e11,lc_1,lc_u))))))$
(%t4)  0
(%i5)  ishow(canform(contract(expand(applyb1(e12,lc_1,lc_u))))))$
(%t5)  0

```

**lc\_u**

Function

Simplification rule used for expressions containing the unevaluated Levi-Civita symbol (`levi_civita`). Along with `lc_u`, it can be used to simplify many expressions more efficiently than the evaluation of `levi_civita`. For details, see `lc_1`.

**canten** (*expr*)

Function

Simplifies *expr* by renaming (see `rename`) and permuting dummy indices. `rename` is restricted to sums of tensor products in which no derivatives are present. As such it is limited and should only be used if `canform` is not capable of carrying out the required simplification.

**29.2.2 Tensor symmetries****allsym**

Variable

Default: `false`. if `true` then all indexed objects are assumed symmetric in all of their covariant and contravariant indices. If `false` then no symmetries of any kind are assumed in these indices. Derivative indices are always taken to be symmetric unless `iframe_flag` is set to `true`.

**decsym** (*tensor*, *m*, *n*, [*cov\_1*, *cov\_2*, ...], [*contr\_1*, *contr\_2*, ...])

Function

Declares symmetry properties for *tensor* of *m* covariant and *n* contravariant indices. The *cov<sub>i</sub>* and *contr<sub>i</sub>* are pseudofunctions expressing symmetry relations among the covariant and contravariant indices respectively. These are of the form `symoper(index_1, index_2, ...)` where `symoper` is one of `sym`, `anti` or `cyc` and the *index<sub>i</sub>* are integers indicating the position of the index in the *tensor*. This will declare *tensor* to be symmetric, antisymmetric or cyclic respectively in the *index<sub>i</sub>*. `symoper(all)` is also an allowable form which indicates all indices obey the symmetry condition. For example, given an object `b` with 5 covariant indices, `decsym(b,5,3,[sym(1,2),anti(3,4)],[cyc(all)])` declares `b` symmetric in its first and second and antisymmetric in its third and fourth covariant indices, and

cyclic in all of its contravariant indices. Either list of symmetry declarations may be null. The function which performs the simplifications is `canform` as the example below illustrates.

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) expr:contract(expand(a([i1,j1,k1],[[]])*kdels([i,j,k],[i1,j1,k1])))$
(%i3) ishow(expr)$
(%t3)      a      + a      + a      + a      + a      + a
           k j i    k i j    j k i    j i k    i k j    i j k
(%i4) decsym(a,3,0,[sym(all)],[]);
(%o4) done
(%i5) ishow(canform(expr))$
(%t5)      6 a
           i j k

(%i6) remsym(a,3,0);
(%o6) done
(%i7) decsym(a,3,0,[anti(all)],[]);
(%o7) done
(%i8) ishow(canform(expr))$
(%t8)      0

(%i9) remsym(a,3,0);
(%o9) done
(%i10) decsym(a,3,0,[cyc(all)],[]);
(%o10) done
(%i11) ishow(canform(expr))$
(%t11)      3 a      + 3 a
           i k j      i j k

(%i12) dispsym(a,3,0);
(%o12)      [[cyc, [[1, 2, 3]], []]]
```

**remsym** (*tensor*, *m*, *n*) Function  
Removes all symmetry properties from *tensor* which has *m* covariant indices and *n* contravariant indices.

**canform** (*expr*) Function  
Simplifies *expr* by renaming dummy indices and reordering all indices as dictated by symmetry conditions imposed on them. If `allsym` is `true` then all indices are assumed symmetric, otherwise symmetry information provided by `decsym` declarations will be used. The dummy indices are renamed in the same manner as in the `rename` function. When `canform` is applied to a large expression the calculation may take a considerable amount of time. This time can be shortened by calling `rename` on the expression first. Also see the example under `decsym`. Note: `canform` may not be able to reduce an expression completely to its simplest form although it will always return a mathematically correct result.

### 29.2.3 Indicial tensor calculus

**diff** (*expr*, *v\_1*, [*n\_1*, [*v\_2*, *n\_2*] ...]) Function  
 is the usual Maxima differentiation function which has been expanded in its abilities for *itensor*. It takes the derivative of *expr* with respect to *v\_1* *n\_1* times, with respect to *v\_2* *n\_2* times, etc. For the tensor package, the function has been modified so that the *v\_i* may be integers from 1 up to the value of the variable *dim*. This will cause the differentiation to be carried out with respect to the *v\_i*th member of the list *vect\_coords*. If *vect\_coords* is bound to an atomic variable, then that variable subscripted by *v\_i* will be used for the variable of differentiation. This permits an array of coordinate names or subscripted names like *x*[1], *x*[2], ... to be used.

**idiff** (*expr*, *v\_1*, [*n\_1*, [*v\_2*, *n\_2*] ...]) Function  
 Indicial differentiation. Unlike *diff*, which differentiates with respect to an independent variable, *idiff* can be used to differentiate with respect to a coordinate. For an indexed object, this amounts to appending the *v\_i* as derivative indices. Subsequently, derivative indices will be sorted, unless *iframe\_flag* is set to *true*.  
*idiff* can also differentiate the determinant of the metric tensor. Thus, if *imetric* has been bound to *G* then *idiff*(*determinant*(*g*),*k*) will return *2\*determinant*(*g*)\**ichr2*([%i,*k*],[%i]) where the dummy index %i is chosen appropriately.

**liediff** (*v*, *ten*) Function  
 Computes the Lie-derivative of the tensorial expression *ten* with respect to the vector field *v*. *ten* should be any indexed tensor expression; *v* should be the name (without indices) of a vector field. For example:

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) ishow(liediff(v,a([i,j],[k],1)))$
      k %2 %2 %2
(%t2) b (v a + v a + v a )
      ,1 i j,%2 ,j i %2 ,i %2 j
      %1 k %1 k %1 k
      + (v b - b v + v b ) a
      ,%1 1 ,1 ,%1 ,1 ,%1 i j
```

**rediff** (*ten*) Function  
 Evaluates all occurrences of the *idiff* command in the tensorial expression *ten*.

**undiff** (*expr*) Function  
 Returns an expression equivalent to *expr* but with all derivatives of indexed objects replaced by the noun form of the *idiff* function. Its arguments would yield that indexed object if the differentiation were carried out. This is useful when it is desired to replace a differentiated indexed object with some function definition resulting in *expr* and then carry out the differentiation by saying *ev*(*expr*, *idiff*).

**evundiff**

Function

Equivalent to the execution of `undiff`, followed by `ev` and `rediff`.

The point of this operation is to easily evaluate expressions that cannot be directly evaluated in derivative form. For instance, the following causes an error:

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) icurvature([i,j,k],[l],m);
Maxima encountered a Lisp error:
```

```
Error in $ICURVATURE [or a callee]: $ICURVATURE [or a callee] requires less t
```

Automatically continuing.

To reenable the Lisp debugger set `*debugger-hook*` to `nil`.

However, if `icurvature` is entered in noun form, it can be evaluated using `evundiff`:

```
(%i3) ishow('icurvature([i,j,k],[l],m))$
(%t3)          icurvature
          i j k,m
(%i4) ishow(evundiff(%))$
(%t4) - 1      1      %1      1      %1
        i k,j m  %1 j  i k,m  %1 j,m  i k
        + 1      1      %1      1      %1
          i j,k m  %1 k  i j,m  %1 k,m  i j
```

Note: In earlier versions of Maxima, derivative forms of the Christoffel-symbols also could not be evaluated. This has been fixed now, so `evundiff` is no longer necessary for expressions like this:

```
(%i5) imetric(g);
(%o5) done
(%i6) ishow(ichr2([i,j],[k],l))$
k %3
g      (g      - g      + g      )
      j %3,i l  i j,%3 l  i %3,j l
(%t6) -----
              2
              k %3
              g      (g      - g      + g      )
              ,l      j %3,i  i j,%3  i %3,j
+ -----
              2
```

**flush** (*expr*, *tensor-1*, *tensor-2*, ...)

Function

Set to zero, in *expr*, all occurrences of the *tensor-*i** that have no derivative indices.

**flushd** (*expr*, *tensor\_1*, *tensor\_2*, ...) Function  
 Set to zero, in *expr*, all occurrences of the *tensor\_i* that have derivative indices.

**flushnd** (*expr*, *tensor*, *n*) Function  
 Set to zero, in *expr*, all occurrences of the differentiated object *tensor* that have *n* or more derivative indices as the following example demonstrates.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) ishow(a([i],[J,r],k,r)+a([i],[j,r,s],k,r,s))$
(%t2)
          J r      j r s
a      + a
i,k r   i,k r s

(%i3) ishow(flushnd(%,a,3))$
(%t3)
          J r
a
          i,k r
```

**coord** (*tensor\_1*, *tensor\_2*, ...) Function  
 Gives *tensor\_i* the coordinate differentiation property that the derivative of contravariant vector whose name is one of the *tensor\_i* yields a Kronecker delta. For example, if **coord**(*x*) has been done then **idiff**(*x*([ ], [i]), *j*) gives **kdelta**([i], [j]). **coord** is a list of all indexed objects having this property.

**remcoord** (*tensor\_1*, *tensor\_2*, ...) Function  
**remcoord** (*all*) Function  
 Removes the coordinate differentiation property from the *tensor\_i* that was established by the function **coord**. **remcoord**(*all*) removes this property from all indexed objects.

**makebox** (*expr*) Function  
 Display *expr* in the same manner as **show**; however, any tensor d'Alembertian occurring in *expr* will be indicated using the symbol [ ]. For example, [ ]p([m], [n]) represents  $g([ ], [i, j]) * p([m], [n], i, j)$ .

**conmetderiv** (*expr*, *tensor*) Function  
 Simplifies expressions containing ordinary derivatives of both covariant and contravariant forms of the metric tensor (the current restriction). For example, **conmetderiv** can relate the derivative of the contravariant metric tensor with the Christoffel symbols as seen from the following:

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) ishow(g([ ], [a,b], c))$
(%t2)
          a b
g
          ,c
```



```
(%i3) ishow(commetderiv(%,g))$
(%t3)          %1 b      a      %1 a      b
          - g      ichr2      - g      ichr2
                    %1 c      %1 c
```

**flush1deriv** (*expr*, *tensor*) Function  
Set to zero, in *expr*, all occurrences of *tensor* that have exactly one derivative index.

## 29.2.4 Tensors in curved spaces

**imetric** (*g*) Function  
Specifies the metric by assigning the variable `imetric:g` in addition, the contraction properties of the metric *g* are set up by executing the commands `defcon(g),defcon(g,g,kdelta)`. The variable `imetric`, default: `[]`, is bound to the metric, assigned by the `imetric(g)` command.

**ichr1** (*[i, j, k]*) Function  
Yields the Christoffel symbol of the first kind via the definition

$$(g_{ik,j} + g_{jk,i} - g_{ij,k})/2 .$$

To evaluate the Christoffel symbols for a particular metric, the variable `imetric` must be assigned a name as in the example under `chr2`.

**ichr2** (*[i, j], [k]*) Function  
Yields the Christoffel symbol of the second kind defined by the relation

$$\text{ichr2}([i, j], [k]) = g^{ks} (g_{is,j} + g_{js,i} - g_{ij,s})/2$$

**icurvature** (*[i, j, k], [h]*) Function  
Yields the Riemann curvature tensor in terms of the Christoffel symbols of the second kind (`ichr2`). The following notation is used:

$$\text{icurvature}_{i j k}^h = - \text{ichr2}_{i k, j}^h - \text{ichr2}_{i j, k}^h + \text{ichr2}_{i j, k}^h + \text{ichr2}_{i k, j}^h$$

**covdiff** (*expr*, *v\_1*, *v\_2*, ...) Function  
Yields the covariant derivative of *expr* with respect to the variables *v\_i* in terms of the Christoffel symbols of the second kind (`ichr2`). In order to evaluate these, one should use `ev(expr, ichr2)`.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) entertensor()$
Enter tensor name: a;
Enter a list of the covariant indices: [i,j];
Enter a list of the contravariant indices: [k];
Enter a list of the derivative indices: [];

(%t2)
          k
          a
          i j

(%i3) ishow(covdiff(%s))$
          k      %1      k      %1      k      %1
(%t3)  - a      ichr2  - a      ichr2  + a      + ichr2  a
          i %1      j s   %1 j      i s   i j,s      %1 s i j
```

**lorentz\_gauge** (*expr*)

Function

Imposes the Lorentz condition by substituting 0 for all indexed objects in *expr* that have a derivative index identical to a contravariant index.

**igeodesic\_coords** (*expr, name*)

Function

Causes undifferentiated Christoffel symbols and first derivatives of the metric tensor vanish in *expr*. The *name* in the **igeodesic\_coords** function refers to the metric *name* (if it appears in *expr*) while the connection coefficients must be called with the names *ichr1* and/or *ichr2*. The following example demonstrates the verification of the cyclic identity satisfied by the Riemann curvature tensor using the **igeodesic\_coords** function.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) ishow(icurvature([r,s,t],[u]))$
          u      u      %1      u      u      %1
(%t2)  - ichr2      - ichr2      ichr2      + ichr2      + ichr2      ichr2
          r t,s      %1 s      r t      r s,t      %1 t      r s
(%i3) ishow(igeodesic_coords(%s,ichr2))$
          u      u
(%t3)      ichr2      - ichr2
          r s,t      r t,s
(%i4) ishow(igeodesic_coords(icurvature([r,s,t],[u]),ichr2)+
          igeodesic_coords(icurvature([s,t,r],[u]),ichr2)+
          igeodesic_coords(icurvature([t,r,s],[u]),ichr2))$
          u      u      u      u      u
(%t4)  - ichr2      + ichr2      + ichr2      - ichr2      - ichr2
          t s,r      t r,s      s t,r      s r,t      r t,s
          u
          + ichr2
```

r s,t

```
(%i5) canform(%);
(%o5) 0
```

### 29.2.5 Moving frames

Maxima now has the ability to perform calculations using moving frames. These can be orthonormal frames (tetrads, vielbeins) or an arbitrary frame.

To use frames, you must first set `iframe_flag` to `true`. This causes the Christoffel-symbols, `ichr1` and `ichr2`, to be replaced by the more general frame connection coefficients `icc1` and `icc2` in calculations. Specially, the behavior of `covdiff` and `icurvature` is changed.

The frame is defined by two tensors: the inverse frame field (`ifri`), and the frame metric `ifg`. The frame metric is the identity matrix for orthonormal frames, or the Lorentz metric for orthonormal frames in Minkowski spacetime. The inverse frame field defines the frame base (unit vectors). Contraction properties are defined for the frame field and the frame metric.

When `iframe_flag` is true, many `itensor` expressions use the frame metric `ifg` instead of the metric defined by `imetric` for raising and lowering indices.

**IMPORTANT:** Setting the variable `iframe_flag` to `true` does NOT undefine the contraction properties of a metric defined by a call to `defcon` or `imetric`. If a frame field is used, it is best to define the metric by assigning its name to the variable `imetric` and NOT invoke the `imetric` function.

Maxima uses these two tensors to define the frame coefficients (`ifc1` and `ifc2`) which form part of the connection coefficients (`icc1` and `icc2`), as the following example demonstrates:

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) iframe_flag:true;
(%o2) true
(%i3) ishow(covdiff(v([],[i]),j))$
(%t3)      i      i      %1
      v  + icc2  v
      ,j      %1 j
(%i4) ishow(ev(%,icc2))$
(%t4)      %1      i      i      i
      v  (ifc2  + ichr2  ) + v
      %1 j      %1 j      ,j
(%i5) ishow(ev(%,ifc2))$
      %1      i %2
      v  ifg  (ifb  - ifb  + ifb  )
      j %2 %1      %2 %1 j      %1 j %2      i
(%t5)  ----- + v
      2
(%i6) ishow(ifb([a,b,c]))$
```

$$(\%t6) \quad \begin{array}{cccc} & \%5 & \%4 & \\ \text{ifri} & \text{ifri} & (\text{ifri} & - \text{ifri} \\ a & b & c \%4,\%5 & c \%5,\%4 \end{array} )$$

An alternate method is used to compute the frame bracket (`ifb`) if the `iframe_bracket_form` flag is set to `false`:

$$(\%i8) \text{ block}([\text{iframe\_bracket\_form}:\text{false}], \text{ishow}(\text{ifb}([a,b,c])))\$$$

$$(\%t8) \quad \begin{array}{cccc} \%7 & \%6 & \%6 & \%7 \\ (\text{ifri} & \text{ifri} & - \text{ifri} & \text{ifri} ) \text{ifri} \\ a & b,\%7 & a,\%7 & b & c \%6 \end{array}$$

**iframes** () Function

Since in this version of Maxima, contraction identities for `ifr` and `ifri` are always defined, as is the frame bracket (`ifb`), this function does nothing.

**ifb** Variable

The frame bracket. The contribution of the frame metric to the connection coefficients is expressed using the frame bracket:

$$\text{ifc1} = \frac{- \text{ifb} \quad + \text{ifb} \quad + \text{ifb}}{\text{abc} \quad \quad \quad 2}$$

The frame bracket itself is defined in terms of the frame field and frame metric. Two alternate methods of computation are used depending on the value of `frame_bracket_form`. If true (the default) or if the `itorsion_flag` is true:

$$\text{ifb} = \begin{array}{cccc} d & e & & f \\ \text{ifri} & \text{ifri} & (\text{ifri} & - \text{ifri} & - \text{ifri} & \text{itr} \\ abc & b & c & a d,e & a e,d & a f & d e \end{array} )$$

Otherwise:

$$\text{ifb} = (\text{ifri} \quad \text{ifri} \quad - \text{ifri} \quad \text{ifri} ) \text{ifri}$$

**icc1** Variable

Connection coefficients of the first kind. In `itensor`, defined as

$$\text{icc1} = \text{ichr1} - \text{ikt1} - \text{inmc1}$$

$$\text{abc} \quad \text{abc} \quad \text{abc} \quad \text{abc}$$

In this expression, if `iframe_flag` is true, the Christoffel-symbol `ichr1` is replaced with the frame connection coefficient `ifc1`. If `itorsion_flag` is false, `ikt1` will be omitted. It is also omitted if a frame base is used, as the torsion is already calculated as part of the frame bracket. Lastly, if `inonmet_flag` is false, `inmc1` will not be present.

**icc2** Variable  
 Connection coefficients of the second kind. In `itensor`, defined as

$$\text{icc2} \quad \text{c} \quad \text{c} \quad \text{c} \quad \text{c}$$

$$= \frac{\text{ichr2}}{\text{ab}} - \frac{\text{ikt2}}{\text{ab}} - \frac{\text{inmc2}}{\text{ab}}$$

In this expression, if `iframe_flag` is true, the Christoffel-symbol `ichr2` is replaced with the frame connection coefficient `ifc2`. If `itorsion_flag` is false, `ikt2` will be omitted. It is also omitted if a frame base is used, as the torsion is already calculated as part of the frame bracket. Lastly, if `inonmet_flag` is false, `inmc2` will not be present.

**ifc1** Variable  
 Frame coefficient of the first kind (also known as Ricci-rotation coefficients.) This tensor represents the contribution of the frame metric to the connection coefficient of the first kind. Defined as:

$$\text{ifc1} \quad \text{c} \quad \text{c} \quad \text{c}$$

$$= \frac{-\text{ifb}_{c a b} + \text{ifb}_{b c a} + \text{ifb}_{a b c}}{2}$$

**ifc2** Variable  
 Frame coefficient of the first kind. This tensor represents the contribution of the frame metric to the connection coefficient of the first kind. Defined as a permutation of the frame bracket (`ifb`) with the appropriate indices raised and lowered as necessary:

$$\text{ifc2} \quad \text{c} \quad \text{cd}$$

$$= \frac{\text{ifg}_{ab}}{\text{ab}} \quad \text{ifc1}_{abd}$$

**ifr** Variable  
 The frame field. Contracts with the inverse frame field (`ifri`) to form the frame metric (`ifg`).

- ifri** Variable
- The inverse frame field. Specifies the frame base (basis vectors). Along with the frame metric, it forms the basis of all calculations based on frames.
- ifg** Variable
- The frame metric. Defaults to `kdelta`, but can be changed using `components`.
- ifgi** Variable
- The inverse frame metric. Contracts with the frame metric (`ifg`) to `kdelta`.
- iframe\_bracket\_form** Variable
- Specifies how the frame bracket (`ifb`) is computed. Default is `true`.

### 29.2.6 Torsion and nonmetricity

Maxima can now take into account torsion and nonmetricity. When the flag `itorsion_flag` is set to `true`, the contribution of torsion is added to the connection coefficients. Similarly, when the flag `inonmet_flag` is true, nonmetricity components are included.

- inm** Variable
- The nonmetricity vector. Conformal nonmetricity is defined through the covariant derivative of the metric tensor. Normally zero, the metric tensor's covariant derivative will evaluate to the following when `inonmet_flag` is set to `true`:

$$g_{ij;k} = -g_{ij} \text{ inm}_k$$

- inmc1** Variable
- Covariant permutation of the nonmetricity vector components. Defined as

$$\text{inmc1}_{abc} = \frac{g_{ab} \text{ inm}_c - \text{inm}_a g_{bc} - g_{ac} \text{ inm}_b}{2}$$

(Substitute `ifg` in place of `g` if a frame metric is used.)

- inmc2** Variable
- Contravariant permutation of the nonmetricity vector components. Used in the connection coefficients if `inonmet_flag` is true. Defined as:

$$\text{inmc2}_{ab} = \frac{-\text{inm}_a \text{ kdelta}_{cb} - \text{ kdelta}_{ca} \text{ inm}_b + g_{cd} \text{ inm}_d g_{ab}}{2}$$

(Substitute `ifg` in place of `g` if a frame metric is used.)

**ikt1**

Variable

Covariant permutation of the torsion tensor (also known as contorsion). Defined as:

$$\text{ikt1} = \frac{-g \begin{matrix} d \\ ad \end{matrix} \text{itr} \begin{matrix} d \\ cb \end{matrix} - g \begin{matrix} d \\ bd \end{matrix} \text{itr} \begin{matrix} d \\ ca \end{matrix} - \text{itr} \begin{matrix} d \\ ab \end{matrix} g \begin{matrix} d \\ cd \end{matrix}}{2}$$

(Substitute ifg in place of g if a frame metric is used.)

**ikt2**

Variable

Contravariant permutation of the torsion tensor (also known as contorsion). Defined as:

$$\text{ikt2} = g \begin{matrix} c \\ ab \end{matrix} \text{ikt1} \begin{matrix} cd \\ abd \end{matrix}$$

(Substitute ifg in place of g if a frame metric is used.)

**itr**

Variable

The torsion tensor. For a metric with torsion, repeated covariant differentiation on a scalar function will not commute, as demonstrated by the following example:

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) imetric:g;
(%o2) g
(%i3) covdiff(covdiff(f([],[]),i),j)-covdiff(covdiff(f([],[]),j),i)$
(%i4) ishow(%)$
(%t4) f ichr2 - f ichr2
      ,%4 j i ,%2 i j
(%i5) canform(%)$
(%o5) 0
(%i6) itorsion_flag:true;
(%o6) true
(%i7) covdiff(covdiff(f([],[]),i),j)-covdiff(covdiff(f([],[]),j),i)$
(%i8) ishow(%)$
(%t8) f icc2 - f icc2 - f + f
      ,%8 j i ,%6 i j ,j i ,i j
(%i9) ishow(canform(%)$
(%t9) f icc2 - f icc2
      ,%1 j i ,%1 i j
```

```
(%i10) ishow(canform(ev(% ,icc2)))$
(%t10)
          %1          %1
          f    ikt2  - f    ikt2
          ,%1    i j  ,%1    j i
(%i11) ishow(canform(ev(% ,ikt2)))$
          %2 %1          %2 %1
          f    g    ikt1  - f    g    ikt1
          ,%2          i j %1  ,%2          j i %1
(%i12) ishow(factor(canform(rename(expand(ev(% ,ikt1))))))$
          %3 %2          %1          %1
          f    g    g    (itr  - itr  )
          ,%3          %2 %1    j i    i j
(%t12)
          -----
                          2
(%i13) decsym(itr,2,1,[anti(all)],[]);
(%o13)
done
(%i14) defcon(g,g,kdelta);
(%o14)
done
(%i15) subst(g,nounify(g),%th(3))$
(%i16) ishow(canform(contract(%)))$
(%t16)
          %1
          - f    itr
          ,%1    i j
```

### 29.2.7 Exterior algebra

The `itensor` package can perform operations on totally antisymmetric covariant tensor fields. A totally antisymmetric tensor field of rank (0,L) corresponds with a differential L-form. On these objects, a multiplication operation known as the exterior product, or wedge product, is defined.

Unfortunately, not all authors agree on the definition of the wedge product. Some authors prefer a definition that corresponds with the notion of antisymmetrization: in these works, the wedge product of two vector fields, for instance, would be defined as

$$a_i \wedge a_j = \frac{a_i a_j - a_j a_i}{2}$$

More generally, the product of a p-form and a q-form would be defined as

$$A_{i_1..i_p} \wedge B_{j_1..j_q} = \frac{1}{(p+q)!} D_{i_1..i_p j_1..j_q}^{k_1..k_p l_1..l_q} A_{k_1..k_p} B_{l_1..l_q}$$

where D stands for the Kronecker-delta.

Other authors, however, prefer a “geometric” definition that corresponds with the notion of the volume element:

$$a_i \wedge a_j = a_i a_j - a_j a_i$$



and, in the general case

$$A \wedge B = \frac{1}{p! q!} \sum_{i_1..i_p, j_1..j_q} A_{i_1..i_p} B_{j_1..j_q} = \frac{1}{p! q!} \sum_{i_1..i_p, j_1..j_q, k_1..k_p, l_1..l_q} A_{i_1..i_p} B_{j_1..j_q} \epsilon^{k_1..k_p l_1..l_q}$$

Since `itensor` is a tensor algebra package, the first of these two definitions appears to be the more natural one. Many applications, however, utilize the second definition. To resolve this dilemma, a flag has been implemented that controls the behavior of the wedge product: if `igeowedge_flag` is `false` (the default), the first, "tensorial" definition is used, otherwise the second, "geometric" definition will be applied.

"~"

special symbol

The wedge product operator is denoted by the tilde `~`. This is a binary operator. Its arguments should be expressions involving scalars, covariant tensors of rank one, or covariant tensors of rank 1 that have been declared antisymmetric in all covariant indices.

The behavior of the wedge product operator is controlled by the `igeowedge_flag` flag, as in the following example:

```
(%i1) load(itensor);
(%o1) /share/tensor/itensor.lisp
(%i2) ishow(a([i])~b([j]))$
(%t2)
      a  b  - b  a
      i  j   i  j
      -----
              2
(%i3) decsym(a,2,0,[anti(all)],[]);
(%o3) done
(%i4) ishow(a([i,j])~b([k]))$
(%t4)
      a  b  + b  a  - a  b
      i  j  k   i  j  k   i  k  j
      -----
              3
(%i5) igeowedge_flag:true;
(%o5) true
(%i6) ishow(a([i])~b([j]))$
(%t6)
      a  b  - b  a
      i  j   i  j
(%i7) ishow(a([i,j])~b([k]))$
(%t7)
      a  b  + b  a  - a  b
      i  j  k   i  j  k   i  k  j
```

"|"

special symbol

The vertical bar `|` denotes the "contraction with a vector" binary operation. When a totally antisymmetric covariant tensor is contracted with a contravariant vector, the result is the same regardless which index was used for the contraction. Thus, it is possible to define the contraction operation in an index-free manner.

In the `itensor` package, contraction with a vector is always carried out with respect to the first index in the literal sorting order. This ensures better simplification of expressions involving the `|` operator. For instance:

```

(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) decsym(a,2,0,[anti(all)],[]);
(%o2)      done
(%i3) ishow(a([i,j],[])|v)$
(%t3)      %1
            v   a
            %1 j
(%i4) ishow(a([j,i],[])|v)$
(%t4)      %1
            - v   a
            %1 j

```

Note that it is essential that the tensors used with the `|` operator be declared totally antisymmetric in their covariant indices. Otherwise, the results will be incorrect.

### **extdiff** (*expr*, *i*)

Function

Computes the exterior derivative of *expr* with respect to the index *i*. The exterior derivative is formally defined as the wedge product of the partial derivative operator and a differential form. As such, this operation is also controlled by the setting of `igeowedge_flag`. For instance:

```

(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) ishow(extdiff(v([i]),j))$
(%t2)      v   - v
            j,i   i,j
            -----
            2
(%i3) decsym(a,2,0,[anti(all)],[]);
(%o3)      done
(%i4) ishow(extdiff(a([i,j]),k))$
(%t4)      a   - a   + a
            j k,i   i k,j   i j,k
            -----
            3
(%i5) igeowedge_flag:true;
(%o5)      true
(%i6) ishow(extdiff(v([i]),j))$
(%t6)      v   - v
            j,i   i,j
(%i7) ishow(extdiff(a([i,j]),k))$
(%t7)      a   - a   + a
            j k,i   i k,j   i j,k

```

### **igeowedge\_flag**

Variable

Controls the behavior of the wedge product and exterior derivative. When set to `false` (the default), the notion of differential forms will correspond with that of a totally antisymmetric covariant tensor field. When set to `true`, differential forms will agree with the notion of the volume element.

### 29.2.8 Exporting TeX expressions

The `itensor` package provides limited support for exporting tensor expressions to TeX. Since `itensor` expressions appear as function calls, the regular Maxima `tex` command will not produce the expected output. You can try instead the `tentex` command, which attempts to translate tensor expressions into appropriately indexed TeX objects.

**tentex** (*expr*) Function

To use the `tentex` function, you must first load `tentex`, as in the following example:

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
(%i2) load(tentex);
(%o2)      /share/tensor/tentex.lisp
(%i3) idummyx:m;
(%o3)
(%i4) ishow(icurvature([j,k,l],[i]))$
(%t4)      m1      i      m1      i      i      i
ichr2      ichr2      - ichr2      ichr2      - ichr2      + ichr2
      j k      m1 l      j l      m1 k      j l,k      j k,l
(%i5) tentex(%)$
$$\Gamma_{j,k}^{m_1}\backslash,\Gamma_{l,m_1}^{i}-\Gamma_{j,l}^{m_1}\backslash,
\Gamma_{k,m_1}^{i}-\Gamma_{j,l,k}^{i}+\Gamma_{j,k,l}^{i}$$
```

Note the use of the `idummyx` assignment, to avoid the appearance of the percent sign in the TeX expression, which may lead to compile errors.

NB: This version of the `tentex` function is somewhat experimental.

### 29.2.9 Interfacing with ctensor

The `itensor` package has the ability to generate Maxima code that can then be executed in the context of the `ctensor` package. The function that performs this task is `ic_convert`.

**ic\_convert** (*eqn*) Function

Converts the `itensor` equation *eqn* to a `ctensor` assignment statement. Implied sums over dummy indices are made explicit while indexed objects are transformed into arrays (the array subscripts are in the order of covariant followed by contravariant indices of the indexed objects). The derivative of an indexed object will be replaced by the noun form of `diff` taken with respect to `ct_coords` subscripted by the derivative index. The Christoffel symbols `ichr1` and `ichr2` will be translated to `lcs` and `mcs`, respectively and if `metricconvert` is `true` then all occurrences of the metric with two covariant (contravariant) indices will be renamed to `lg` (`ug`). In addition, `do` loops will be introduced summing over all free indices so that the transformed assignment statement can be evaluated by just doing `ev`. The following examples demonstrate the features of this function.

```
(%i1) load(itensor);
(%o1)      /share/tensor/itensor.lisp
```

```

(%i2) eqn:ishow(t([i,j],[k])=f([],[])*g([1,m],[])*a([],[m],j)*b([i],[1,k]))$
          k      m  l k
(%t2)      t      = f a  b  g
          i j      ,j i  l m

(%i3) ic_convert(eqn);
(%o3) for i thru dim do (for j thru dim

do (for k thru dim do t      : f sum(sum(diff(a , ct_coords ) b
          i, j, k          m          j i, l, k

g      , l, 1, dim), m, 1, dim))
  l, m
(%i4) imetric(g);
(%o4)                                     done
(%i5) metricconvert:true;
(%o5)                                     true
(%i6) ic_convert(eqn);
(%o6) for i thru dim do (for j thru dim

do (for k thru dim do t      : f sum(sum(diff(a , ct_coords ) b
          i, j, k          m          j i, l, k

lg      , l, 1, dim), m, 1, dim))
  l, m

```

### 29.2.10 Reserved words

The following Maxima words are used by the `itensor` package internally and should not be redefined:

Keyword	Comments
indices2()	Internal version of indices()
conti	Lists contravariant indices
covi	Lists covariant indices of a indexed object
deri	Lists derivative indices of an indexed object
name	Returns the name of an indexed object
concan	
irpmon	
lc0	
_lc2kdt0	
_lcprod	
_extlc	



## 30 ctensor

### 30.1 Introduction to ctensor

`ctensor` is a component tensor manipulation package. To use the `ctensor` package, type `load(ctensor)`. To begin an interactive session with `ctensor`, type `csetup()`. You are first asked to specify the dimension of the manifold. If the dimension is 2, 3 or 4 then the list of coordinates defaults to `[x,y]`, `[x,y,z]` or `[x,y,z,t]` respectively. These names may be changed by assigning a new list of coordinates to the variable `ct_coords` (described below) and the user is queried about this. **\*\* Care must be taken to avoid the coordinate names conflicting with other object definitions \*\***.

Next, the user enters the metric either directly or from a file by specifying its ordinal position. As an example of a file of common metrics, see `share/tensor/metrics.mac`. The metric is stored in the matrix `LG`. Finally, the metric inverse is computed and stored in the matrix `UG`. One has the option of carrying out all calculations in a power series.

A sample protocol is begun below for the static, spherically symmetric metric (standard coordinates) which will be applied to the problem of deriving Einstein's vacuum equations (which lead to the Schwarzschild solution) as an example. Many of the functions in `ctensor` will be displayed for the standard metric as examples.

```
(%i1) load(ctensor);
(%o1)      /usr/local/lib/maxima/share/tensor/ctensor.mac
(%i2) csetup();
Enter the dimension of the coordinate system:
4;
Do you wish to change the coordinate names?
n;
Do you want to
1. Enter a new metric?

2. Enter a metric from a file?

3. Approximate a metric with a Taylor series?
1;

Is the matrix  1. Diagonal  2. Symmetric  3. Antisymmetric  4. General
Answer 1, 2, 3 or 4
1;
Row 1 Column 1:
a;
Row 2 Column 2:
x^2;
Row 3 Column 3:
x^2*sin(y)^2;
Row 4 Column 4:
-d;

Matrix entered.
```

Enter functional dependencies with the DEPENDS function or 'N' if none depends([a,d],x);

Do you wish to see the metric?

y;

```
[ a 0      0      0 ]
[      ]
[      2      ]
[ 0 x      0      0 ]
[      ]
[      2      2      ]
[ 0 0 x sin (y) 0 ]
[      ]
[ 0 0      0      - d ]
```

(%o2)

done

(%i3) christof(mcs);

(%t3) 
$$\text{mcs}_{1, 1, 1} = \frac{a}{2x}$$

(%t4) 
$$\text{mcs}_{1, 2, 2} = -\frac{1}{x}$$

(%t5) 
$$\text{mcs}_{1, 3, 3} = -\frac{1}{x}$$

(%t6) 
$$\text{mcs}_{1, 4, 4} = \frac{d}{2x}$$

(%t7) 
$$\text{mcs}_{2, 2, 1} = -\frac{x}{a}$$

(%t8) 
$$\text{mcs}_{2, 3, 3} = \frac{\cos(y)}{\sin(y)}$$

(%t9) 
$$\text{mcs}_{3, 3, 1} = -\frac{x \sin^2(y)}{a}$$

(%t10) 
$$\text{mcs}_{3, 3, 2} = -\cos(y) \sin(y)$$

```
(%t11)
mcs          = ---
              d
              x
              2 a
(%o11)
done
```

## 30.2 Definitions for ctensor

### 30.2.1 Initialization and setup

**csetup** () Function  
 A function in the **ctensor** (component tensor) package which initializes the package and allows the user to enter a metric interactively. See **ctensor** for more details.

**cmetric** (*dis*) Function  
**cmetric** () Function

A function in the **ctensor** (component tensor) package that computes the metric inverse and sets up the package for further calculations.

If **cframe\_flag** is false, the function computes the inverse metric **ug** from the (user-defined) matrix **lg**. The metric determinant is also computed and stored in the variable **gdet**. Furthermore, the package determines if the metric is diagonal and sets the value of **diagmetric** accordingly. If the optional argument *dis* is present and not equal to false, the user is prompted to see the metric inverse.

If **cframe\_flag** is true, the function expects that the values of **fri** (the inverse frame matrix) and **lfg** (the frame metric) are defined. From these, the frame matrix **fr** and the inverse frame metric **ufg** are computed.

**ct\_coordsys** (*coordinate\_system, extra\_arg*) Function  
**ct\_coordsys** (*coordinate\_system*) Function

Sets up a predefined coordinate system and metric. The argument *coordinate\_system* can be one of the following symbols:

SYMBOL	Dim	Coordinates	Description/comments
<b>cartesian2d</b>	2	[x,y]	Cartesian 2D coordinate system
<b>polar</b>	2	[r,phi]	Polar coordinate system
<b>elliptic</b>	2	[u,v]	
<b>confocalelliptic</b>	2	[u,v]	
<b>bipolar</b>	2	[u,v]	
<b>parabolic</b>	2	[u,v]	
<b>cartesian3d</b>	3	[x,y,z]	Cartesian 3D coordinate system
<b>polarcylindrical</b>	3	[r,theta,z]	
<b>ellipticcylindrical</b>	3	[u,v,z]	Elliptic 2D with cylindrical Z
<b>confocalellipsoidal</b>	3	[u,v,w]	



bipolarcylindrical	3	[u,v,z]	Bipolar 2D with cylindrical Z
paraboliccylindrical	3	[u,v,z]	Parabolic 2D with cylindrical Z
paraboloidal	3	[u,v,phi]	
conical	3	[u,v,w]	
toroidal	3	[u,v,phi]	
spherical	3	[r,theta,phi]	Spherical coordinate system
oblatespheroidal	3	[u,v,phi]	
oblatespheroidalsqrt	3	[u,v,phi]	
prolatespheroidal	3	[u,v,phi]	
prolatespheroidalsqrt	3	[u,v,phi]	
ellipsoidal	3	[r,theta,phi]	
cartesian4d	4	[x,y,z,t]	Cartesian 4D coordinate system
spherical4d	4	[r,theta,eta,phi]	
exteriorschwarzschild	4	[t,r,theta,phi]	Schwarzschild metric
interiorschwarzschild	4	[t,z,u,v]	Interior Schwarzschild metric
kerr_newman	4	[t,r,theta,phi]	Charged axially symmetric metric

`coordinate_system` can also be a list of transformation functions, followed by a list containing the coordinate variables. For instance, you can specify a spherical metric as follows:

```
(%i1) load(ctensor);
(%o1) /share/tensor/ctensor.mac
(%i2) ct_coordsys([r*cos(theta)*cos(phi),r*cos(theta)*sin(phi),
r*sin(theta),[r,theta,phi]]);
(%o2) done
(%i3) lg:trigsimp(lg);
(%o3)
[ 1  0      0      ]
[                ]
[      2          ]
[ 0  r      0      ]
[                ]
[                2  2 ]
[ 0  0  r  cos(theta) ]

(%i4) ct_coords;
(%o4) [r, theta, phi]
(%i5) dim;
(%o5) 3
```

Transformation functions can also be used when `cframe_flag` is true:

```
(%i1) load(ctensor);
(%o1) /share/tensor/ctensor.mac
(%i2) cframe_flag:true;
(%o2) true
(%i3) ct_coordsys([r*cos(theta)*cos(phi),r*cos(theta)*sin(phi),
r*sin(theta),[r,theta,phi]]);
(%o3) done
```



### uriem

`ctensor` can also work using moving frames. When `cframe_flag` is set to `true`, the following tensors can be calculated:

```

lfg -- ufg
 \
fri -- fr -- lcs -- mcs -- lriem -- ric -- uric
  \
   lg -- ug
           | \
           |  weyl  \
           |   tracer - ein -- lein
           | \
           |  riem
           |
           | \
           |  uriem

```

### christof (*dis*)

Function

A function in the `ctensor` (component tensor) package. It computes the Christoffel symbols of both kinds. The argument *dis* determines which results are to be immediately displayed. The Christoffel symbols of the first and second kinds are stored in the arrays `lcs[i,j,k]` and `mcs[i,j,k]` respectively and defined to be symmetric in the first two indices. If the argument to `christof` is `lcs` or `mcs` then the unique non-zero values of `lcs[i,j,k]` or `mcs[i,j,k]`, respectively, will be displayed. If the argument is `all` then the unique non-zero values of `lcs[i,j,k]` and `mcs[i,j,k]` will be displayed. If the argument is `false` then the display of the elements will not occur. The array elements `mcs[i,j,k]` are defined in such a manner that the final index is contravariant.

### ricci (*dis*)

Function

A function in the `ctensor` (component tensor) package. `ricci` computes the covariant (symmetric) components `ric[i,j]` of the Ricci tensor. If the argument *dis* is `true`, then the non-zero components are displayed.

### uricci (*dis*)

Function

This function first computes the covariant components `ric[i,j]` of the Ricci tensor. Then the mixed Ricci tensor is computed using the contravariant metric tensor. If the value of the argument *dis* is `true`, then these mixed components, `uric[i,j]` (the index *i* is covariant and the index *j* is contravariant), will be displayed directly. Otherwise, `uricci(false)` will simply compute the entries of the array `uric[i,j]` without displaying the results.

### scurvature ()

Function

returns the scalar curvature (obtained by contracting the Ricci tensor) of the Riemannian manifold with the given metric.

**einstein** (*dis*)

Function

A function in the `ctensor` (component tensor) package. `einstein` computes the mixed Einstein tensor after the Christoffel symbols and Ricci tensor have been obtained (with the functions `christof` and `ricci`). If the argument `dis` is `true`, then the non-zero values of the mixed Einstein tensor `ein[i,j]` will be displayed where `j` is the contravariant index. The variable `rateinstein` will cause the rational simplification on these components. If `ratfac` is `true` then the components will also be factored.

**leinstein** (*dis*)

Function

Covariant Einstein-tensor. `leinstein` stores the values of the covariant Einstein tensor in the array `lein`. The covariant Einstein-tensor is computed from the mixed Einstein tensor `ein` by multiplying it with the metric tensor. If the argument `dis` is `true`, then the non-zero values of the covariant Einstein tensor are displayed.

**riemann** (*dis*)

Function

A function in the `ctensor` (component tensor) package. `riemann` computes the Riemann curvature tensor from the given metric and the corresponding Christoffel symbols. The following index conventions are used:

$$R[i,j,k,l] = R \begin{matrix} l \\ ijk \end{matrix} = \begin{matrix} \_l \\ ij,k \end{matrix} - \begin{matrix} \_l \\ ik,j \end{matrix} + \begin{matrix} \_l \ \_m \\ mk \ ij \end{matrix} - \begin{matrix} \_l \ \_m \\ mj \ ik \end{matrix}$$

This notation is consistent with the notation used by the `ITENSOR` package and its `icurvature` function. If the optional argument `dis` is `true`, the non-zero components `riem[i,j,k,l]` will be displayed. As with the Einstein tensor, various switches set by the user control the simplification of the components of the Riemann tensor. If `ratriemann` is `true`, then rational simplification will be done. If `ratfac` is `true` then each of the components will also be factored.

If the variable `cframe_flag` is `false`, the Riemann tensor is computed directly from the Christoffel-symbols. If `cframe_flag` is `false`, the covariant Riemann-tensor is computed first from the frame field coefficients.

**lriemann** (*dis*)

Function

Covariant Riemann-tensor (`lriem[]`).

Computes the covariant Riemann-tensor as the array `lriem`. If the argument `dis` is `true`, unique nonzero values are displayed.

If the variable `cframe_flag` is `true`, the covariant Riemann tensor is computed directly from the frame field coefficients. Otherwise, the (3,1) Riemann tensor is computed first.

For information on index ordering, see `riemann`.

**uriemann** (*dis*)

Function

Computes the contravariant components of the Riemann curvature tensor as array elements `uriem[i,j,k,l]`. These are displayed if `dis` is `true`.



```

(%i5) lg:matrix([-1,0,0,0],[0,1,0,0],[0,0,r^2,0],[0,0,0,r^2*sin(theta)^2]);
      [ - 1  0  0      0      ]
      [                               ]
      [  0  1  0      0      ]
      [                               ]
      [                               ]
(%o5) [                               ]
      [  0  0  r^2      0      ]
      [                               ]
      [                               ]
      [  0  0  0  r^2  sin^2(theta) ]
(%i6) h:matrix([h11,0,0,0],[0,h22,0,0],[0,0,h33,0],[0,0,0,h44]);
      [ h11  0  0  0 ]
      [                               ]
      [  0  h22  0  0 ]
      [                               ]
      [  0  0  h33  0 ]
      [                               ]
      [  0  0  0  h44 ]
(%i7) depends(l,r);
(%o7) [l(r)]
(%i8) lg:lg+l*h;
      [ h11 l - 1      0      0      0      ]
      [                               ]
      [  0      h22 l + 1      0      0      ]
      [                               ]
      [                               ]
(%o8) [  0      0      r^2 + h33 l      0      ]
      [                               ]
      [                               ]
      [  0      0      0      r^2 sin^2(theta) + h44 l ]
(%i9) cmetric(false);
(%o9) done
(%i10) einstein(false);
(%o10) done
(%i11) ntermst(ein);
[[1, 1], 62]
[[1, 2], 0]
[[1, 3], 0]
[[1, 4], 0]
[[2, 1], 0]
[[2, 2], 24]
[[2, 3], 0]
[[2, 4], 0]
[[3, 1], 0]
[[3, 2], 0]
[[3, 3], 46]
[[3, 4], 0]
[[4, 1], 0]
[[4, 2], 0]

```

```

[[4, 3], 0]
[[4, 4], 46]
(%o12) done

```

However, if we recompute this example as an approximation that is linear in the variable  $l$ , we get much simpler expressions:

```

(%i14) ctayswitch:true;
(%o14) true
(%i15) ctayvar:l;
(%o15) l
(%i16) ctaypov:1;
(%o16) 1
(%i17) ctaypt:0;
(%o17) 0
(%i18) christof(false);
(%o18) done
(%i19) ricci(false);
(%o19) done
(%i20) einstein(false);
(%o20) done
(%i21) ntermst(ein);
[[1, 1], 6]
[[1, 2], 0]
[[1, 3], 0]
[[1, 4], 0]
[[2, 1], 0]
[[2, 2], 13]
[[2, 3], 2]
[[2, 4], 0]
[[3, 1], 0]
[[3, 2], 2]
[[3, 3], 9]
[[3, 4], 0]
[[4, 1], 0]
[[4, 2], 0]
[[4, 3], 0]
[[4, 4], 9]
(%o21) done
(%i22) ratsimp(ein[1,1]);
(%o22) - (((h11 h22 - h11 ) (l ) r2 - 2 h33 l r2 ) sin (theta)
          r r
          - 2 h44 l r2 - h33 h44 (l ) ) / (4 r4 sin (theta))

```

This capability can be useful, for instance, when working in the weak field limit far from a gravitational source.

### 30.2.4 Frame fields

When the variable `cframe_flag` is set to true, the `ctensor` package performs its calculations using a moving frame.

**frame\_bracket** (*fr, fri, diagframe*)

Function

The frame bracket (`fb[]`).

Computes the frame bracket according to the following definition:

$$\text{ifb}_{ab} = \begin{pmatrix} c & c & c & d & e \\ \text{ifri} & -\text{ifri} & & \text{ifr} & \text{ifr} \\ d,e & e,d & a & b \end{pmatrix}$$

### 30.2.5 Algebraic classification

A new feature (as of November, 2004) of `ctensor` is its ability to compute the Petrov classification of a 4-dimensional spacetime metric. For a demonstration of this capability, see the file `share/tensor/petrov.dem`.

**nptetrad** ()

Function

Computes a Newman-Penrose null tetrad (`np`) and its raised-index counterpart (`npi`). See `petrov` for an example.

The null tetrad is constructed on the assumption that a four-dimensional orthonormal frame metric with metric signature  $(-,+,+,+)$  is being used. The components of the null tetrad are related to the inverse frame matrix as follows:

$$\begin{aligned} \text{np}_1 &= (\text{fri}_1 + \text{fri}_2) / \text{sqrt}(2) \\ \text{np}_2 &= (\text{fri}_1 - \text{fri}_2) / \text{sqrt}(2) \\ \text{np}_3 &= (\text{fri}_3 + \%i \text{fri}_4) / \text{sqrt}(2) \\ \text{np}_4 &= (\text{fri}_3 - \%i \text{fri}_4) / \text{sqrt}(2) \end{aligned}$$

**psi** (*dis*)

Function

Computes the five Newman-Penrose coefficients `psi[0]...psi[4]`. If `psi` is set to `true`, the coefficients are displayed. See `petrov` for an example.

These coefficients are computed from the Weyl-tensor in a coordinate base. If a frame base is used, the Weyl-tensor is first converted to a coordinate base, which can be a



computationally expensive procedure. For this reason, in some cases it may be more advantageous to use a coordinate base in the first place before the Weyl tensor is computed. Note however, that constructing a Newman-Penrose null tetrad requires a frame base. Therefore, a meaningful computation sequence may begin with a frame base, which is then used to compute `lg` (computed automatically by `cmetric` and then `ug`). At this point, you can switch back to a coordinate base by setting `cframe_flag` to false before beginning to compute the Christoffel symbols. Changing to a frame base at a later stage could yield inconsistent results, as you may end up with a mixed bag of tensors, some computed in a frame base, some in a coordinate base, with no means to distinguish between the two.

**petrov ()** Function

Computes the Petrov classification of the metric characterized by `psi[0]...psi[4]`.

For example, the following demonstrates how to obtain the Petrov-classification of the Kerr metric:

```
(%i1) load(ctensor);
(%o1) /share/tensor/ctensor.mac
(%i2) (cframe_flag:true,gcd:smod,ctrgsimp:true,ratfac:true);
(%o2) true
(%i3) ct_coordsys(exteriorschwarzschild,all);
(%o3) done
(%i4) ug:invert(lg)$
(%i5) weyl(false);
(%o5) done
(%i6) nptetrad(true);
(%t6) np =

[ sqrt(r - 2 m)          sqrt(r)
[ -----  -----  0      0
[ sqrt(2) sqrt(r)  sqrt(2) sqrt(r - 2 m)
[
[ sqrt(r - 2 m)          sqrt(r)
[ -----  -----  0      0
[ sqrt(2) sqrt(r)  sqrt(2) sqrt(r - 2 m)
[
[
[          r      %i r sin(theta)
[          0      0      -----  -----
[          sqrt(2)      sqrt(2)
[
[          r      %i r sin(theta)
[          0      0      -----  -----
[          sqrt(2)      sqrt(2)

(%t7) npi = matrix([- sqrt(r)          sqrt(r - 2 m)
                    sqrt(2) sqrt(r - 2 m)  sqrt(2) sqrt(r), 0, 0],
                    sqrt(r)          sqrt(r - 2 m)
```

```

[- -----, - -----, 0, 0],
  sqrt(2) sqrt(r - 2 m)   sqrt(2) sqrt(r)

[0, 0, -----, -----],
          1                %i
          sqrt(2) r        sqrt(2) r sin(theta)

[0, 0, -----, - -----])
          1                %i
          sqrt(2) r        sqrt(2) r sin(theta)

(%o7)                                     done
(%i7) psi(true);
(%t8)                                     psi = 0
                                           0

(%t9)                                     psi = 0
                                           1

(%t10)                                    psi = --
                                           2   3
                                           r

(%t11)                                    psi = 0
                                           3

(%t12)                                    psi = 0
                                           4
(%o12)                                     done
(%i12) petrov();
(%o12)                                     D

```

The Petrov classification function is based on the algorithm published in "Classifying geometries in general relativity: III Classification in practice" by Pollney, Skea, and d'Inverno, *Class. Quant. Grav.* 17 2885-2902 (2000). Except for some simple test cases, the implementation is untested as of December 19, 2004, and is likely to contain errors.

### 30.2.6 Torsion and nonmetricity

`ctensor` has the ability to compute and include torsion and nonmetricity coefficients in the connection coefficients.

The torsion coefficients are calculated from a user-supplied tensor `tr`, which should be a rank (2,1) tensor. From this, the torsion coefficients `kt` are computed according to the following formulae:

$$k^m_{\quad n} = \Gamma^m_{\quad [n\quad p]} - \Gamma^m_{\quad [p\quad n]}$$

$$kt_{ijk} = \frac{-g_{im} tr_{kj} - g_{jm} tr_{ki} - tr_{ij} g_{km}}{2}$$

$$kt_{ij} = g_{ij} - kt_{ijm}$$

Note that only the mixed-index tensor is calculated and stored in the array `kt`.

The nonmetricity coefficients are calculated from the user-supplied nonmetricity vector `nm`. From this, the nonmetricity coefficients `nmc` are computed as follows:

$$nmc_{ij} = \frac{-nm_{ik} D_{kj} - D_{ij} nm_{km} + g_{im} nm_{kj}}{2}$$

where `D` stands for the Kronecker-delta.

When `ctorsion_flag` is set to `true`, the values of `kt` are subtracted from the mixed-indexed connection coefficients computed by `christof` and stored in `mcs`. Similarly, if `cnonmet_flag` is set to `true`, the values of `nmc` are subtracted from the mixed-indexed connection coefficients.

If necessary, `christof` calls the functions `contortion` and `nonmetricity` in order to compute `kt` and `nm`.

**contortion** (*tr*) Function  
 Computes the (2,1) contortion coefficients from the torsion tensor *tr*.

**nonmetricity** (*nm*) Function  
 Computes the (2,1) nonmetricity coefficients from the nonmetricity vector *nm*.

### 30.2.7 Miscellaneous features

**ctransform** (*M*) Function  
 A function in the `ctensor` (component tensor) package which will perform a coordinate transformation upon an arbitrary square symmetric matrix *M*. The user must input the functions which define the transformation. (Formerly called `transform`.)

**findde** (*A, n*) Function  
 returns a list of the unique differential equations (expressions) corresponding to the elements of the *n* dimensional square array *A*. Presently, *n* may be 2 or 3. `deindex` is a global list containing the indices of *A* corresponding to these unique differential

equations. For the Einstein tensor (`ein`), which is a two dimensional array, if computed for the metric in the example below, `findde` gives the following independent differential equations:

```
(%i1) load(ctensor);
(%o1)      /share/tensor/ctensor.mac
(%i2) derivabbrev:true;
(%o2)
           true
(%i3) dim:4;
(%o3)
           4
(%i4) lg:matrix([a,0,0,0],[0,x^2,0,0],[0,0,x^2*sin(y)^2,0],[0,0,0,-d]);
           [ a  0    0    0 ]
           [          ]
           [    2          ]
(%o4)      [ 0  x    0    0 ]
           [          ]
           [          2    2 ]
           [ 0  0  x  sin (y)  0 ]
           [          ]
           [ 0  0    0    - d ]

(%i5) depends([a,d],x);
(%o5)
           [a(x), d(x)]
(%i6) ct_coords:[x,y,z,t];
(%o6)
           [x, y, z, t]
(%i7) cmetric();
(%o7)
           done
(%i8) einstein(false);
(%o8)
           done
(%i9) findde(ein,2);
(%o9) [d x - a d + d, 2 a d d x - a (d ) x - a d d x + 2 a d d
           x          x x          x          x          x          x
           - 2 a d , a x + a - a]
           x          x

(%i10) deindex;
(%o10)
           [[1, 1], [2, 2], [4, 4]]
```

### **cograd ()**

Function

Computes the covariant gradient of a scalar function allowing the user to choose the corresponding vector name as the example under `contragrad` illustrates.

### **contragrad ()**

Function

Computes the contravariant gradient of a scalar function allowing the user to choose the corresponding vector name as the example below for the Schwarzschild metric illustrates:

```

(%i1) load(ctensor);
(%o1)      /share/tensor/ctensor.mac
(%i2) derivabbrev:true;
(%o2)      true
(%i3) ct_coordsys(exteriorschwarzschild,all);
(%o3)      done
(%i4) depends(f,r);
(%o4)      [f(r)]
(%i5) cograd(f,g1);
(%o5)      done
(%i6) listarray(g1);
(%o6)      [0, f , 0, 0]
           r
(%i7) contragrad(f,g2);
(%o7)      done
(%i8) listarray(g2);
           f r - 2 f m
           r      r
(%o8)      [0, -----, 0, 0]
           r

```

**dscalar ()**

Function

computes the tensor d'Alembertian of the scalar function once dependencies have been declared upon the function. For example:

```

(%i1) load(ctensor);
(%o1)      /share/tensor/ctensor.mac
(%i2) derivabbrev:true;
(%o2)      true
(%i3) ct_coordsys(exteriorschwarzschild,all);
(%o3)      done
(%i4) depends(p,r);
(%o4)      [p(r)]
(%i5) factor(dscalar(p));
           2
           p  r  - 2 m p  r + 2 p  r - 2 m p
           r r      r r      r      r
(%o5)      -----
           2
           r

```

**checkdiv ()**

Function

computes the covariant divergence of the mixed second rank tensor (whose first index must be covariant) by printing the corresponding  $n$  components of the vector field (the divergence) where  $n = \text{dim}$ . If the argument to the function is  $\mathbf{g}$  then the divergence of the Einstein tensor will be formed and must be zero. In addition, the divergence (vector) is given the array name `div`.

**cgeodesic** (*dis*) Function

A function in the `ctensor` (component tensor) package. `cgeodesic` computes the geodesic equations of motion for a given metric. They are stored in the array `geod[i]`. If the argument *dis* is `true` then these equations are displayed.

**bdvac** (*f*) Function

generates the covariant components of the vacuum field equations of the Brans- Dicke gravitational theory. The scalar field is specified by the argument *f*, which should be a (quoted) function name with functional dependencies, e.g., `'p(x)`.

The components of the second rank covariant field tensor are represented by the array `bd`.

**invariant1** () Function

generates the mixed Euler- Lagrange tensor (field equations) for the invariant density of  $R^2$ . The field equations are the components of an array named `inv1`.

**invariant2** () Function

\*\*\* NOT YET IMPLEMENTED \*\*\*

generates the mixed Euler- Lagrange tensor (field equations) for the invariant density of `ric[i,j]*uriem[i,j]`. The field equations are the components of an array named `inv2`.

**bimetric** () Function

\*\*\* NOT YET IMPLEMENTED \*\*\*

generates the field equations of Rosen's bimetric theory. The field equations are the components of an array named `rosen`.

### 30.2.8 Utility functions

**diagmatrixp** (*M*) Function

Returns `true` if *M* is a diagonal matrix or (2D) array.

**symmetricp** (*M*) Function

Returns `true` if *M* is a symmetric matrix or (2D) array.

**ntermst** (*f*) Function

gives the user a quick picture of the "size" of the doubly subscripted tensor (array) *f*. It prints two element lists where the second element corresponds to `NTERMS` of the components specified by the first elements. In this way, it is possible to quickly find the non-zero expressions and attempt simplification.

**cdisplay** (*ten*) Function

displays all the elements of the tensor *ten*, as represented by a multidimensional array. Tensors of rank 0 and 1, as well as other types of variables, are displayed as with `ldisplay`. Tensors of rank 2 are displayed as 2-dimensional matrices, while tensors of higher rank are displayed as a list of 2-dimensional matrices. For instance, the Riemann-tensor of the Schwarzschild metric can be viewed as:

```
(%i1) load(ctensor);
(%o1) /share/tensor/ctensor.mac
(%i2) ratfac:true;
(%o2) true
(%i3) ct_coordsys(exteriorschwarzschild,all);
(%o3) done
(%i4) riemann(false);
(%o4) done
(%i5) cdisplay(riem);
```

$$\begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & -\frac{3m(r-2m)}{4r^4} + \frac{m}{3r^3} - \frac{2m}{4r^4} & 0 & 0 \\ 0 & 0 & \frac{m(r-2m)}{4r} & 0 \\ 0 & 0 & 0 & \frac{m(r-2m)}{4r} \end{bmatrix}$$

```
riem
1, 1
```

$$\begin{bmatrix} 2m(r-2m) & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$

```
riem
1, 2
```

$$\begin{bmatrix} m(r-2m) & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \end{bmatrix}$$

```
riem
1, 3
```

$$\begin{aligned}
 \text{riem}_{1,4} &= \begin{bmatrix} m(r-2m) \\ 0 & 0 & 0 & -\frac{\quad}{4} \\ r \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}
 \end{aligned}$$

$$\begin{aligned}
 \text{riem}_{2,1} &= \begin{bmatrix} 0 & 0 & 0 & 0 \\ 2m \\ -\frac{\quad}{2} & 0 & 0 & 0 \\ r & (r-2m) \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}
 \end{aligned}$$

$$\begin{aligned}
 \text{riem}_{2,2} &= \begin{bmatrix} 2m \\ -\frac{\quad}{2} & 0 & 0 & 0 \\ r & (r-2m) \\ 0 & 0 & 0 & 0 \\ 0 & 0 & -\frac{m}{2} & 0 \\ \quad & r & (r-2m) \\ 0 & 0 & 0 & -\frac{m}{2} \\ \quad & r & (r-2m) \end{bmatrix}
 \end{aligned}$$

$$\begin{aligned}
 \text{riem}_{2,3} &= \begin{bmatrix} 0 & 0 & 0 & 0 \\ m \\ 0 & 0 & -\frac{\quad}{2} & 0 \\ r & (r-2m) \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix}
 \end{aligned}$$



$$\text{riem}_{2,4} = \begin{bmatrix} 0 & 0 & 0 & 0 \\ [ & & & ] \\ [ & & m & ] \\ [ 0 & 0 & 0 & \text{-----} ] \\ [ & & 2 & ] \\ [ & & r & (r - 2 m) ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \end{bmatrix}$$

$$\text{riem}_{3,1} = \begin{bmatrix} [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ m & & & ] \\ [ - & 0 & 0 & 0 ] \\ [ r & & & ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \end{bmatrix}$$

$$\text{riem}_{3,2} = \begin{bmatrix} [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ m & & & ] \\ [ 0 & - & 0 & 0 ] \\ [ r & & & ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \end{bmatrix}$$

$$\text{riem}_{3,3} = \begin{bmatrix} [ m & & & ] \\ [ - & - & 0 & 0 & 0 ] \\ [ r & & & ] \\ [ & & & ] \\ [ m & & & ] \\ [ 0 & - & - & 0 & 0 ] \\ [ r & & & ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ & & & ] \\ [ 0 & 0 & 0 & \frac{2 m - r}{r} + 1 ] \\ [ & & & ] \end{bmatrix}$$

$$\begin{bmatrix} [ 0 & 0 & 0 & 0 ] \\ [ & & & ] \\ [ 0 & 0 & 0 & 0 ] \end{bmatrix}$$

$$\text{riem}_{3,4} = \begin{bmatrix} [ & & & & ] \\ [ & & & 2 m & ] \\ [ 0 & 0 & 0 & - & ] \\ [ & & & r & ] \\ [ & & & & ] \\ [ 0 & 0 & 0 & 0 & ] \end{bmatrix}$$

$$\text{riem}_{4,1} = \begin{bmatrix} [ & & 0 & 0 & 0 & 0 ] \\ [ & & & & & ] \\ [ & & 0 & 0 & 0 & 0 ] \\ [ & & & & & ] \\ [ & & 0 & 0 & 0 & 0 ] \\ [ & & & & & ] \\ [ & & 2 & & & ] \\ [ m \sin(\theta) & & & & & ] \\ [ - & & & & & ] \\ [ & & r & & & ] \end{bmatrix}$$

$$\text{riem}_{4,2} = \begin{bmatrix} [ 0 & & 0 & 0 & 0 ] \\ [ & & & & ] \\ [ 0 & & 0 & 0 & 0 ] \\ [ & & & & ] \\ [ 0 & & 0 & 0 & 0 ] \\ [ & & & & ] \\ [ & & 2 & & ] \\ [ m \sin(\theta) & & & & ] \\ [ 0 & - & & & 0 & 0 ] \\ [ & & r & & & ] \end{bmatrix}$$

$$\text{riem}_{4,3} = \begin{bmatrix} [ 0 & 0 & & 0 & 0 ] \\ [ & & & & ] \\ [ 0 & 0 & & 0 & 0 ] \\ [ & & & & ] \\ [ 0 & 0 & & 0 & 0 ] \\ [ & & & & ] \\ [ & & & 2 & ] \\ [ & & & 2 m \sin(\theta) & ] \\ [ 0 & 0 & - & & 0 ] \\ [ & & & r & ] \end{bmatrix}$$

$$\text{riem}_{4,4} = \begin{bmatrix} [ & & 2 & & ] \\ [ m \sin(\theta) & & & & ] \\ [ - & & & & 0 & 0 ] \\ [ & & r & & & ] \\ [ & & & & & ] \\ [ & & & 2 & & ] \\ [ & & & m \sin(\theta) & & ] \\ [ 0 & - & & & 0 & 0 ] \\ [ & & & r & & ] \end{bmatrix}$$

```

          [
          [
          [      2
          [      0      0      -----  0 ]
          [      r
          [      0      0      0      0 ]
          ]
          ]
          ]
(%o5)
done
```

**deleten** (*L*, *n*)

Function

Returns a new list consisting of *L* with the *n*'th element deleted.

**30.2.9 Variables used by ctensor****dim**

Variable

Default value: 4

An option in the **ctensor** (component tensor) package. **dim** is the dimension of the manifold with the default 4. The command **dim: n** will reset the dimension to any other value **n**.

**diagmetric**

Variable

Default value: **false**

An option in the **ctensor** (component tensor) package. If **diagmetric** is **true** special routines compute all geometrical objects (which contain the metric tensor explicitly) by taking into consideration the diagonality of the metric. Reduced run times will, of course, result. Note: this option is set automatically by **csetup** if a diagonal metric is specified.

**ctrjsimp**

Variable

Causes trigonometric simplifications to be used when tensors are computed. Presently, **ctrjsimp** affects only computations involving a moving frame.

**cframe\_flag**

Variable

Causes computations to be performed relative to a moving frame as opposed to a holonomic metric. The frame is defined by the inverse frame array **fri** and the frame metric **lfg**. For computations using a Cartesian frame, **lfg** should be the unit matrix of the appropriate dimension; for computations in a Lorentz frame, **lfg** should have the appropriate signature.

**ctorsion\_flag**

Variable

Causes the contortion tensor to be included in the computation of the connection coefficients. The contortion tensor itself is computed by **contortion** from the user-supplied tensor **tr**.

- cnonmet\_flag** Variable  
 Causes the nonmetricity coefficients to be included in the computation of the connection coefficients. The nonmetricity coefficients are computed from the user-supplied nonmetricity vector `nm` by the function `nonmetricity`.
- ctayswitch** Variable  
 If set to `true`, causes some `ctensor` computations to be carried out using Taylor-series expansions. Presently, `christof`, `ricci`, `uricci`, `einstein`, and `weyl` take into account this setting.
- ctayvar** Variable  
 Variable used for Taylor-series expansion if `ctayswitch` is set to `true`.
- ctaypov** Variable  
 Maximum power used in Taylor-series expansion when `ctayswitch` is set to `true`.
- ctaypt** Variable  
 Point around which Taylor-series expansion is carried out when `ctayswitch` is set to `true`.
- gdet** Variable  
 The determinant of the metric tensor `lg`. Computed by `cmetric` when `cframe_flag` is set to `false`.
- ratchristof** Variable  
 Causes rational simplification to be applied by `christof`.
- rateinstein** Variable  
 Default value: `true`  
 If `true` rational simplification will be performed on the non-zero components of Einstein tensors; if `ratfac` is `true` then the components will also be factored.
- ratriemann** Variable  
 Default value: `true`  
 One of the switches which controls simplification of Riemann tensors; if `true`, then rational simplification will be done; if `ratfac` is `true` then each of the components will also be factored.
- ratweyl** Variable  
 Default value: `true`  
 If `true`, this switch causes the `weyl` function to apply rational simplification to the values of the Weyl tensor. If `ratfac` is `true`, then the components will also be factored.

<b>lfg</b>	Variable
The covariant frame metric. By default, it is initialized to the 4-dimensional Lorentz frame with signature (+,+,+,-). Used when <code>cframe_flag</code> is <code>true</code> .	
<b>ufg</b>	Variable
The inverse frame metric. Computed from <code>lfg</code> when <code>cmetric</code> is called while <code>cframe_flag</code> is set to <code>true</code> .	
<b>riem</b>	Variable
The (3,1) Riemann tensor. Computed when the function <code>riemann</code> is invoked. For information about index ordering, see the description of <code>riemann</code> . if <code>cframe_flag</code> is <code>true</code> , <code>riem</code> is computed from the covariant Riemann-tensor <code>lriem</code> .	
<b>lriem</b>	Variable
The covariant Riemann tensor. Computed by <code>lriemann</code> .	
<b>uriem</b>	Variable
The contravariant Riemann tensor. Computed by <code>uriemann</code> .	
<b>ric</b>	Variable
The mixed Ricci-tensor. Computed by <code>ricci</code> .	
<b>uric</b>	Variable
The contravariant Ricci-tensor. Computed by <code>uricci</code> .	
<b>lg</b>	Variable
The metric tensor. This tensor must be specified (as a <code>dim</code> by <code>dim</code> matrix) before other computations can be performed.	
<b>ug</b>	Variable
The inverse of the metric tensor. Computed by <code>cmetric</code> .	
<b>weyl</b>	Variable
The Weyl tensor. Computed by <code>weyl</code> .	
<b>fb</b>	Variable
Frame bracket coefficients, as computed by <code>frame_bracket</code> .	
<b>kinvariant</b>	Variable
The Kretschmann invariant. Computed by <code>rinvariant</code> .	
<b>np</b>	Variable
A Newman-Penrose null tetrad. Computed by <code>nptetrad</code> .	

<b>npi</b>	Variable
The raised-index Newman-Penrose null tetrad. Computed by <code>nptetrad</code> . Defined as <code>ug.np</code> . The product <code>np.transpose(npi)</code> is constant:	
<pre>(%i39) trigsimp(np.transpose(npi));           [ 0  -1  0  0 ]           [          ]           [ -1  0  0  0 ] (%o39)    [          ]           [ 0  0  0  1 ]           [          ]           [ 0  0  1  0 ]</pre>	
<b>tr</b>	Variable
User-supplied rank-3 tensor representing torsion. Used by <code>contortion</code> .	
<b>kt</b>	Variable
The contortion tensor, computed from <code>tr</code> by <code>contortion</code> .	
<b>nm</b>	Variable
User-supplied nonmetricity vector. Used by <code>nonmetricity</code> .	
<b>nmc</b>	Variable
The nonmetricity coefficients, computed from <code>nm</code> by <code>nonmetricity</code> .	
<b>tensorkill</b>	Variable
Variable indicating if the tensor package has been initialized. Set and used by <code>csetup</code> , reset by <code>init_ctensor</code> .	
<b>ct_coords</b>	Variable
Default value: <code>[]</code>	
An option in the <code>ctensor</code> (component tensor) package. <code>ct_coords</code> contains a list of coordinates. While normally defined when the function <code>csetup</code> is called, one may redefine the coordinates with the assignment <code>ct_coords: [j1, j2, ..., jn]</code> where the <code>j</code> 's are the new coordinate names. See also <code>csetup</code> .	

### 30.2.10 Reserved names

The following names are used internally by the `ctensor` package and should not be redefined:

Name	Description
-----	
<code>_lg()</code>	Evaluates to <code>lfg</code> if frame metric used, <code>lg</code> otherwise
<code>_ug()</code>	Evaluates to <code>ufg</code> if frame metric used, <code>ug</code> otherwise
<code>cleanup()</code>	Removes items from the deindex list
<code>contract4()</code>	Used by <code>psi()</code>
<code>filemet()</code>	Used by <code>csetup()</code> when reading the metric from a file
<code>findde1()</code>	Used by <code>findde()</code>

findde2()	Used by findde()
findde3()	Used by findde()
kdelt()	Kronecker-delta (not generalized)
newmet()	Used by csetup() for setting up a metric interactively
setflags()	Used by init_ctensor()
readvalue()	
resimp()	
sermet()	Used by csetup() for entering a metric as Taylor-series
txyzsum()	
tmetric()	Frame metric, used by cmetric() when cframe_flag:true
triemann()	Riemann-tensor in frame base, used when cframe_flag:true
tricci()	Ricci-tensor in frame base, used when cframe_flag:true
trrc()	Ricci rotation coefficients, used by christof()
yesp()	

### 30.2.11 Changes

In November, 2004, the `ctensor` package was extensively rewritten. Many functions and variables have been renamed in order to make the package compatible with the commercial version of Macsyma.

New Name	Old Name	Description
ctaylor()	DLGTAYLOR()	Taylor-series expansion of an expression
lgeod[]	EM	Geodesic equations
ein[]	G[]	Mixed Einstein-tensor
ric[]	LR[]	Mixed Ricci-tensor
ricci()	LRICCOM()	Compute the mixed Ricci-tensor
ctaypov	MINP	Maximum power in Taylor-series expansion
cgeodesic()	MOTION	Compute geodesic equations
ct_coords	OMEGA	Metric coordinates
ctayvar	PARAM	Taylor-series expansion variable
lriem[]	R[]	Covariant Riemann-tensor
uriemann()	RAISERIEMANN()	Compute the contravariant Riemann-tensor
ratriemann	RATRIEMAN	Rational simplification of the Riemann-tensor
uric[]	RICCI[]	Contravariant Ricci-tensor
uricci()	RICCOM()	Compute the contravariant Ricci-tensor
cmetric()	SETMETRIC()	Set up the metric
ctaypt	TAYPT	Point for Taylor-series expansion
ctayswitch	TAYSWITCH	Taylor-series setting switch
csetup()	TSETUP()	Start interactive setup session
ctransform()	TTRANSFORM()	Interactive coordinate transformation
uriem[]	UR[]	Contravariant Riemann-tensor
weyl[]	W[]	(3,1) Weyl-tensor





```
(%o10)
[
[
[      1      2      1      2 ]
[
[ v      - 1      v . v      - v ]
[      1      1      2      2 ]
[
[ v      - v . v      - 1      v ]
[      2      1      2      1 ]
[
[ v . v      v      - v      - 1 ]
[      1      2      2      1 ]
]
```

`atensor` recognizes as base vectors indexed symbols, where the symbol is that stored in `asymbol` and the index runs between 1 and `adim`. For indexed symbols, and indexed symbols only, the bilinear forms `sf`, `af`, and `av` are evaluated. The evaluation substitutes the value of `aform[i,j]` in place of `fun(v[i],v[j])` where `v` represents the value of `asymbol` and `fun` is either `af` or `sf`; or, it substitutes `v[aform[i,j]]` in place of `av(v[i],v[j])`.

Needless to say, the functions `sf`, `af` and `av` can be redefined.

When the `atensor` package is loaded, the following flags are set:

```
dotscrules:true;
dotdistrib:true;
dotexptsimp:false;
```

If you wish to experiment with a nonassociative algebra, you may also consider setting `dotassoc` to `false`. In this case, however, `atensimp` will not always be able to obtain the desired simplifications.

## 31.2 Definitions for `atensor`

**init\_atensor** (*alg\_type*, *opt\_dims*)

Function

**init\_atensor** (*alg\_type*)

Function

Initializes the `atensor` package with the specified algebra type. *alg\_type* can be one of the following:

**universal**: The universal algebra has no commutation rules.

**grassmann**: The Grassman algebra is defined by the commutation relation  $u.v+v.u=0$ .

**clifford**: The Clifford algebra is defined by the commutation relation  $u.v+v.u=-2*sf(u,v)$  where `sf` is a symmetric scalar-valued function. For this algebra, *opt\_dims* can be up to three nonnegative integers, representing the number of positive, degenerate, and negative dimensions of the algebra, respectively. If any *opt\_dims* values are supplied, `atensor` will configure the values of `adim` and `aform` appropriately. Otherwise, `adim` will default to 0 and `aform` will not be defined.

**symmetric**: The symmetric algebra is defined by the commutation relation  $u.v-v.u=0$ .

**symplectic**: The symplectic algebra is defined by the commutation relation  $u.v-v.u=2*af(u,v)$  where `af` is an antisymmetric scalar-valued function. For the symplectic algebra, *opt\_dims* can be up to two nonnegative integers, representing the nondegenerate and degenerate dimensions, respectively. If any *opt\_dims* values are

supplied, `atensor` will configure the values of `adim` and `aform` appropriately. Otherwise, `adim` will default to 0 and `aform` will not be defined.

`lie_envelop`: The algebra of the Lie envelope is defined by the commutation relation  $u.v - v.u = 2*av(u,v)$  where `av` is an antisymmetric function.

The `init_atensor` function also recognizes several predefined algebra types:

`complex` implements the algebra of complex numbers as the Clifford algebra  $Cl(0,1)$ . The call `init_atensor(complex)` is equivalent to `init_atensor(clifford,0,0,1)`.

`quaternion` implements the algebra of quaternions. The call `init_atensor(quaternion)` is equivalent to `init_atensor(clifford,0,0,2)`.

`pauli` implements the algebra of Pauli-spinors as the Clifford-algebra  $Cl(3,0)$ . A call to `init_atensor(pauli)` is equivalent to `init_atensor(clifford,3)`.

`dirac` implements the algebra of Dirac-spinors as the Clifford-algebra  $Cl(3,1)$ . A call to `init_atensor(dirac)` is equivalent to `init_atensor(clifford,3,0,1)`.

**atensimp** (*expr*) Function

Simplifies an algebraic tensor expression *expr* according to the rules configured by a call to `init_atensor`. Simplification includes recursive application of commutation relations and resolving calls to `sf`, `af`, and `av` where applicable. A safeguard is used to ensure that the function always terminates, even for complex expressions.

**alg\_type** Function

The algebra type. Valid values are `universal`, `grassmann`, `clifford`, `symmetric`, `symplectic` and `lie_envelop`.

**adim** Variable

The dimensionality of the algebra. `atensor` uses the value of `adim` to determine if an indexed object is a valid base vector. Defaults to 0.

**aform** Variable

Default values for the bilinear forms `sf`, `af`, and `av`. The default is the identity matrix `ident(3)`.

**asymbol** Variable

The symbol for base vectors. Defaults to `v`.

**sf** (*u*, *v*) Function

A symmetric scalar function that is used in commutation relations. The default implementation checks if both arguments are base vectors using `abasep` and if that is the case, substitutes the corresponding value from the matrix `aform`.

**af** (*u*, *v*) Function

An antisymmetric scalar function that is used in commutation relations. The default implementation checks if both arguments are base vectors using `abasep` and if that is the case, substitutes the corresponding value from the matrix `aform`.

**av** (*u*, *v*) Function

An antisymmetric function that is used in commutation relations. The default implementation checks if both arguments are base vectors using **abasep** and if that is the case, substitutes the corresponding value from the matrix **aform**.

For instance:

```
(%i1) load(atensor);
(%o1)      /share/tensor/atensor.mac
(%i2) adim:3;
(%o2)      3
(%i3) aform:matrix([0,3,2],[3,0,1],[2,1,0]);
(%o3)      [ 0 3 2 ]
           [      ]
           [ 3 0 1 ]
           [      ]
           [ 2 1 0 ]

(%i4) asymbol:x;
(%o4)      x
(%i5) av(x[1],x[2]);
(%o5)      x
           3
```

**abasep** (*v*) Function

Checks if its argument is an **atensor** base vector. That is, if it is an indexed symbol, with the symbol being the same as the value of **asymbol**, and the index having a numeric value between 1 and **adim**.

## 32 Series

### 32.1 Introduction to Series

Maxima contains functions `taylor` and `powerseries` for finding the series of differentiable functions. It also has tools such as `nusum` capable of finding the closed form of some series. Operations such as addition and multiplication work as usual on series. This section presents the various global variables which control the expansion.

### 32.2 Definitions for Series

#### `cauchysum`

Variable

Default value: `false`

When multiplying together sums with `inf` as their upper limit, if `sumexpand` is `true` and `cauchysum` is `true` then the Cauchy product will be used rather than the usual product. In the Cauchy product the index of the inner summation is a function of the index of the outer one rather than varying independently.

Example:

```
(%i1) sumexpand: false$
(%i2) cauchysum: false$
(%i3) s: sum (f(i), i, 0, inf) * sum (g(j), j, 0, inf);
      inf      inf
      =====
      \        \
      ( >  f(i)) >  g(j)
      /        /
      =====
      i = 0    j = 0

(%o3)

(%i4) sumexpand: true$
(%i5) cauchysum: true$
(%i6) ''s;
      inf      i1
      =====
      \        \
      >        >    g(i1 - i2) f(i2)
      /        /
      =====
      i1 = 0 i2 = 0

(%o6)
```

#### `deftaylor` ( $f_1(x_1), \text{expr}_1, \dots, f_n(x_n), \text{expr}_n$ )

Function

For each function  $f_i$  of one variable  $x_i$ , `deftaylor` defines  $\text{expr}_i$  as the Taylor series about zero.  $\text{expr}_i$  is typically a polynomial in  $x_i$  or a summation; more general expressions are accepted by `deftaylor` without complaint.

`powerseries (f_i(x_i), x_i, 0)` returns the series defined by `deftaylor`.

`deftaylor` returns a list of the functions  $f_1, \dots, f_n$ . `deftaylor` evaluates its arguments.

Example:

```
(%i1) defaylor (f(x), x^2 + sum(x^i/(2^i*i!^2), i, 4, inf));
(%o1) [f]
(%i2) powerseries (f(x), x, 0);
      inf
      ====
      \      i1
      >      x      2
      /      ----- + x
      /      i1      2
      ==== 2  i1!
      i1 = 4
(%i3) taylor (exp (sqrt (f(x))), x, 0, 4);
      2      3      4
      x      3073 x      12817 x
(%o3)/T/ 1 + x + -- + ----- + ----- + . . .
          2      18432      307200
```

### **maxtayorder**

Variable

Default value: true

When `maxtayorder` is true, then during algebraic manipulation of (truncated) Taylor series, `taylor` tries to retain as many terms as are known to be correct.

### **niceindices** (*expr*)

Function

Renames the indices of sums and products in *expr*. `niceindices` attempts to rename each index to the value of `niceindicespref[1]`, unless that name appears in the summand or multiplicand, in which case `niceindices` tries the succeeding elements of `niceindicespref` in turn, until an unused variable is found. If the entire list is exhausted, additional indices are constructed by appending integers to the value of `niceindicespref[1]`, e.g., `i0, i1, i2, ...`

`niceindices` returns an expression. `niceindices` evaluates its argument.

Example:

```
(%i1) niceindicespref;
(%o1) [i, j, k, l, m, n]
(%i2) product (sum (f (foo + i*j*bar), foo, 1, inf), bar, 1, inf);
      inf      inf
      /====\  ====
      ! !    \
      ! !    >      f(bar i j + foo)
      ! !    /
      bar = 1  ====
              foo = 1
(%i3) niceindices (%);
      inf      inf
      /====\  ====
      ! !    \
```

```
(%o3)          !! > f(i j l + k)
              !! /
              l = 1 =====
              k = 1
```

**niceindicespref**

Variable

Default value: [i, j, k, l, m, n]

`niceindicespref` is the list from which `niceindices` takes the names of indices for sums and products.

The elements of `niceindicespref` are typically names of variables, although that is not enforced by `niceindices`.

Example:

```
(%i1) niceindicespref: [p, q, r, s, t, u]$
(%i2) product (sum (f (foo + i*j*bar), foo, 1, inf), bar, 1, inf);
          inf  inf
          /====\ =====
          !! \
(%o2)          !! > f(bar i j + foo)
          !! /
          bar = 1 =====
          foo = 1
(%i3) niceindices (%);
          inf  inf
          /====\ =====
          !! \
(%o3)          !! > f(i j q + p)
          !! /
          q = 1 =====
          p = 1
```

**nusum** (*expr*, *x*, *i\_0*, *i\_1*)

Function

Carries out indefinite hypergeometric summation of *expr* with respect to *x* using a decision procedure due to R.W. Gosper. *expr* and the result must be expressible as products of integer powers, factorials, binomials, and rational functions.

The terms "definite" and "indefinite summation" are used analogously to "definite" and "indefinite integration". To sum indefinitely means to give a symbolic result for the sum over intervals of variable length, not just e.g. 0 to inf. Thus, since there is no formula for the general partial sum of the binomial series, `nusum` can't do it.

`nusum` and `unsum` know a little about sums and differences of finite products. See also `unsum`.

Examples:

```
(%i1) nusum (n*n!, n, 0, n);
```

Dependent equations eliminated: (1)

```
(%o1)          (n + 1)! - 1
```

```
(%i2) nusum (n^4*4^n/binomial(2*n,n), n, 0, n);
```

```

      4      3      2      n
      2 (n + 1) (63 n + 112 n + 18 n - 22 n + 3) 4      2
(%o2) ----- - -----
      693 binomial(2 n, n)                        3 11 7
(%i3) unsum (% , n);

      4 n
      n 4
      -----
      binomial(2 n, n)
(%i4) unsum (prod (i^2, i, 1, n), n);
      n - 1
      /===\
      ! ! 2
(%o4) ( ! ! i ) (n - 1) (n + 1)
      ! !
      i = 1
(%i5) nusum (% , n, 1, n);

Dependent equations eliminated: (2 3)
      n
      /===\
      ! ! 2
(%o5) ! ! i - 1
      ! !
      i = 1

```

**pade** (*taylor\_series*, *numer\_deg\_bound*, *denom\_deg\_bound*) Function

Returns a list of all rational functions which have the given Taylor series expansion where the sum of the degrees of the numerator and the denominator is less than or equal to the truncation level of the power series, i.e. are "best" approximants, and which additionally satisfy the specified degree bounds.

*taylor\_series* is a univariate Taylor series. *numer\_deg\_bound* and *denom\_deg\_bound* are positive integers specifying degree bounds on the numerator and denominator.

*taylor\_series* can also be a Laurent series, and the degree bounds can be `inf` which causes all rational functions whose total degree is less than or equal to the length of the power series to be returned. Total degree is defined as *numer\_deg\_bound* + *denom\_deg\_bound*. Length of a power series is defined as "truncation level" + 1 - `min(0, "order of series")`.

```

(%i1) taylor (1 + x + x^2 + x^3, x, 0, 3);
      2      3
(%o1)/T/      1 + x + x + x + . . .
(%i2) pade (% , 1, 1);
      1
(%o2)      [- -----]
      x - 1
(%i3) t: taylor(-(83787*x^10 - 45552*x^9 - 187296*x^8
      + 387072*x^7 + 86016*x^6 - 1507328*x^5
      + 1966080*x^4 + 4194304*x^3 - 25165824*x^2

```

```

+ 67108864*x - 134217728)
/134217728, x, 0, 10);
      2      3      4      5      6      7
      x  3 x  x  15 x  23 x  21 x  189 x
(%o3)/T/ 1 - - + ---- - - - - ---- + ---- - ---- - ----
      2    16   32  1024  2048  32768  65536

      8      9      10
      5853 x  2847 x  83787 x
+ ---- + ---- - ---- + . . .
 4194304  8388608  134217728

(%i4) pade (t, 4, 4);
(%o4) []

```

There is no rational function of degree 4 numerator/denominator, with this power series expansion. You must in general have degree of the numerator and degree of the denominator adding up to at least the degree of the power series, in order to have enough unknown coefficients to solve.

```

(%i5) pade (t, 5, 5);
(%o5) [- (520256329 x5 - 96719020632 x4 - 489651410240 x3
- 1619100813312 x2 - 2176885157888 x - 2386516803584)
/(47041365435 x5 + 381702613848 x4 + 1360678489152 x3
+ 2856700692480 x2 + 3370143559680 x + 2386516803584)]

```

### powerdisp

Variable

Default value: `false`

When `powerdisp` is `true`, a sum is displayed with its terms in order of increasing power. Thus a polynomial is displayed as a truncated power series, with the constant term first and the highest power last.

By default, terms of a sum are displayed in order of decreasing power.

### powerseries (expr, x, a)

Function

Returns the general form of the power series expansion for `expr` in the variable `x` about the point `a` (which may be `inf` for infinity).

If `powerseries` is unable to expand `expr`, `taylor` may give the first several terms of the series.

When `verbose` is `true`, `powerseries` prints progress messages.

```

(%i1) verbose: true$
(%i2) powerseries (log(sin(x)/x), x, 0);
can't expand

```



```

                                log(sin(x))
so we'll try again after applying the rule:
                                d
                                / -- (sin(x))
                                [ dx
log(sin(x)) = i ----- dx
                                ]   sin(x)
                                /
in the first simplification we have returned:
                                /
                                [
                                i cot(x) dx - log(x)
                                ]
                                /
inf
====
\      i1  2 i1      2 i1
 (- 1)  2      bern(2 i1) x
 > -----
 /      i1 (2 i1)!
====
i1 = 1
(%o2) -----
                                2

```

**psexpand**

Variable

Default value: `false`

When `psexpand` is `true`, an extended rational function expression is displayed fully expanded. The switch `ratexpand` has the same effect.

When `psexpand` is `false`, a multivariate expression is displayed just as in the rational function package.

When `psexpand` is `multi`, then terms with the same total degree in the variables are grouped together.

**revert** (*expr*, *x*)

Function

**revert2** (*expr*, *x*, *n*)

Function

These functions return the reversion of *expr*, a Taylor series about zero in the variable *x*. `revert` returns a polynomial of degree equal to the highest power in *expr*. `revert2` returns a polynomial of degree *n*, which may be greater than, equal to, or less than the degree of *expr*.

`load ("revert")` loads these functions.

Examples:

```

(%i1) load ("revert")$
(%i2) t: taylor (exp(x) - 1, x, 0, 6);
                                2   3   4   5   6
                                x   x   x   x   x
(%o2)/T/      x + -- + -- + -- + --- + --- + . . .
                                2   6   24  120  720

```

```
(%i3) revert (t, x);
          6      5      4      3      2
      10 x  - 12 x  + 15 x  - 20 x  + 30 x  - 60 x
(%o3)/R/ -----
                      60

(%i4) ratexpand (%);
          6      5      4      3      2
          x      x      x      x      x
(%o4)      - -- + -- - -- + -- - -- + x
          6      5      4      3      2

(%i5) taylor (log(x+1), x, 0, 6);
          2      3      4      5      6
          x      x      x      x      x
(%o5)/T/      x - -- + -- - -- + -- - -- + . . .
          2      3      4      5      6

(%i6) ratsimp (revert (t, x) - taylor (log(x+1), x, 0, 6));
(%o6)
          0

(%i7) revert2 (t, x, 4);
          4      3      2
          x      x      x
(%o7)      - -- + -- - -- + x
          4      3      2
```

**taylor** (*expr*, *x*, *a*, *n*) Function  
**taylor** (*expr*, [*x*<sub>1</sub>, *x*<sub>2</sub>, ...], *a*, *n*) Function  
**taylor** (*expr*, [*x*, *a*, *n*, 'asympt]) Function  
**taylor** (*expr*, [*x*<sub>1</sub>, *x*<sub>2</sub>, ...], [*a*<sub>1</sub>, *a*<sub>2</sub>, ...], [*n*<sub>1</sub>, *n*<sub>2</sub>, ...]) Function

**taylor** (*expr*, *x*, *a*, *n*) expands the expression *expr* in a truncated Taylor or Laurent series in the variable *x* around the point *a*, containing terms through  $(x - a)^n$ .

If *expr* is of the form  $f(x)/g(x)$  and  $g(x)$  has no terms up to degree *n* then **taylor** attempts to expand  $g(x)$  up to degree  $2n$ . If there are still no nonzero terms, **taylor** doubles the degree of the expansion of  $g(x)$  so long as the degree of the expansion is less than or equal to  $n 2^{\text{taylordepth}}$ .

**taylor** (*expr*, [*x*<sub>1</sub>, *x*<sub>2</sub>, ...], *a*, *n*) returns a truncated power series of degree *n* in all variables *x*<sub>1</sub>, *x*<sub>2</sub>, ... about the point (*a*, *a*, ...).

**taylor** (*expr*, [*x*<sub>1</sub>, *a*<sub>1</sub>, *n*<sub>1</sub>], [*x*<sub>2</sub>, *a*<sub>2</sub>, *n*<sub>2</sub>], ...) returns a truncated power series in the variables *x*<sub>1</sub>, *x*<sub>2</sub>, ... about the point (*a*<sub>1</sub>, *a*<sub>2</sub>, ...), truncated at *n*<sub>1</sub>, *n*<sub>2</sub>, ...

**taylor** (*expr*, [*x*<sub>1</sub>, *x*<sub>2</sub>, ...], [*a*<sub>1</sub>, *a*<sub>2</sub>, ...], [*n*<sub>1</sub>, *n*<sub>2</sub>, ...]) returns a truncated power series in the variables *x*<sub>1</sub>, *x*<sub>2</sub>, ... about the point (*a*<sub>1</sub>, *a*<sub>2</sub>, ...), truncated at *n*<sub>1</sub>, *n*<sub>2</sub>, ...

**taylor** (*expr*, [*x*, *a*, *n*, 'asympt]) returns an expansion of *expr* in negative powers of  $x - a$ . The highest order term is  $(x - a)^{-n}$ .

When **maxtayorder** is **true**, then during algebraic manipulation of (truncated) Taylor series, **taylor** tries to retain as many terms as are known to be correct.

When **psexpand** is **true**, an extended rational function expression is displayed fully expanded. The switch **ratexpand** has the same effect. When **psexpand** is **false**, a

multivariate expression is displayed just as in the rational function package. When `psexpand` is `multi`, then terms with the same total degree in the variables are grouped together.

See also the `taylor_logexpand` switch for controlling expansion.

Examples:

```
(%i1) taylor (sqrt (sin(x) + a*x + 1), x, 0, 3);
```

```
(%o1)/T/ 1 + 
$$\frac{(a + 1) x^2}{2} - \frac{(a^2 + 2 a + 1) x^4}{8} + \frac{(3 a^3 + 9 a^2 + 9 a - 1) x^6}{48} + \dots$$

```

```
(%i2) %^2;
```

```
(%o2)/T/ 
$$1 + (a + 1) x^3 - \frac{x^6}{6} + \dots$$

```

```
(%i3) taylor (sqrt (x + 1), x, 0, 5);
```

```
(%o3)/T/ 
$$1 + \frac{x}{2} - \frac{x^2}{8} + \frac{x^3}{16} - \frac{5 x^4}{128} + \frac{7 x^5}{256} + \dots$$

```

```
(%i4) %^2;
```

```
(%o4)/T/ 
$$1 + x + \dots$$

```

```
(%i5) product ((1 + x^i)^2.5, i, 1, inf)/(1 + x^2);
```

```
(%o5) 
$$\frac{\prod_{i=1}^{\infty} (1 + x^i)^{2.5}}{1 + x^2}$$

```

```
(%o5)
```

```
(%i6) ev (taylor(%, x, 0, 3), keepfloat);
```

```
(%o6)/T/ 
$$1 + 2.5 x + 3.375 x^2 + 6.5625 x^3 + \dots$$

```

```
(%i7) taylor (1/log (x + 1), x, 0, 3);
```

```
(%o7)/T/ 
$$-\frac{1}{x} + \frac{1}{2} - \frac{x}{12} + \frac{x^2}{24} - \frac{19 x^3}{720} + \dots$$

```

```
(%i8) taylor (cos(x) - sec(x), x, 0, 5);
```

```
(%o8) 
$$\frac{x^4}{2} - x^2$$

```

```
(%o8)/T/          - x - -- + . . .
                    6
(%i9) taylor ((cos(x) - sec(x))^3, x, 0, 5);
(%o9)/T/          0 + . . .
(%i10) taylor (1/(cos(x) - sec(x))^3, x, 0, 5);
                    2          4
(%o10)/T/ - -- + ---- + ----- - ---- - ---- - ----
            6      4      2      15120  604800  7983360
            x      2 x      120 x
+ . . .

(%i11) taylor (sqrt (1 - k^2*sin(x)^2), x, 0, 6);
            2 2      4      2 4
            k x      (3 k - 4 k ) x
(%o11)/T/ 1 - ---- - ----
              2          24
                    6      4      2 6
                    (45 k - 60 k + 16 k ) x
                    - ---- + . . .
                      720

(%i12) taylor ((x + 1)^n, x, 0, 4);
            2      2      3      2      3
            (n - n) x      (n - 3 n + 2 n) x
(%o12)/T/ 1 + n x + ---- + ----
                  2          6
                    4      3      2      4
                    (n - 6 n + 11 n - 6 n) x
                    + ---- + . . .
                      24

(%i13) taylor (sin (y + x), x, 0, 3, y, 0, 3);
            3          2
            y          y
(%o13)/T/ y - -- + . . . + (1 - -- + . . .) x
            6          2
                    3          2
                    y y      2      1 y      3
                    + (- - + -- + . . .) x + (- - + -- + . . .) x + . . .
                    2 12      6 12

(%i14) taylor (sin (y + x), [x, y], 0, 3);
            3      2      2      3
            x + 3 y x + 3 y x + y
(%o14)/T/ y + x - ---- + . . .
                    6

(%i15) taylor (1/sin (y + x), x, 0, 3, y, 0, 3);
            1 y      1 1      1      2
            1 y      1 1      1      2
```

```
(%o15)/T/ - + - + . . . + (- -- + - + . . .) x + (-- + . . .) x
          y 6          2 6          3
                                     1          3
                                     + (- -- + . . .) x + . . .
                                     4
(%i16) taylor (1/sin (y + x), [x, y], 0, 3);
          y
          3          2          3
(%o16)/T/ ----- + ----- + ----- + . . .
          x + y      6          360
```

**taylordepth**

Variable

Default value: 3

If there are still no nonzero terms, `taylor` doubles the degree of the expansion of  $g(x)$  so long as the degree of the expansion is less than or equal to  $n 2^{\text{taylordepth}}$ .

**taylorinfo** (*expr*)

Function

Returns information about the Taylor series *expr*. The return value is a list of lists. Each list comprises the name of a variable, the point of expansion, and the degree of the expansion.

`taylorinfo` returns `false` if *expr* is not a Taylor series.

Example:

```
(%i1) taylor ((1 - y^2)/(1 - x), x, 0, 3, [y, a, inf]);
(%o1)/T/ - (y - a) - 2 a (y - a) + (1 - a)
          2          2
          + (1 - a - 2 a (y - a) - (y - a) ) x
          2          2 2
          + (1 - a - 2 a (y - a) - (y - a) ) x
          2          2 3
          + (1 - a - 2 a (y - a) - (y - a) ) x + . . .
(%i2) taylorinfo(%);
(%o2) [[y, a, inf], [x, 0, 3]]
```

**taylorp** (*expr*)

Function

Returns `true` if *expr* is a Taylor series, and `false` otherwise.

**taylor\_logexpand**

Variable

Default value: `true`

`taylor_logexpand` controls expansions of logarithms in `taylor` series.

When `taylor_logexpand` is `true`, all logarithms are expanded fully so that zero-recognition problems involving logarithmic identities do not disturb the expansion process. However, this scheme is not always mathematically correct since it ignores branch information.

When `taylor_logexpand` is set to `false`, then the only expansion of logarithms that occur is that necessary to obtain a formal power series.

**taylor\_order\_coefficients** Variable

Default value: `true`

`taylor_order_coefficients` controls the ordering of coefficients in a Taylor series.

When `taylor_order_coefficients` is `true`, coefficients of Taylor series are ordered canonically.

**taylor\_simplifier** (*expr*) Function

Simplifies coefficients of the power series *expr*. `taylor` calls this function.

**taylor\_truncate\_polynomials** Variable

Default value: `true`

When `taylor_truncate_polynomials` is `true`, polynomials are truncated based upon the input truncation levels.

Otherwise, polynomials input to `taylor` are considered to have infinite precision.

**taylorat** (*expr*) Function

Converts *expr* from `taylor` form to canonical rational expression (CRE) form. The effect is the same as `rat` (`ratdisrep` (*expr*)), but faster.

**trunc** (*expr*) Function

Annotates the internal representation of the general expression *expr* so that it is displayed as if its sums were truncated Taylor series. *expr* is not otherwise modified.

Example:

```
(%i1) expr: x^2 + x + 1;
(%o1)          2
              x  + x + 1
(%i2) trunc (expr);
(%o2)          2
              1 + x + x  + . . .
(%i3) is (expr = trunc (expr));
(%o3)          true
```

**unsum** (*f*, *n*) Function

Returns the first backward difference  $f(n) - f(n - 1)$ . Thus `unsum` in a sense is the inverse of `sum`.

See also `nusum`.

Examples:

```
(%i1) g(p) := p*4^n/binomial(2*n,n);
(%o1)          g(p) := 
$$\frac{p^4}{\text{binomial}(2n, n)}$$

(%i2) g(n^4);
(%o2)          
$$\frac{n^4}{\text{binomial}(2n, n)}$$

(%i3) nusum (%, n, 0, n);
(%o3) 
$$\frac{2(n+1)(63n^4 + 112n^3 + 18n^2 - 22n + 3)}{693 \text{binomial}(2n, n)}$$

(%i4) unsum (%, n);
(%o4)          
$$\frac{n^4}{\text{binomial}(2n, n)}$$

```

**verbose**

Variable

Default value: false

When `verbose` is true, `powerseries` prints progress messages.

## 33 Number Theory

### 33.1 Definitions for Number Theory

**bern** (*n*) Function

Returns the *n*'th Bernoulli number for integer *n*. Bernoulli numbers equal to zero are suppressed if `zerobern` is `false`.

See also `burn`.

```
(%i1) zerobern: true$
(%i2) map (bern, [0, 1, 2, 3, 4, 5, 6, 7, 8]);
          1 1      1      1      1
(%o2) [1, - -, -, 0, - --, 0, --, 0, - --]
          2 6      30      42      30
(%i3) zerobern: false$
(%i4) map (bern, [0, 1, 2, 3, 4, 5, 6, 7, 8]);
          1 1      1 5      691 7      3617 43867
(%o4) [1, - -, -, - --, --, - ----, -, - ----, ----]
          2 6      30 66      2730 6      510      798
```

**bernpoly** (*x*, *n*) Function

Returns the *n*'th Bernoulli polynomial in the variable *x*.

**bfzeta** (*s*, *n*) Function

Returns the Riemann zeta function for the argument *s*. The return value is a big float (bfloat); *n* is the number of digits in the return value.

`load ("bffac")` loads this function.

**bfhzeta** (*s*, *h*, *n*) Function

Returns the Hurwitz zeta function for the arguments *s* and *h*. The return value is a big float (bfloat); *n* is the number of digits in the return value.

The Hurwitz zeta function is defined as

$$\sum ((k+h)^{-s}, k, 0, \text{inf})$$

`load ("bffac")` loads this function.

**binomial** (*x*, *y*) Function

The binomial coefficient  $(x + y)! / (x! y!)$ . If *x* and *y* are integers, then the numerical value of the binomial coefficient is computed. If *y*, or *x* - *y*, is an integer, the binomial coefficient is expressed as a polynomial.

**burn** (*n*) Function

Returns the *n*'th Bernoulli number for integer *n*. `burn` may be more efficient than `bern` for large, isolated *n* (perhaps *n* greater than 105 or so), as `bern` computes all the Bernoulli numbers up to index *n* before returning.

`burn` exploits the observation that (rational) Bernoulli numbers can be approximated by (transcendental) zetas with tolerable efficiency.

`load ("bffac")` loads this function.



**cf** (*expr*)

Function

Converts *expr* into a continued fraction. *expr* is an expression comprising continued fractions and square roots of integers. Operands in the expression may be combined with arithmetic operators. Aside from continued fractions and square roots, factors in the expression must be integer or rational numbers. Maxima does not know about operations on continued fractions outside of **cf**.

**cf** evaluates its arguments after binding **listarith** to **false**. **cf** returns a continued fraction, represented as a list.

A continued fraction  $a + 1/(b + 1/(c + \dots))$  is represented by the list  $[a, b, c, \dots]$ . The list elements  $a, b, c, \dots$  must evaluate to integers. *expr* may also contain **sqrt** (*n*) where *n* is an integer. In this case **cf** will give as many terms of the continued fraction as the value of the variable **cflength** times the period.

A continued fraction can be evaluated to a number by evaluating the arithmetic representation returned by **cfdisrep**. See also **cfexpand** for another way to evaluate a continued fraction.

See also **cfdisrep**, **cfexpand**, and **cflength**.

Examples:

- expr* is an expression comprising continued fractions and square roots of integers.

```
(%i1) cf ([5, 3, 1]*[11, 9, 7] + [3, 7]/[4, 3, 2]);
(%o1) [59, 17, 2, 1, 1, 1, 27]
(%i2) cf ((3/17)*[1, -2, 5]/sqrt(11) + (8/13));
(%o2) [0, 1, 1, 1, 3, 2, 1, 4, 1, 9, 1, 9, 2]
```

- cflength** controls how many periods of the continued fraction are computed for algebraic, irrational numbers.

```
(%i1) cflength: 1$
(%i2) cf ((1 + sqrt(5))/2);
(%o2) [1, 1, 1, 1, 2]
(%i3) cflength: 2$
(%i4) cf ((1 + sqrt(5))/2);
(%o4) [1, 1, 1, 1, 1, 1, 1, 2]
(%i5) cflength: 3$
(%i6) cf ((1 + sqrt(5))/2);
(%o6) [1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 2]
```

- A continued fraction can be evaluated by evaluating the arithmetic representation returned by **cfdisrep**.

```
(%i1) cflength: 3$
(%i2) cfdisrep (cf (sqrt (3)))$
(%i3) ev (% , numer);
(%o3) 1.731707317073171
```

- Maxima does not know about operations on continued fractions outside of **cf**.

```
(%i1) cf ([1,1,1,1,1,2] * 3);
(%o1) [4, 1, 5, 2]
(%i2) cf ([1,1,1,1,1,2]) * 3;
(%o2) [3, 3, 3, 3, 3, 6]
```

**cfdisrep** (*list*) Function

Constructs and returns an ordinary arithmetic expression of the form  $a + 1/(b + 1/(c + \dots))$  from the list representation of a continued fraction  $[a, b, c, \dots]$ .

```
(%i1) cf ([1, 2, -3] + [1, -2, 1]);
(%o1) [1, 1, 1, 2]
(%i2) cfdisrep (%);
(%o2) 1 + -----
              1
              1 + -----
                    1
                    1 + -
                          2
```

**cfexpand** (*x*) Function

Returns a matrix of the numerators and denominators of the last (column 1) and next-to-last (column 2) convergents of the continued fraction *x*.

```
(%i1) cf (rat (ev (%pi, numer)));
'rat' replaced 3.141592653589793 by 103993//33102 = 3.141592653011902
(%o1) [3, 7, 15, 1, 292]
(%i2) cfexpand (%);
(%o2) [ 103993  355 ]
      [          ]
      [ 33102   113 ]
(%i3) %[1,1]/[%2,1], numer;
(%o3) 3.141592653011902
```

**cflength** Variable

Default value: 1

**cflength** controls the number of terms of the continued fraction the function **cf** will give, as the value **cflength** times the period. Thus the default is to give one period.

```
(%i1) cflength: 1$
(%i2) cf ((1 + sqrt(5))/2);
(%o2) [1, 1, 1, 1, 2]
(%i3) cflength: 2$
(%i4) cf ((1 + sqrt(5))/2);
(%o4) [1, 1, 1, 1, 1, 1, 1, 2]
(%i5) cflength: 3$
(%i6) cf ((1 + sqrt(5))/2);
(%o6) [1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 2]
```

**divsum** (*n*, *k*) Function

**divsum** (*n*) Function

**divsum** (*n*, *k*) returns the sum of the divisors of *n* raised to the *k*'th power.

**divsum** (*n*) returns the sum of the divisors of *n*.

```
(%i1) divsum (12);
(%o1) 28
(%i2) 1 + 2 + 3 + 4 + 6 + 12;
(%o2) 28
(%i3) divsum (12, 2);
(%o3) 210
(%i4) 1^2 + 2^2 + 3^2 + 4^2 + 6^2 + 12^2;
(%o4) 210
```

**euler** (*n*) Function

Returns the *n*'th Euler number for nonnegative integer *n*.

For the Euler-Mascheroni constant, see `%gamma`.

```
(%i1) map (euler, [0, 1, 2, 3, 4, 5, 6, 7, 8, 9, 10]);
(%o1) [1, 0, -1, 0, 5, 0, -61, 0, 1385, 0, -50521]
```

**%gamma** Variable

The Euler-Mascheroni constant, 0.5772156649015329 ....

**factorial** (*x*) Function

Represents the factorial function. Maxima treats `factorial (x)` the same as `x!`. See `!`.

**fib** (*n*) Function

Returns the *n*'th Fibonacci number. `fib(0)` equal to 0 and `fib(1)` equal to 1, and `fib (-n)` equal to  $(-1)^{(n+1)} * \text{fib}(n)$ .

After calling `fib`, `prevfib` is equal to `fib (x - 1)`, the Fibonacci number preceding the last one computed.

```
(%i1) map (fib, [0, 1, 2, 3, 4, 5, 6, 7, 8, 9, 10]);
(%o1) [0, 1, 1, 2, 3, 5, 8, 13, 21, 34, 55]
```

**fibtophi** (*expr*) Function

Expresses Fibonacci numbers in terms of the constant `%phi`, which is  $(1 + \sqrt{5})/2$ , approximately 1.61803399.

By default, Maxima does not know about `%phi`. After executing `tellrat (%phi^2 - %phi - 1)` and `algebraic: true`, `ratsimp` can simplify some expressions containing `%phi`.

```
(%i1) fibtophi (fib (n));
(%o1) 
$$\frac{\%phi^n - (1 - \%phi)^n}{2 \%phi - 1}$$

(%i2) fib (n-1) + fib (n) - fib (n+1);
(%o2) - fib(n + 1) + fib(n) + fib(n - 1)
(%i3) ratsimp (fibtophi (%));
(%o3) 0
```

**inrt** (*x*, *n*) Function

Returns the integer *n*'th root of the absolute value of *x*.

```
(%i1) 1: [1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12]$
(%i2) map (lambda ([a], inrt (10^a, 3)), 1);
(%o2) [2, 4, 10, 21, 46, 100, 215, 464, 1000, 2154, 4641, 10000]
```

**jacobi** (*p*, *q*) Function

Returns the Jacobi symbol of *p* and *q*.

```
(%i1) 1: [1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12]$
(%i2) map (lambda ([a], jacobi (a, 9)), 1);
(%o2) [1, 1, 0, 1, 1, 0, 1, 1, 0, 1, 1, 0]
```

**lcm** (*expr\_1*, ..., *expr\_n*) Function

Returns the least common multiple of its arguments. The arguments may be general expressions as well as integers.

load ("functs") loads this function.

**minfactorial** (*expr*) Function

Examines *expr* for occurrences of two factorials which differ by an integer. **minfactorial** then turns one into a polynomial times the other.

```
(%i1) n!/(n+2)!;
(%o1)
          n!
-----
      (n + 2)!
(%i2) minfactorial (%);
(%o2)
          1
-----
      (n + 1) (n + 2)
```

**partfrac** (*expr*, *var*) Function

Expands the expression *expr* in partial fractions with respect to the main variable *var*. **partfrac** does a complete partial fraction decomposition. The algorithm employed is based on the fact that the denominators of the partial fraction expansion (the factors of the original denominator) are relatively prime. The numerators can be written as linear combinations of denominators, and the expansion falls out.

```
(%i1) 1/(1+x)^2 - 2/(1+x) + 2/(2+x);
(%o1)
          2      2      1
----- - ----- + -----
      x + 2   x + 1   (x + 1)^2
(%i2) ratsimp (%);
(%o2)
          x
-----
          3      2
         x  + 4 x  + 5 x + 2
(%i3) partfrac (% , x);
          2      2      1
```

```
(%o3)          ----- - ----- + -----
              x + 2   x + 1   (x + 1)2
```

**primep** (*n*) Function

Returns true if *n* is a prime, false if not.

**qunit** (*n*) Function

Returns the principal unit of the real quadratic number field `sqrt` (*n*) where *n* is an integer, i.e., the element whose norm is unity. This amounts to solving Pell's equation  $a^2 - n b^2 = 1$ .

```
(%i1) qunit (17);
(%o1)          sqrt(17) + 4
(%i2) expand (% * (sqrt(17) - 4));
(%o2)          1
```

**totient** (*n*) Function

Returns the number of integers less than or equal to *n* which are relatively prime to *n*.

**zerobern** Variable

Default value: true

When `zerobern` is false, `bern` excludes the Bernoulli numbers which are equal to zero. See `bern`.

**zeta** (*n*) Function

Returns the Riemann zeta function if *x* is a negative integer, 0, 1, or a positive even number, and returns a noun form `zeta` (*n*) for all other arguments, including rational noninteger, floating point, and complex arguments.

See also `bfzeta` and `zeta%pi`.

```
(%i1) map (zeta, [-4, -3, -2, -1, 0, 1, 2, 3, 4, 5]);
(%o1) [0, ---, 0, - ---, - -, inf, ---, zeta(3), ---, zeta(5)]
        120      12   2     6      90
```

**zeta%pi** Variable

Default value: true

When `zeta%pi` is true, `zeta` returns an expression proportional to  $\%pi^n$  for even integer *n*. Otherwise, `zeta` returns a noun form `zeta` (*n*) for even integer *n*.

```
(%i1) zeta%pi: true$
(%i2) zeta (4);
(%o2)          4
              %pi
              ----
              90
```

```
(%i3) zeta%pi: false$
(%i4) zeta (4);
(%o4) zeta(4)
```



## 34 Symmetries

### 34.1 Definitions for Symmetries

**comp2pui** ( $n, l$ ) Function

re'alise le passage des fonctions syme'triques comple'tes, donne'es dans la liste  $l$ , aux fonctions syme'triques e'le'mentaires de 0 a'  $n$ . Si la liste  $l$  contient moins de  $n+1$  e'le'ments les valeurs formelles viennent la completer. Le premier e'le'ment de la liste  $l$  donne le cardinal de l'alphabet si il existe, sinon on le met e'gal a  $n$ .

```
(%i1) comp2pui (3, [4, g]);
      2
      2
(%o1) [4, g, 2 h2 - g , 3 h3 - g h2 + g (g - 2 h2)]
```

**cont2part** ( $pc, lvar$ ) Function

rend le polyno'me partitionne' associe' a' la forme contracte'e  $pc$  dont les variables sont dans  $lvar$ .

```
(%i1) pc: 2*a^3*b*x^4*y + x^5;
      3 4 5
      2 a b x y + x
(%i2) cont2part (pc, [x, y]);
      3
(%o2) [[1, 5, 0], [2 a b, 4, 1]]
```

Autres fonctions de changements de repre'sentations :

contract, explose, part2cont, partpol, tcontract, tpartpol.

**contract** ( $psym, lvar$ ) Function

rend une forme contracte'e (i.e. un mono'me par orbite sous l'action du groupe syme'trique) du polyno'me  $psym$  en les variables contenues dans la liste  $lvar$ . La fonction **explose** re'alise l'ope'ration inverse. La fonction **tcontract** teste en plus la syme'trie du polyno'me.

```
(%i1) psym: explose (2*a^3*b*x^4*y, [x, y, z]);
      3 4 3 4 3 4 3 4
      2 a b y z + 2 a b x z + 2 a b y z + 2 a b x z
      3 4 3 4
      + 2 a b x y + 2 a b x y
(%i2) contract (psym, [x, y, z]);
      3 4
(%o2) 2 a b x y
```

Autres fonctions de changements de repre'sentations :

cont2part, explose, part2cont, partpol, tcontract, tpartpol.

**direct** ( $[p_1, \dots, p_n], y, f, [lvar_1, \dots, lvar_n]$ ) Function

calcul l'image directe (voir M. GIUSTI, D. LAZARD et A. VALIBOUZE, ISSAC 1988, Rome) associe'e a' la fonction  $f$ , en les listes de variables  $lvar_1, \dots, lvar_n$ , et



aux polynômes  $p_1, \dots, p_n$  d'une variable  $y$ . l'arité de la fonction  $f$  est importante pour le calcul. Ainsi, si l'expression de  $f$  ne depend pas d'une variable, non seulement il est inutile de donner cette variable mais cela diminue considerablement les calculs si on ne le fait pas.

```
(%i1) direct ([z^2 - e1*z + e2, z^2 - f1*z + f2],
             z, b*v + a*u, [[u, v], [a, b]]);
```

```
(%o1) y^2 - e1 f1 y
```

$$+ \frac{-4 e_2 f_2 - (e_1^2 - 2 e_2) (f_1^2 - 2 f_2) + e_1^2 f_1^2}{2}$$

```
(%i2) ratsimp (%);
```

```
(%o2) y^2 - e1 f1 y + (e1^2 - 4 e2) f2 + e2 f1^2
```

```
(%i3) ratsimp (direct ([z^3-e1*z^2+e2*z-e3,z^2 - f1*z + f2],
                       z, b*v + a*u, [[u, v], [a, b]]));
```

```
(%o3) y^6 - 2 e1 f1 y^5 + ((2 e1^2 - 6 e2) f2 + (2 e2 + e1^2) f1^2) y^4
```

$$+ ((9 e_3 + 5 e_1 e_2 - 2 e_1^3) f_1 f_2 + (-2 e_3 - 2 e_1 e_2) f_1^3) y^3$$

$$+ ((9 e_2^2 - 6 e_1^2 e_2 + e_1^4) f_2^2$$

$$+ (-9 e_1 e_3 - 6 e_2^2 + 3 e_1^2 e_2) f_1^2 f_2 + (2 e_1 e_3 + e_2^2) f_1^4$$

$$y^2 + (((9 e_1^2 - 27 e_2) e_3 + 3 e_1 e_2^2 - e_1^3 e_2) f_1^2 f_2^2$$

$$+ ((15 e_2^2 - 2 e_1^2) e_3 - e_1^2 e_2^3) f_1^2 f_2^3 - 2 e_2 e_3 f_1^5) y$$

$$+ (-27 e_3^2 + (18 e_1 e_2 - 4 e_1^3) e_3 - 4 e_2^3 + e_1^2 e_2^2) f_2^3$$

$$+ (27 e_3^2 + (e_1^3 - 9 e_1 e_2) e_3 + e_2^3) f_1^3 f_2^2$$

$$+ (e_1 e_2 e_3 - 9 e_3^2) f_1^4 f_2 + e_3^6 f_1^6$$

Recherche du polynôme dont les racines sont les somme  $a+u$  ou  $a$  est racine de  $z^2 - e1*z + e2$  et  $u$  est racine de  $z^2 - f1*z + f2$

```
(%i1) ratsimp (direct ([z^2 - e1*z + e2, z^2 - f1*z + f2],
                       z, a + u, [[u], [a]]));
```

```

(%o1) y4 + (- 2 f1 - 2 e1) y3 + (2 f2 + f12 + 3 e1 f1 + 2 e2
+ e12) y2 + ((- 2 f1 - 2 e1) f2 - e1 f12 + (- 2 e2 - e12) f1
- 2 e1 e2) y + f22 + (e1 f1 - 2 e2 + e12) f2 + e2 f12 + e1 e2 f1
+ e22

```

`direct` peut prendre deux drapeaux possibles : `elementaires` et `puissances` (valeur par de'faut) qui permettent de de'composer les polyno^mes syme'triques apparaissant dans ce calcul par les fonctions syme'triques e'le'mentaires ou les fonctions puissances respectivement.

Fonctions de `sym` utilis'ees dans cette fonction :

`multi_orbit` (donc `orbit`), `pui_direct`, `multi_elem` (donc `elem`), `multi_pui` (donc `pui`), `pui2ele`, `ele2pui` (si le drapeau `direct` est a' `puissances`).

### `ele2comp` ( $m, l$ )

Function

passer des fonctions syme'triques e'le'mentaires aux fonctions comple'tes. Similaire a' `comp2ele` et `comp2pui`.

Autres fonctions de changements de bases :

`comp2ele`, `comp2pui`, `ele2pui`, `elem`, `mon2schur`, `multi_elem`, `multi_pui`, `pui`, `pui2comp`, `pui2ele`, `puireduc`, `schur2comp`.

### `ele2polynome` ( $l, z$ )

Function

donne le polyno^me en  $z$  dont les fonctions syme'triques e'le'mentaires des racines sont dans la liste  $l$ .  $l = [n, e_1, \dots, e_n]$  ou'  $n$  est le degre' du polyno^me et  $e_i$  la  $i$ -ie'me fonction syme'trique e'le'mentaire.

```

(%i1) ele2polynome ([2, e1, e2], z);
(%o1) z2 - e1 z + e2
(%i2) polynome2ele (x^7 - 14*x^5 + 56*x^3 - 56*x + 22, x);
(%o2) [7, 0, -14, 0, 56, 0, -56, -22]
(%i3) ele2polynome ([7, 0, -14, 0, 56, 0, -56, -22], x);
(%o3) x7 - 14 x5 + 56 x3 - 56 x + 22

```

La re'ciproque: `polynome2ele` ( $P, z$ )

Autres fonctions a' voir :

`polynome2ele`, `pui2polynome`.

### `ele2pui` ( $m, l$ )

Function

passer des fonctions syme'triques e'le'mentaires aux fonctions comple'tes. Similaire a' `comp2ele` et `comp2pui`.

Autres fonctions de changements de bases :

comp2ele, comp2pui, ele2comp, elem, mon2schur, multi\_elem, multi\_pui, pui, pui2comp, pui2ele, puireduc, schur2comp.

**elem** (*ele, sym, lvar*)

Function

decompose le polynôme symétrique *sym*, en les variables contenues de la liste *lvar*, par les fonctions symétriques élémentaires contenues dans la liste *ele*. Si le premier élément de *ele* est donné ce sera le cardinal de l'alphabet sinon on prendra le degré du polynôme *sym*. Si il manque des valeurs à la liste *ele* des valeurs formelles du type "ei" sont rajoutées. Le polynôme *sym* peut être donné sous 3 formes différentes : contractée (**elem** doit alors valoir 1 sa valeur par défaut), partitionnée (**elem** doit alors valoir 3) ou étendue (i.e. le polynôme en entier) (**elem** doit alors valoir 2). L'utilisation de la fonction **pui** se réalise sur le même modèle.

Sur un alphabet de cardinal 3 avec e1, la première fonction symétrique élémentaire, valant 7, le polynôme symétrique en 3 variables dont la forme contractée (ne dépendant ici que de deux de ses variables) est  $x^4 - 2x^2y$  se décompose ainsi en les fonctions symétriques élémentaires :

```
(%i1) elem ([3, 7], x^4 - 2*x*y, [x, y]);
(%o1) 7 (e3 - 7 e2 + 7 (49 - e2)) + 21 e3
                                     + (- 2 (49 - e2) - 2) e2
(%i2) ratsimp (%);
(%o2) 28 e3 + 2 e2^2 - 198 e2 + 2401
```

Autres fonctions de changements de bases :

comp2ele, comp2pui, ele2comp, ele2pui, mon2schur, multi\_elem, multi\_pui, pui, pui2comp, pui2ele, puireduc, schur2comp.

**explode** (*pc, lvar*)

Function

rend le polynôme symétrique associé à la forme contractée *pc*. La liste *lvar* contient les variables.

```
(%i1) explode (a*x + 1, [x, y, z]);
(%o1) a z + a y + a x + 1
```

Autres fonctions de changements de représentations :

contract, cont2part, part2cont, partpol, tcontract, tpartpol.

**kostka** (*part\_1, part\_2*)

Function

écrite par P. ESPERET, calcule le nombre de Kostka associé aux partitions *part\_1* et *part\_2*.

```
(%i1) kostka ([3, 3, 3], [2, 2, 2, 1, 1, 1]);
(%o1) 6
```

**lgtreillis** (*n, m*)

Function

rend la liste des partitions de poids *n* et de longueur *m*.

```
(%i1) lgtreillis (4, 2);
(%o1)          [[3, 1], [2, 2]]
```

Voir e'galement : `ltreillis`, `treillis` et `treinat`.

### **ltreillis** (*n*, *m*)

Function

rend la liste des partitions de poids *n* et de longueur inférieure ou égale à *m*.

```
(%i1) ltreillis (4, 2);
(%o1)          [[4, 0], [3, 1], [2, 2]]
```

Voir e'galement : `lgtreillis`, `treillis` et `treinat`.

### **mon2schur** (*l*)

Function

la liste *l* représente la fonction de Schur  $S_l$ : On a  $l = [i_1, i_2, \dots, i_q]$  avec  $i_1 \leq i_2 \leq \dots \leq i_q$ . La fonction de Schur est  $S_{[i_1, i_2, \dots, i_q]}$  est le mineur de la matrice infinie  $(h_{i-j})_{i \geq 1, j \geq 1}$  composé des *q* premières lignes et des colonnes  $i_1 + 1, i_2 + 2, \dots, i_q + q$ .

On écrit cette fonction de Schur en fonction des formes monomiales en utilisant les fonctions `treinat` et `kostka`. La forme rendue est un polynôme symétrique dans une de ses représentations contractées avec les variables  $x_1, x_2, \dots$

```
(%i1) mon2schur ([1, 1, 1]);
(%o1)          x1 x2 x3
(%i2) mon2schur ([3]);
(%o2)          x1 x2 x3 + x1^2 x2 + x1^3
(%i3) mon2schur ([1, 2]);
(%o3)          2 x1 x2 x3 + x1^2 x2
```

ce qui veut dire que pour 3 variables cela donne :

$$2 x_1 x_2 x_3 + x_1^2 x_2 + x_2^2 x_1 + x_1^2 x_3 + x_3^2 x_1 + x_2^2 x_3 + x_3^2 x_2$$

Autres fonctions de changements de bases :

`comp2ele`, `comp2pui`, `ele2comp`, `ele2pui`, `elem`, `multi_elem`, `multi_pui`, `pui`, `pui2comp`, `pui2ele`, `puireduc`, `schur2comp`.

### **multi\_elem** (*lelem*, *multi\_pc*, *Lvar*)

Function

decompose un polynôme multi-symétrique sous la forme multi-contractée *multi\_pc* en les groupes de variables contenue dans la liste de listes *Lvar* sur les groupes de fonctions symétriques élémentaires contenues dans *lelem*.

```
(%i1) multi_elem ([[2, e1, e2], [2, f1, f2]], a*x + a^2 + x^3, [[x, y], [a, b]])
(%o1)          - 2 f2 + f1 (f1 + e1) - 3 e1 e2 + e1^3
(%i2) ratsimp (%);
(%o2)          - 2 f2 + f1^2 + e1 f1 - 3 e1 e2 + e1^3
```

Autres fonctions de changements de bases :

`comp2ele`, `comp2pui`, `ele2comp`, `ele2pui`, `elem`, `mon2schur`, `multi_pui`, `pui`, `pui2comp`, `pui2ele`, `puireduc`, `schur2comp`.

**multi\_orbit** ( $P$ , [ $lvar_1$ ,  $lvar_2$ , ...,  $lvar_p$ ]) Function

$P$  est un polynôme en l'ensemble des variables contenues dans les listes  $lvar_1$ ,  $lvar_2$ , ...,  $lvar_p$ . Cette fonction ramène l'orbite du polynôme  $P$  sous l'action du produit des groupes symétriques des ensembles de variables représentés par ces  $p$  listes.

```
(%i1) multi_orbit (a*x + b*y, [[x, y], [a, b]]);
(%o1)          [b y + a x, a y + b x]
(%i2) multi_orbit (x + y + 2*a, [[x, y], [a, b, c]]);
(%o2)          [y + x + 2 c, y + x + 2 b, y + x + 2 a]
```

Voir également : `orbit` pour l'action d'un seul groupe symétrique.

**multi\_pui** Function

est à la fonction `pui` ce que la fonction `multi_elem` est à la fonction `elem`.

```
(%i1) multi_pui ([[2, p1, p2], [2, t1, t2]], a*x + a^2 + x^3, [[x, y], [a, b]])
3
(%o1)          t2 + p1 t1 + ----- - ----
                2          2
```

**multinomial** ( $r$ ,  $part$ ) Function

où  $r$  est le poids de la partition  $part$ . Cette fonction ramène le coefficient multinomial associé : si les parts de la partitions  $part$  sont  $i_1, i_2, \dots, i_k$ , le résultat de `multinomial` est  $r!/(i_1! i_2! \dots i_k!)$ .

**multsym** ( $ppart_1$ ,  $ppart_2$ ,  $n$ ) Function

réalise le produit de deux polynômes symétriques de  $n$  variables en ne travaillant que modulo l'action du groupe symétrique d'ordre  $n$ . Les polynômes sont dans leur représentation partitionnée.

Soient les 2 polynômes symétriques en  $x, y$ :  $3*(x + y) + 2*x*y$  et  $5*(x^2 + y^2)$  dont les formes partitionnées sont respectivement  $[[3, 1], [2, 1, 1]]$  et  $[[5, 2]]$ , alors leur produit sera donné par :

```
(%i1) multsym ([[3, 1], [2, 1, 1]], [[5, 2]], 2);
(%o1)          [[10, 3, 1], [15, 3, 0], [15, 2, 1]]
```

soit  $10*(x^3*y + y^3*x) + 15*(x^2*y + y^2*x) + 15*(x^3 + y^3)$ .

Fonctions de changements de représentations d'un polynôme symétrique :

`contract`, `cont2part`, `explose`, `part2cont`, `partpol`, `tcontract`, `tpartpol`.

**orbit** ( $P$ ,  $lvar$ ) Function

calcul l'orbite du polynôme  $P$  en les variables de la liste  $lvar$  sous l'action du groupe symétrique de l'ensemble des variables contenues dans la liste  $lvar$ .

```
(%i1) orbit (a*x + b*y, [x, y]);
(%o1)          [a y + b x, b y + a x]
(%i2) orbit (2*x + x^2, [x, y]);
(%o2)          [y + 2 y, x + 2 x]
```

Voir également : `multi_orbit` pour l'action d'un produit de groupes symétriques sur un polynôme.

**part2cont** (*ppart*, *lvar*) Function  
 passe de la forme partitionne'e a' la forme contracte'e d'un polynome syme'trique.  
 La forme contracte'e est rendue avec les variables contenues dans *lvar*.

```
(%i1) part2cont ([[2*a^3*b, 4, 1]], [x, y]);
              3      4
(%o1)          2 a b x y
```

Autres fonctions de changements de repre'sentations :

`contract`, `cont2part`, `explode`, `partpol`, `tcontract`, `tpartpol`.

**partpol** (*psym*, *lvar*) Function  
*psym* est un polynome syme'trique en les variables de *lvar*. Cette fonction rame'ne  
 sa repre'sentation partitionne'e.

```
(%i1) partpol (-a*(x + y) + 3*x*y, [x, y]);
(%o1)          [[3, 1, 1], [- a, 1, 0]]
```

Autres fonctions de changements de repre'sentations :

`contract`, `cont2part`, `explode`, `part2cont`, `tcontract`, `tpartpol`.

**permut** (*l*) Function  
 rame'ne la liste des permutations de la liste *l*.

**polynome2ele** (*P*, *x*) Function  
 donne la liste  $l = [n, e_1, \dots, e_n]$  ou'  $n$  est le degre' du polynome  $P$  en la variable  
 $x$  et  $e_i$  la  $i$ -ieme fonction syme'trique e'le'mentaire des racines de  $P$ .

```
(%i1) polynome2ele (x^7 - 14*x^5 + 56*x^3 - 56*x + 22, x);
(%o1)          [7, 0, - 14, 0, 56, 0, - 56, - 22]
(%i2) ele2polynome ([7, 0, -14, 0, 56, 0, -56, -22], x);
              7      5      3
(%o2)          x - 14 x + 56 x - 56 x + 22
```

La re'ciproque : `ele2polynome` (*l*, *x*)

**prodrac** (*l*, *k*) Function  
*l* est une liste contenant les fonctions syme'triques e'le'mentaires sur un ensemble  $A$ .  
`prodrac` rend le polynome dont les racines sont les produits  $k$  a'  $k$  des e'le'ments de  
 $A$ .

**pui** (*l*, *sym*, *lvar*) Function  
 de'compose le polynome syme'trique *sym*, en les variables contenues de la liste *lvar*,  
 par les fonctions puissances contenues dans la liste *l*. Si le premier e'le'ment de *l* est  
 donne' ce sera le cardinal de l'alphabet sinon on prendra le degre' du polynome *sym*.  
 Si il manque des valeurs a' la liste *l*, des valeurs formelles du type "pi" sont rajoute'es.  
 Le polynome *sym* peut etre donne' sous 3 formes diffe'rentes : contracte'e (`pui` doit  
 alors valoir 1 sa valeur par de'faut), partitionne'e (`pui` doit alors valoir 3) ou e'tendue  
 (i.e. le polynome en entier) (`pui` doit alors valoir 2). La fonction `elem` s'utilise de  
 la me'me manie're.

```
(%i1) pui;
(%o1)
(%i2) pui ([3, a, b], u*x*y*z, [x, y, z]);
          1
          2
          a (a - b) u  (a b - p3) u
(%o2)  ----- - -----
          6          3
(%i3) ratsimp (%);
          3
          (2 p3 - 3 a b + a ) u
(%o3)  -----
          6
```

Autres fonctions de changements de bases :

comp2ele, comp2pui, ele2comp, ele2pui, elem, mon2schur, multi\_elem, multi\_pui, pui2comp, pui2ele, puireduc, schur2comp.

### **pui2comp** (*n*, *lpui*)

Function

rend la liste des *n* premières fonctions complètes (avec en tête le cardinal) en fonction des fonctions puissance données dans la liste *lpui*. Si la liste *lpui* est vide le cardinal est N sinon c'est son premier élément similaire à `comp2ele` et `comp2pui`.

```
(%i1) pui2comp (2, []);
          2
          p2 + p1
(%o1)  [2, p1, -----]
          2
(%i2) pui2comp (3, [2, a1]);
          2
          a1 (p2 + a1 )
          2  p3 + ----- + a1 p2
          2
(%o2)  [2, a1, -----, -----]
          2          3
(%i3) ratsimp (%);
          2          3
          p2 + a1  2 p3 + 3 a1 p2 + a1
(%o3)  [2, a1, -----, -----]
          2          6
```

Autres fonctions de changements de bases :

comp2ele, comp2pui, ele2comp, ele2pui, elem, mon2schur, multi\_elem, multi\_pui, pui, pui2ele, puireduc, schur2comp.

### **pui2ele** (*n*, *lpui*)

Function

réalise le passage des fonctions puissances aux fonctions symétriques élémentaires. Si le drapeau `pui2ele` est `girard`, on récupère la liste des fonctions symétriques élémentaires de 1 à *n*, et s'il est égal à `close`, la *n*-ième fonction symétrique élémentaire.

Autres fonctions de changements de bases :

comp2ele, comp2pui, ele2comp, ele2pui, elem, mon2schur, multi\_elem, multi\_pui, pui, pui2comp, puireduc, schur2comp.

**pui2polynome** (*x*, *lpui*)

Function

calcul le polynôme en *x* dont les fonctions puissances des racines sont données dans la liste *lpui*.

```
(%i1) pui;
(%o1) 1
(%i2) kill(labels);
(%o0) done
(%i1) polynome2ele (x^3 - 4*x^2 + 5*x - 1, x);
(%o1) [3, 4, 5, 1]
(%i2) ele2pui (3, %);
(%o2) [3, 4, 6, 7]
(%i3) pui2polynome (x, %);
(%o3) x^3 - 4 x^2 + 5 x - 1
```

Autres fonctions à voir : polynome2ele, ele2polynome.

**pui\_direct** (*orbite*, [*lvar\_1*, ..., *lvar\_n*], [*d\_1*, *d\_2*, ..., *d\_n*])

Function

Soit *f* un polynôme en *n* blocs de variables *lvar\_1*, ..., *lvar\_n*. Soit *c\_i* le nombre de variables dans *lvar\_i*. Et *SC* le produit des *n* groupes symétriques de degré *c\_1*, ..., *c\_n*. Ce groupe agit naturellement sur *f*. La liste *orbite* est l'orbite, notée *SC(f)*, de la fonction *f* sous l'action de *SC*. (Cette liste peut être obtenue avec la fonction : multi\_orbit). Les *d\_i* sont des entiers tels que *c\_1* ≤ *d\_1*, *c\_2* ≤ *d\_2*, ..., *c\_n* ≤ *d\_n*. Soit *SD* le produit des groupes symétriques *S\_d1* × *S\_d2* × ... × *S\_dn*.

La fonction *pui\_direct* ramène les *n* premières fonctions puissances de *SD(f)* de'duites des fonctions puissances de *SC(f)* ou *n* est le cardinal de *SD(f)*.

Le résultat est rendu sous forme multi-contractée par rapport à *SD*. i.e. on ne conserve qu'un élément par orbite sous l'action de *SD*).

```
(%i1) l: [[x, y], [a, b]];
(%o1) [[x, y], [a, b]]
(%i2) pui_direct (multi_orbit (a*x + b*y, 1), 1, [2, 2]);
(%o2) [a x, 4 a b x y + a x ]
(%i3) pui_direct (multi_orbit (a*x + b*y, 1), 1, [3, 2]);
(%o3) [2 a x, 4 a b x y + 2 a x , 3 a b x y + 2 a x ,
      2 2 2 2      3 3      4 4
      12 a b x y + 4 a b x y + 2 a x ,
      3 2 3 2      4 4      5 5
      10 a b x y + 5 a b x y + 2 a x ,
      3 3 3 3      4 2 4 2      5 5      6 6
      40 a b x y + 15 a b x y + 6 a b x y + 2 a x ]
```



```
(%i4) pui_direct ([y + x + 2*c, y + x + 2*b, y + x + 2*a], [[x, y], [a, b, c]])
(%o4) [3 x + 2 a, 6 x y + 3 x2 + 4 a x + 4 a2,
      9 x2 y + 12 a x y + 3 x3 + 6 a x2 + 12 a2 x + 8 a3]
```

**puireduc** (*n*, *lpui*)

Function

*lpui* est une liste dont le premier e'le'ment est un entier *m*. **puireduc** donne les *n* preme'eres fonctions puissances en fonction des *m* preme'eres.

```
(%i1) puireduc (3, [2]);
```

```
(%o1) [2, p1, p2, p1 p2 -  $\frac{p1^2 (p1^2 - p2^2)}{2}$ ]
```

```
(%i2) ratsimp (%);
```

```
(%o2) [2, p1, p2,  $\frac{3 p1 p2^3 - p1^3}{2}$ ]
```

**resolvante** (*P*, *x*, *f*, [*x*<sub>1</sub>, ..., *x*<sub>*d*</sub>])

Function

calcule la re'solvante du polyno^me *P* de la variable *x* et de degre' *n* >= *d* par la fonction *f* exprime'e en les variables *x*<sub>1</sub>, ..., *x*<sub>*d*</sub>. Il est important pour l'efficacite' des calculs de ne pas mettre dans la liste [*x*<sub>1</sub>, ..., *x*<sub>*d*</sub>] les variables n'intervenant pas dans la fonction de transformation *f*.

Afin de rendre plus efficaces les calculs on peut mettre des drapeaux a' la variable **resolvante** afin que des algorithmes ade'quates soient utilise's :

Si la fonction *f* est unitaire :

- un polyno^me d'une variable,
- line'aire ,
- alterne'e,
- une somme de variables,
- syme'trique en les variables qui apparaissent dans son expression,
- un produit de variables,
- la fonction de la re'solvante de Cayley (utilisable qu'en degre' 5)

$$(x_1 x_2 + x_2 x_3 + x_3 x_4 + x_4 x_5 + x_5 x_1 - (x_1 x_3 + x_3 x_5 + x_5 x_2 + x_2 x_4 + x_4 x_1))^2$$

generale,

le drapeau de **resolvante** pourra e^tre respectivement :

- unitaire,
- lineaire,
- alternee,
- somme,

- produit,
- cayley,
- generale.

```
(%i1) resolvante: unitaire$
(%i2) resolvante (x^7 - 14*x^5 + 56*x^3 - 56*x + 22, x, x^3 - 1, [x]);■

" resolvante unitaire " [7, 0, 28, 0, 168, 0, 1120, - 154, 7840, - 2772, 56448
413952, - 352352, 3076668, - 3363360, 23114112, - 30494464,
175230832, - 267412992, 1338886528, - 2292126760]
      3      6      3      9      6      3
[x  - 1, x  - 2 x  + 1, x  - 3 x  + 3 x  - 1,
      12      9      6      3      15      12      9      6      3
x  - 4 x  + 6 x  - 4 x  + 1, x  - 5 x  + 10 x  - 10 x  + 5 x
      18      15      12      9      6      3
- 1, x  - 6 x  + 15 x  - 20 x  + 15 x  - 6 x  + 1,
      21      18      15      12      9      6      3
x  - 7 x  + 21 x  - 35 x  + 35 x  - 21 x  + 7 x  - 1]
[- 7, 1127, - 6139, 431767, - 5472047, 201692519, - 3603982011]
      7      6      5      4      3      2
(%o2) y  + 7 y  - 539 y  - 1841 y  + 51443 y  + 315133 y
      + 376999 y + 125253

(%i3) resolvante: lineaire$
(%i4) resolvante (x^4 - 1, x, x1 + 2*x2 + 3*x3, [x1, x2, x3]);

" resolvante lineaire "
      24      20      16      12      8
(%o4) y  + 80 y  + 7520 y  + 1107200 y  + 49475840 y
      + 344489984 y  + 655360000

(%i5) resolvante: general$
(%i6) resolvante (x^4 - 1, x, x1 + 2*x2 + 3*x3, [x1, x2, x3]);

" resolvante generale "
      24      20      16      12      8
(%o6) y  + 80 y  + 7520 y  + 1107200 y  + 49475840 y
      + 344489984 y  + 655360000

(%i7) resolvante (x^4 - 1, x, x1 + 2*x2 + 3*x3, [x1, x2, x3, x4]);

" resolvante generale "
```

```

(%o7)  $y^{24} + 80 y^{20} + 7520 y^{16} + 1107200 y^{12} + 49475840 y^8$ 
+ 344489984  $y^4 + 655360000$ 
(%i8) direct ([x^4 - 1], x, x1 + 2*x2 + 3*x3, [[x1, x2, x3]]);
(%o8)  $y^{24} + 80 y^{20} + 7520 y^{16} + 1107200 y^{12} + 49475840 y^8$ 
+ 344489984  $y^4 + 655360000$ 
(%i9) resolvante :lineaire$
(%i10) resolvante (x^4 - 1, x, x1 + x2 + x3, [x1, x2, x3]);
" resolvante lineaire "
(%o10)  $y^4 - 1$ 
(%i11) resolvante: symetrique$
(%i12) resolvante (x^4 - 1, x, x1 + x2 + x3, [x1, x2, x3]);
" resolvante symetrique "
(%o12)  $y^4 - 1$ 
(%i13) resolvante (x^4 + x + 1, x, x1 - x2, [x1, x2]);
" resolvante symetrique "
(%o13)  $y^6 - 4 y^2 - 1$ 
(%i14) resolvante: alternee$
(%i15) resolvante (x^4 + x + 1, x, x1 - x2, [x1, x2]);
" resolvante alternee "
(%o15)  $y^{12} + 8 y^8 + 26 y^6 - 112 y^4 + 216 y^2 + 229$ 
(%i16) resolvante: produit$
(%i17) resolvante (x^7 - 7*x + 3, x, x1*x2*x3, [x1, x2, x3]);
" resolvante produit "
(%o17)  $y^{35} - 7 y^{33} - 1029 y^{29} + 135 y^{28} + 7203 y^{27} - 756 y^{26}$ 
+ 1323  $y^{24} + 352947 y^{23} - 46305 y^{22} - 2463339 y^{21} + 324135 y^{20}$ 
- 30618  $y^{19} - 453789 y^{18} - 40246444 y^{17} + 282225202 y^{15}$ 
- 44274492  $y^{14} + 155098503 y^{12} + 12252303 y^{11} + 2893401 y^{10}$ 

```

```

          9          8          7          6
- 171532242 y + 6751269 y + 2657205 y - 94517766 y

          5          3
- 3720087 y + 26040609 y + 14348907
(%i18) resolvante: symetrique$
(%i19) resolvante (x^7 - 7*x + 3, x, x1*x2*x3, [x1, x2, x3]);

" resolvante symetrique "
          35          33          29          28          27          26
(%o19) y - 7 y - 1029 y + 135 y + 7203 y - 756 y

          24          23          22          21          20
+ 1323 y + 352947 y - 46305 y - 2463339 y + 324135 y

          19          18          17          15
- 30618 y - 453789 y - 40246444 y + 282225202 y

          14          12          11          10
- 44274492 y + 155098503 y + 12252303 y + 2893401 y

          9          8          7          6
- 171532242 y + 6751269 y + 2657205 y - 94517766 y

          5          3
- 3720087 y + 26040609 y + 14348907
(%i20) resolvante: cayley$
(%i21) resolvante (x^5 - 4*x^2 + x + 1, x, a, []);

" resolvante de Cayley "
          6          5          4          3          2
(%o21) x - 40 x + 4080 x - 92928 x + 3772160 x + 37880832 x

+ 93392896

```

Pour la re'solvante de Cayley, les 2 derniers arguments sont neutres et le polyno<sup>me</sup> donne' en entre'e doit ne'cessairement e<sup>tre</sup> de degre' 5.

Voir e'galement :

resolvante\_bipartite, resolvante\_produit\_sym, resolvante\_unitaire,  
 resolvante\_alternee1, resolvante\_klein, resolvante\_klein3, resolvante\_  
 vierer, resolvante\_diedrale.

### resolvante\_alternee1 ( $P, x$ )

Function

calcule la transformation de  $P(x)$  de degre  $n$  par la fonction  $\prod_{1 \leq i < j \leq n-1} (x_i - x_j)$ .

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante , resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale, resolvante\_bipartite.

**resolvante\_bipartite** ( $P, x$ ) Function

calcule la transformation de  $P(x)$  de degre  $n$  ( $n$  pair) par la fonction  $x_1 x_2 \dots x_{n/2} + x_{n/2+1} \dots x_n$

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante , resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale, resolvante\_alternee1.

```
(%i1) resolvante_bipartite (x^6 + 108, x);
              10      8      6      4
(%o1)      y  - 972 y  + 314928 y  - 34012224 y
```

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante, resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale, resolvante\_alternee1.

**resolvante\_diedrale** ( $P, x$ ) Function

calcule la transformation de  $P(x)$  par la fonction  $x_1 x_2 + x_3 x_4$ .

```
(%i1) resolvante_diedrale (x^5 - 3*x^4 + 1, x);
      15      12      11      10      9      8      7
(%o1) x  - 21 x  - 81 x  - 21 x  + 207 x  + 1134 x  + 2331 x
      6      5      4      3      2
      - 945 x  - 4970 x  - 18333 x  - 29079 x  - 20745 x  - 25326 x
      - 697
```

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante\_alternee1, resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante.

**resolvante\_klein** ( $P, x$ ) Function

calcule la transformation de  $P(x)$  par la fonction  $x_1 x_2 x_4 + x_4$ .

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante\_alternee1, resolvante, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale.

**resolvante\_klein3** ( $P, x$ ) Function

calcule la transformation de  $P(x)$  par la fonction  $x_1 x_2 x_4 + x_4$ .

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante\_alternee1, resolvante\_klein, resolvante, resolvante\_vierer, resolvante\_diedrale.

**resolvante\_produit\_sym** ( $P, x$ ) Function

calcule la liste toutes les r'esolvantes produit du polyn\^ome  $P(x)$ .

```
(%i1) resolvante_produit_sym (x^5 + 3*x^4 + 2*x - 1, x);
(%o1) [y^5 + 3 y^4 + 2 y^3 - 1, y^10 - 2 y^8 - 21 y^7 - 31 y^6 - 14 y^5
      - y^4 + 14 y^3 + 3 y^2 + 1, y^10 + 3 y^8 + 14 y^7 - y^6 - 14 y^5 - 31 y^4
      - 21 y^3 - 2 y^2 + 1, y^5 - 2 y^4 - 3 y^3 - 1, y - 1]
(%i2) resolvante: produit$
(%i3) resolvante (x^5 + 3*x^4 + 2*x - 1, x, a*b*c, [a, b, c]);

" resolvante produit "
(%o3) y^10 + 3 y^8 + 14 y^7 - y^6 - 14 y^5 - 31 y^4 - 21 y^3 - 2 y^2 + 1
```

Voir e'galement :

resolvante, resolvante\_unitaire, resolvante\_alternee1, resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale.

**resolvante\_unitaire** ( $P, Q, x$ ) Function

calcule la r'esolvante du polyn\^ome  $P(x)$  par le polyn\^ome  $Q(x)$ .

Voir e'galement :

resolvante\_produit\_sym, resolvante, resolvante\_alternee1, resolvante\_klein, resolvante\_klein3, resolvante\_vierer, resolvante\_diedrale.

**resolvante\_vierer** ( $P, x$ ) Function

calcule la transformation de  $P(x)$  par la fonction  $x_1 x_2 - x_3 x_4$ .

Voir e'galement :

resolvante\_produit\_sym, resolvante\_unitaire, resolvante\_alternee1, resolvante\_klein, resolvante\_klein3, resolvante, resolvante\_diedrale.

**schur2comp** ( $P, Lvar$ ) Function

$P$  est un polyno\^mes en les variables contenues dans la liste  $Lvar$ . Chacune des variables de  $Lvar$  repre'sente une fonction syme'trique comple'te. On repre'sente dans  $Lvar$  la ie'me fonction syme'trique comple'te comme la concate'nation de la lettre  $h$  avec l'entier  $i$  :  $h_i$ . Cette fonction donne l'expression de  $P$  en fonction des fonctions de Schur.

```
(%i1) schur2comp (h1*h2 - h3, [h1, h2, h3]);
(%o1) s
      1, 2
(%i2) schur2comp (a*h3, [h3]);
(%o2) s a
      3
```

**somrac** ( $l, k$ ) Function  
 la liste  $l$  contient les fonctions syme'triques e'l'ementaires d'un polyno^me  $P$ . On  
 calcul le polyno^mes dont les racines sont les sommes  $K$  a'  $K$  distinctes des racines  
 de  $P$ .

Voir e'galement `prodrac`.

**tcontract** ( $pol, lvar$ ) Function  
 teste si le polyno^me  $pol$  est syme'trique en les variables contenues dans la liste  $lvar$ .  
 Si oui il rend une forme contracte'e comme la fonction `contract`.

Autres fonctions de changements de repre'sentations :

`contract, cont2part, expose, part2cont, partpol, tpartpol`.

**tpartpol** ( $pol, lvar$ ) Function  
 teste si le polyno^me  $pol$  est syme'trique en les variables contenues dans la liste  $lvar$ .  
 Si oui il rend sa forme partitionne'e comme la fonction `partpol`.

Autres fonctions de changements de repre'sentations :

`contract, cont2part, expose, part2cont, partpol, tcontract`.

**treillis** ( $n$ ) Function  
 rame'ne toutes les partitions de poids  $n$ .

```
(%i1) treillis (4);
```

```
(%o1) [[4], [3, 1], [2, 2], [2, 1, 1], [1, 1, 1, 1]]
```

Voir e'galement : `lgtreillis, ltreillis` et `treinat`.

**treinat** ( $part$ ) Function  
 rame'ne la liste des partitions infe'rieures a' la partition  $part$  pour l'ordre naturel.

```
(%i1) treinat ([5]);
```

```
(%o1) [[5]]
```

```
(%i2) treinat ([1, 1, 1, 1, 1]);
```

```
(%o2) [[5], [4, 1], [3, 2], [3, 1, 1], [2, 2, 1], [2, 1, 1, 1],
```

```
[1, 1, 1, 1, 1]]
```

```
(%i3) treinat ([3, 2]);
```

```
(%o3) [[5], [4, 1], [3, 2]]
```

Voir e'galement : `lgtreillis, ltreillis` et `treillis`.

## 35 Groups

### 35.1 Definitions for Groups

**todd\_coxeter** (*relations, subgroup*)

Function

**todd\_coxeter** (*relations*)

Function

Find the order of  $G/H$  where  $G$  is the Free Group modulo *relations*, and  $H$  is the subgroup of  $G$  generated by *subgroup*. *subgroup* is an optional argument, defaulting to  $[]$ . In doing this it produces a multiplication table for the right action of  $G$  on  $G/H$ , where the cosets are enumerated  $[H, Hg_2, Hg_3, \dots]$ . This can be seen internally in the `$todd_coxeter_state`.

The multiplication tables for the variables are in `table:todd_coxeter_state[2]`. Then `table[i]` gives the table for the  $i$ th variable. `mulcoset(coset,i) := table[varnum][coset];`

Example:

```
(%i1) symet(n):=create_list(
      if (j - i) = 1 then (p(i,j))3 else
      if (not i = j) then (p(i,j))2 else
      p(i,i) , j, 1, n-1, i, 1, j);
      <3>
(%o1) symet(n) := create_list(if j - i = 1 then p(i, j)
      <2>
      else (if not i = j then p(i, j) else p(i, i)), j, 1, n - 1,
i, 1, j)
(%i2) p(i,j) := concat(x,i).concat(x,j);
(%o2) p(i, j) := concat(x, i) . concat(x, j)
(%i3) symet(5);
      <2> <3> <2> <2> <3>
(%o3) [x1 , (x1 . x2) , x2 , (x1 . x3) , (x2 . x3) ,
      <2> <2> <2> <3> <2>
      x3 , (x1 . x4) , (x2 . x4) , (x3 . x4) , x4 ]
(%i4) todd_coxeter(%o3);

Rows tried 426
(%o4) 120
(%i5) todd_coxeter(%o3,[x1]);

Rows tried 213
(%o5) 60
(%i6) todd_coxeter(%o3,[x1,x2]);

Rows tried 71
(%o6) 20
```



```
(%i7) table:todd_coxeter_state[2]$
(%i8) table[1];
(%o8) {Array: (SIGNED-BYTE 30) #(0 2 1 3 7 6 5 4 8 11 17 9 12 14 #
13 20 16 10 18 19 15 0 0 0 0 0 0 0 0 0 0 0 0
0 0 0)}
```

Note only the elements 1 thru 20 of this array %o8 are meaningful. `table[1][4] = 7` indicates `coset4.var1 = coset7`

## 36 Runtime Environment

### 36.1 Introduction for Runtime Environment

`maxima-init.mac` is a file which is loaded automatically when Maxima starts. You can use `maxima-init.mac` to customize your Maxima environment. `maxima-init.mac`, if it exists, is typically placed in the directory named by `:lisp (default-userdir)`, although it can be in any directory searched by the function `file_search`.

Here is an example `maxima-init.mac` file:

```
setup_autoload ("specfun.mac", ultraspherical, assoc_legendre_p);
showtime:all;
```

In this example, `setup_autoload` tells Maxima to load the specified file (`specfun.mac`) if any of the functions (`ultraspherical`, `assoc_legendre_p`) are called but not yet defined. Thus you needn't remember to load the file before calling the functions.

The statement `showtime: all` tells Maxima to set the `showtime` variable. The `maxima-init.mac` file can contain any other assignments or other Maxima statements.

### 36.2 Interrupts

The user can stop a time-consuming computation with the `^C` (control-C) character. The default action is to stop the computation and print another user prompt. In this case, it is not possible to restart a stopped computation.

If the variable `*debugger-hook*` is set to `nil`, by executing

```
:lisp (setq *debugger-hook* nil)
```

then upon receiving `^C`, Maxima will enter the Lisp debugger, and the user may use the debugger to inspect the Lisp environment. The stopped computation can be restarted by entering `continue` in the Lisp debugger. The means of returning to Maxima from the Lisp debugger (other than running the computation to completion) is different for each version of Lisp.

On Unix systems, the character `^Z` (control-Z) causes Maxima to stop altogether, and control is returned to the shell prompt. The `fg` command causes Maxima to resume from the point at which it was stopped.

### 36.3 Definitions for Runtime Environment

#### feature

declaration

Maxima understands two distinct types of features, system features and features which apply to mathematical expressions. See also `status` for information about system features. See also `features` and `featurep` for information about mathematical features.

`feature` itself is not the name of a function or variable.

**featurep** (*a*, *f*) Function

Attempts to determine whether the object *a* has the feature *f* on the basis of the facts in the current database. If so, it returns **true**, else **false**. See also **features**.

```
(%i1) declare (j, even)$
(%i2) featurep (j, integer);
(%o2)                                     true
```

**room** () Function

**room** (*true*) Function

**room** (*false*) Function

Prints out a description of the state of storage and stack management in Maxima.

**room** calls the Lisp function of the same name.

- **room** () prints out a moderate description.
- **room** (**true**) prints out a verbose description.
- **room** (**false**) prints out a terse description.

**status** (*feature*) Function

**status** (*feature*, *putative\_feature*) Function

**status** (*status*) Function

Returns information about the presence or absence of certain system-dependent features.

- **status** (*feature*) returns a list of system features. These include Lisp version, operating system type, etc. The list may vary from one Lisp type to another.
- **status** (*feature*, *putative\_feature*) returns **true** if *putative\_feature* is on the list of items returned by **status** (*feature*) and **false** otherwise. **status** quotes the argument *putative\_feature*. The double single quotes operator, `''`, defeats the quotation. A feature whose name contains a special character, such as a hyphen, must be given as a string argument. For example, **status** (*feature*, "ansi-cl").
- **status** (*status*) returns a two-element list [*feature*, *status*]. *feature* and *status* are the two arguments accepted by the **status** function; it is unclear if this list has additional significance.

The variable **features** contains a list of features which apply to mathematical expressions. See **features** and **featurep** for more information.

**time** (%o1, %o2, %o3, ...) Function

Returns a list of the times, in seconds, taken to compute the output lines %o1, %o2, %o3, .... The time returned is Maxima's estimate of the internal computation time, not the elapsed time. **time** can only be applied to output line variables; for any other variables, **time** returns **unknown**.

Set **showtime: true** to make Maxima print out the computation time and elapsed time with each output line.

## 37 Miscellaneous Options

### 37.1 Introduction to Miscellaneous Options

In this section various options are discussed which have a global effect on the operation of Maxima. Also various lists such as the list of all user defined functions, are discussed.

### 37.2 Share

The Maxima "share" directory contains programs and other files of interest to Maxima users, but not part of the core implementation of Maxima. These programs are typically loaded via `load` or `setup_autoload`.

`:lisp *maxima-sharedir*` displays the location of the share directory within the user's file system.

`printfile ("share.usg")` prints an out-of-date list of share packages. Users may find it more informative to browse the share directory using a file system browser.

### 37.3 Definitions for Miscellaneous Options

#### aliases

Variable

Default value: []

`aliases` is the list of atoms which have a user defined alias (set up by the `alias`, `ordergreat`, `orderless` functions or by declaring the atom a noun with `declare`).

#### alphabetic

declaration

`declare (char, alphabetic)` adds `char` to Maxima's alphabet, which initially contains the letters A through Z, a through z, % and `_`. `char` is specified as a string of length 1, e.g., "~".

```
(%i1) declare ("~", alphabetic);
(%o1)                                     done
(%i2) foo~bar;
(%o2)                                     foo~bar
(%i3) atom (foo~bar);
(%o3)                                     true
```

#### apropos (string)

Function

Searches for Maxima names which have `string` appearing anywhere within them. Thus, `apropos (exp)` returns a list of all the flags and functions which have `exp` as part of their names, such as `expand`, `exp`, and `exponentialize`. Thus if you can only remember part of the name of something you can use this command to find the rest of the name. Similarly, you could say `apropos (tr_)` to find a list of many of the switches relating to the translator, most of which begin with `tr_`.

- args** (*expr*) Function  
 Returns the list of arguments of *expr*, which may be any kind of expression other than an atom. Only the arguments of the top-level operator are extracted; subexpressions of *expr* appear as elements or subexpressions of elements of the list of arguments. The order of the items in the list may depend on the global flag `inflag`.  
*args* (*expr*) is equivalent to `substpart ("[" , expr, 0)`. See also `substpart`.  
 See also `op`.
- genindex** Variable  
 Default value: `i`  
*genindex* is the alphabetic prefix used to generate the next variable of summation when necessary.
- gensumnum** Variable  
 Default value: `0`  
*gensumnum* is the numeric suffix used to generate the next variable of summation. If it is set to `false` then the index will consist only of *genindex* with no numeric suffix.
- inf** Variable  
 Real positive infinity.
- infinity** Variable  
 Complex infinity, an infinite magnitude of arbitrary phase angle. See also `inf` and `minf`.
- infolists** Variable  
 Default value: `[]`  
*infolists* is a list of the names of all of the information lists in Maxima. These are:  
`labels` - all bound `%i`, `%o`, and `%t` labels.  
`values` - all bound atoms which are user variables, not Maxima options or switches, created by `:` or `::` or functional binding.  
`functions` - all user-defined functions, created by `:=`.  
`arrays` - declared and undeclared arrays, created by `:`, `::`, or `:=`.  
`macros` - any macros defined by the user.  
`myoptions` - all options ever reset by the user (whether or not they are later reset to their default values).  
`rules` - user-defined pattern matching and simplification rules, created by `tellsimp`, `tellsimpafter`, `defmatch`, or `defrule`.  
`aliases` - atoms which have a user-defined alias, created by the `alias`, `ordergreat`, `orderless` functions or by declaring the atom as a noun with `declare`.  
`dependencies` - atoms which have functional dependencies, created by the `depends` or `gradef` functions.

**gradefs** - functions which have user-defined derivatives, created by the **gradef** function.

**props** - atoms which have any property other than those mentioned above, such as **atvalues**, **matchdeclares**, etc., as well as properties specified in the **declare** function.

**let\_rule\_packages** - a list of all the user-defined let rule packages plus the special package **default\_let\_rule\_package**. (**default\_let\_rule\_package** is the name of the rule package used when one is not explicitly set by the user.)

**integerp** (*expr*) Function

Returns **true** if *expr* is a literal numeric integer, otherwise **false**.

**integerp** returns false if its argument is a symbol, even if the argument is declared **integer**.

Examples:

```
(%i1) integerp (0);
(%o1) true
(%i2) integerp (1);
(%o2) true
(%i3) integerp (-17);
(%o3) true
(%i4) integerp (0.0);
(%o4) false
(%i5) integerp (1.0);
(%o5) false
(%i6) integerp (%pi);
(%o6) false
(%i7) integerp (n);
(%o7) false
(%i8) declare (n, integer);
(%o8) done
(%i9) integerp (n);
(%o9) false
```

**m1pbranch** Variable

Default value: **false**

**m1pbranch** is the principal branch for  $-1$  to a power. Quantities such as  $(-1)^{(1/3)}$  (that is, an "odd" rational exponent) and  $(-1)^{(1/4)}$  (that is, an "even" rational exponent) are handled as follows:

domain:real

```
(-1)^(1/3):    -1
(-1)^(1/4):    (-1)^(1/4)
```

domain:complex

```
m1pbranch:false      m1pbranch:true
(-1)^(1/3)           1/2+%i*sqrt(3)/2
(-1)^(1/4)           sqrt(2)/2+%i*sqrt(2)/2
```

**numberp** (*expr*) Function

Returns **true** if *expr* is a literal integer, rational number, floating point number, or bigfloat, otherwise **false**.

**numberp** returns false if its argument is a symbol, even if the argument is a symbolic number such as `%pi` or `%i`, or declared to be even, odd, integer, rational, irrational, real, imaginary, or complex.

Examples:

```
(%i1) numberp (42);
(%o1) true
(%i2) numberp (-13/19);
(%o2) true
(%i3) numberp (3.14159);
(%o3) true
(%i4) numberp (-1729b-4);
(%o4) true
(%i5) map (numberp, [%e, %pi, %i, %phi, inf, minf]);
(%o5) [false, false, false, false, false, false]
(%i6) declare (a, even, b, odd, c, integer, d, rational,
e, irrational, f, real, g, imaginary, h, complex);
(%o6) done
(%i7) map (numberp, [a, b, c, d, e, f, g, h]);
(%o7) [false, false, false, false, false, false, false, false]
```

**properties** (*a*) Function

Returns a list of the names of all the properties associated with the atom *a*.

**props** special symbol

**props** are atoms which have any property other than those explicitly mentioned in **infolists**, such as **atvalues**, **matchdeclares**, etc., as well as properties specified in the **declare** function.

**propvars** (*prop*) Function

Returns a list of those atoms on the **props** list which have the property indicated by *prop*. Thus **propvars** (**atvalue**) returns a list of atoms which have **atvalues**.

**put** (*atom, value, indicator*) Function

Assigns *value* to the property (specified by *indicator*) of *atom*. *indicator* may be the name of any property, not just a system-defined property.

**put** evaluates its arguments. **put** returns *value*.

Examples:

```
(%i1) put (foo, (a+b)^5, expr);
(%o1) (b + a)^5
(%i2) put (foo, "Hello", str);
(%o2) Hello
(%i3) properties (foo);
```

```

(%o3)          [[user properties, str, expr]]
(%i4) get (foo, expr);

(%o4)          5
              (b + a)
(%i5) get (foo, str);
(%o5)          Hello

```

**qput** (*atom, value, indicator*) Function

Assigns *value* to the property (specified by *indicator*) of *atom*. This is the same as **put**, except that the arguments are quoted.

Example:

```

(%i1) foo: aa$
(%i2) bar: bb$
(%i3) baz: cc$
(%i4) put (foo, bar, baz);
(%o4)          bb
(%i5) properties (aa);
(%o5)          [[user properties, cc]]
(%i6) get (aa, cc);
(%o6)          bb
(%i7) qput (foo, bar, baz);
(%o7)          bar
(%i8) properties (foo);
(%o8)          [value, [user properties, baz]]
(%i9) get ('foo, 'baz);
(%o9)          bar

```

**rem** (*atom, indicator*) Function

Removes the property indicated by *indicator* from *atom*.

**remove** (*atom\_1, property\_1, ..., atom\_n, property\_n*) Function

**remove** (*[atom\_1, ..., atom\_m], [property\_1, ..., property\_n], ...*) Function

Removes properties associated with atoms.

**remove** (*atom\_1, property\_1, ..., atom\_n, property\_n*) removes *property\_k* from *atom\_k*.

**remove** (*[atom\_1, ..., atom\_m], [property\_1, ..., property\_n], ...*) removes all properties *property\_1, ..., property\_n* from all atoms *atom\_1, ..., atom\_m*. There may be more than one pair of lists.

The removed properties may be system-defined properties such as **function** or **mode\_declare**, or user-defined properties.

A property may be **transfun** to remove the translated Lisp version of a function. After executing this, the Maxima version of the function is executed rather than the translated version.

A property may be **op** or **operator** to remove a syntax extension given to an atom.

If an atom is "all" then the property is removed from all atoms which have it.



`remove` always returns `done` whether or not an atom has a specified property. This behavior is unlike the more specific remove functions (`remvalue`, `remarray`, `remfunction`, and `remrule`).

**remvalue** (*name\_1*, ..., *name\_n*) Function  
**remvalue** (*all*) Function  
 Removes the values of user variables *name\_1*, ..., *name\_n* (which can be subscripted) from the system.  
**remvalue** (*all*) removes the values of all variables in *values*, the list of all variables given names by the user (as opposed to those which are automatically assigned by Maxima).  
 See also *values*.

**rncombine** (*expr*) Function  
 Transforms *expr* by combining all terms of *expr* that have identical denominators or denominators that differ from each other by numerical factors only. This is slightly different from the behavior of `combine`, which collects terms that have identical denominators.  
 Setting `pformat: true` and using `combine` yields results similar to those that can be obtained with `rncombine`, but `rncombine` takes the additional step of cross-multiplying numerical denominator factors. This results in neater forms, and the possibility of recognizing some cancellations.

**scalarp** (*expr*) Function  
 Returns `true` if *expr* is a number, constant, or variable declared `scalar` with `declare`, or composed entirely of numbers, constants, and such variables, but not containing matrices or lists.

**setup\_autoload** (*filename*, *function\_1*, ..., *function\_n*) Function  
 Specifies that if any of *function\_1*, ..., *function\_n* are referenced and not yet defined, *filename* is loaded via `load`. *filename* usually contains definitions for the functions specified, although that is not enforced.  
**setup\_autoload** does not work for array functions.  
**setup\_autoload** quotes its arguments.

Example:

```
(%i1) legendre_p (1, %pi);
(%o1)          legendre_p(1, %pi)
(%i2) setup_autoload ("specfun.mac", legendre_p, ultraspherical);
(%o2)          done
(%i3) ultraspherical (2, 1/2, %pi);
Warning - you are redefining the Macsyma function ultraspherical
Warning - you are redefining the Macsyma function legendre_p
          2
          3 (%pi - 1)
(%o3)    ----- + 3 (%pi - 1) + 1
```

```

                2
(%i4) legendre_p (1, %pi);
(%o4)          %pi
(%i5) legendre_q (1, %pi);
                %pi + 1
                %pi log(-----)
(%o5)          ----- - 1
                2

```



## 38 Rules and Patterns

### 38.1 Introduction to Rules and Patterns

This section discusses user defined pattern matching and simplification rules (set up by `tellsimp`, `tellsimpafter`, `defmatch`, or, `defrule`.) You may affect the main simplification procedures, or else have your rules applied explicitly using `apply1` and `apply2`. There are additional mechanisms for polynomials rules under `tellrat`, and for commutative and non commutative algebra in chapter on `affine`.

### 38.2 Definitions for Rules and Patterns

**apply1** (*expr*, *rule\_1*, ..., *rule\_n*) Function  
 repeatedly applies the first rule to *exp* until it fails, then repeatedly applies the same rule to all subexpressions of *exp*, left-to-right, until the first rule has failed on all subexpressions. Call the result of transforming *exp* in this manner *exp'*. Then the second rule is applied in the same fashion starting at the top of *exp'*. When the final rule fails on the final subexpression, the application is finished.

**apply2** (*expr*, *rule\_1*, ..., *rule\_n*) Function  
 differs from `apply1` in that if the first rule fails on a given subexpression, then the second rule is repeatedly applied, etc. Only if they all fail on a given subexpression is the whole set of rules repeatedly applied to the next subexpression. If one of the rules succeeds, then the same subexpression is reprocessed, starting with the first rule. `maxapplydepth` is the maximum depth to which `apply1` and `apply2` will delve.

**applyb1** (*expr*, *rule\_1*, ..., *rule\_n*) Function  
 is similar to `apply1` but works from the "bottom up" instead of from the "top down". That is, it processes the smallest subexpression of *exp*, then the next smallest, etc. `maxapplyheight` is the maximum height to which `applyb1` will reach before giving up.

**current\_let\_rule\_package** Variable  
 default:[`default_let_rule_package`] - the name of the rule package that is presently being used. The user may reset this variable to the name of any rule package previously defined via the `let` command. Whenever any of the functions comprising the `let` package are called with no package name the value of  
`current_let_rule_package`  
 is used. If a call such as `letsimp (expr, rule_pkg_name)` is made, the rule package `rule_pkg_name` is used for that `letsimp` command only, i.e. the value of `current_let_rule_package` is not changed.

**default\_let\_rule\_package** Variable  
 - the name of the rule package used when one is not explicitly set by the user with `let` or by changing the value of `current_let_rule_package`.

**defmatch** (*progname*, *pattern*, *x<sub>1</sub>*, ..., *x<sub>n</sub>*) Function

Creates a function *progname* (*expr*, *y<sub>1</sub>*, ..., *y<sub>n</sub>*) which tests *expr* to see if it matches *pattern*.

*pattern* is an expression containing the pattern variables *x<sub>1</sub>*, ..., *x<sub>n</sub>* and pattern parameters, if any. The pattern variables are given explicitly as arguments to **defmatch** while the pattern parameters are declared by the **matchdeclare** function.

The first argument to the created function *progname* is an expression to be matched against the pattern and the other arguments are the actual variables *y<sub>1</sub>*, ..., *y<sub>n</sub>* in the expression which correspond to the dummy variables *x<sub>1</sub>*, ..., *x<sub>n</sub>* in the pattern.

If the match is successful, *progname* returns a list of equations whose left sides are the pattern variables and pattern parameters, and whose right sides are the expressions which the pattern variables and parameters matched. The pattern parameters, but not the variables, are assigned the subexpressions they match. If the match fails, *progname* returns **false**.

Any variables not declared as pattern parameters in **matchdeclare** or as variables in **defmatch** match only themselves.

A pattern which contains no pattern variables or parameters returns **true** if the match succeeds.

Examples:

```
(%i1) matchdeclare (a, freeof(x), b, freeof(x))$
(%i2) defmatch (linearp, a*x + b, x)$
```

This **defmatch** defines the function **linearp** (*expr*, *y*), which tests *expr* to see if it is of the form  $a*y + b$  such that *a* and *b* do not contain *y*.

```
(%i3) linearp (3*z + (y+1)*z + y^2, z);
(%o3)          2
          [b = y , a = y + 4, x = z]
(%i4) a;
(%o4)          y + 4
(%i5) b;
(%o5)          2
          y
```

If the third argument to the **defmatch** in (%i4) had been omitted, then **linear** would only match expressions linear in *X*, not in any other variable.

```
(%i1) matchdeclare ([a, f], true)$
(%i2) constinterval (l, h) := constantp (h - l)$
(%i3) matchdeclare (b, constinterval (a))$
(%i4) matchdeclare (x, atom)$
(%i5) (remove (integrate, outative),
      defmatch (checklimits, 'integrate (f, x, a, b)),
      declare (integrate, outative))$
(%i6) 'integrate (sin(t), t, %pi + x, 2*%pi + x);
(%o6)          x + 2 %pi
          /
          [
          I          sin(t) dt
          ]
```

```

          /
          x + %pi
(%i7) checklimits (%);
(%o7) [b = x + 2 %pi, a = x + %pi, x = t, f = sin(t)]
(%i8) a;
(%o8) x + %pi
(%i9) b;
(%o9) x + 2 %pi
(%i10) f;
(%o10) sin(t)
(%i11) x;
(%o11) t

```

**defrule** (*rulename*, *pattern*, *replacement*) Function

defines and names a replacement rule for the given pattern. If the rule named *rulename* is applied to an expression (by one of the **apply** functions below), every subexpression matching the pattern will be replaced by the replacement. All variables in the replacement which have been assigned values by the pattern match are assigned those values in the replacement which is then simplified. The rules themselves can be treated as functions which will transform an expression by one operation of the pattern match and replacement. If the pattern fails, the original expression is returned.

**disprule** (*rulename\_1*, *rulename\_2*, ...) Function

will display rules with the names *rulename1*, *rulename2*, as were given by **defrule**, **tellsimp**, or **tellsimpafter** or a pattern defined by **defmatch**. For example, the first rule modifying **sin** will be called **sinrule1**. **disprule (all)** will display all rules.

**let** (*prod*, *repl*, *predname*, *arg\_1*, ..., *arg\_n*) Function

**let** (*[prod, repl, predname, arg\_1, ..., arg\_n]*, *package\_name*) Function

Defines a substitution rule for **letsimp** such that *prod* is replaced by *repl*. *prod* is a product of positive or negative powers of the following terms:

- Atoms which **letsimp** will search for literally unless previous to calling **letsimp** the **matchdeclare** function is used to associate a predicate with the atom. In this case **letsimp** will match the atom to any term of a product satisfying the predicate.
- Kernels such as **sin(x)**, **n!**, **f(x,y)**, etc. As with atoms above **letsimp** will look for a literal match unless **matchdeclare** is used to associate a predicate with the argument of the kernel.

A term to a positive power will only match a term having at least that power in the expression being **letsimp**'ed. A term to a negative power on the other hand will only match a term with a power at least as negative. In the case of negative powers in *prod* the switch **letrat** must be set to **true**. See below for more on **letrat**.

If a predicate is included in the **let** function followed by a list of arguments, a tentative match (i.e. one that would be accepted if the predicate were omitted) will

be accepted only if `predname (arg1', ..., argn')` evaluates to `true` where `argi'` is the value matched to `argi`. The `argi` may be the name of any atom or the argument of any kernel appearing in `prod`. `repl` may be any rational expression. If any of the atoms or arguments from `prod` appear in `repl` the appropriate substitutions will be made.

`letrat` when `false`, `letsimp` will simplify the numerator and denominator of `expr` independently and return the result. Substitutions such as `n!/n` goes to `(n-1)!` will fail. To handle such situations `letrat` should be set to `true`, then the numerator, denominator, and their quotient will be simplified in that order.

These substitution functions allow you to work with several rule packages at once. Each rule package can contain any number of `let`'ed rules and is referred to by a user supplied name. `let ([prod, repl, predname, arg_1, ..., arg_n], package_name)` adds the rule `predname` to the rule package `package_name`. `letsimp (expr, package_name)` applies the rules in `package_name`. `letsimp (expr, package_name1, package_name2, ...)` is equivalent to `letsimp (expr, package_name1)` followed by `letsimp (% , package_name2), ...`

`current_let_rule_package` is the name of the rule package that is presently being used. The user may reset this variable to the name of any rule package previously defined via the `let` command. Whenever any of the functions comprising the `let` package are called with no package name, the `current_let_rule_package` is used. If a call such as `letsimp (expr, rule_pkg_name)` is made, the rule package `rule_pkg_name` is used for that `letsimp` command only, and `current_let_rule_package` is not changed. If not otherwise specified, `current_let_rule_package` defaults to `default_let_rule_package`.

```
(%i1) matchdeclare ([a, a1, a2], true)$
(%i2) oneless (x, y) := is (x = y-1)$
(%i3) let (a1*a2!, a1!, oneless, a2, a1);
(%o3)      a1 a2! --> a1! where oneless(a2, a1)
(%i4) letrat: true$
(%i5) let (a1!/a1, (a1-1)!);
          a1!
(%o5)      --- --> (a1 - 1)!
          a1
(%i6) letsimp (n*m!*(n-1)!/m);
(%o6)      (m - 1)! n!
(%i7) let (sin(a)^2, 1 - cos(a)^2);
          2          2
(%o7)      sin (a) --> 1 - cos (a)
(%i8) letsimp (sin(x)^4);
          4          2
(%o8)      cos (x) - 2 cos (x) + 1
```

## letrat

Variable

Default value: `false`

When `letrat` is `false`, `letsimp` simplifies the numerator and denominator of a ratio separately, and does not simplify the quotient.

When `letrat` is `true`, the numerator, denominator, and their quotient will be simplified in that order.

```
(%i1) matchdeclare (n, true)$
(%i2) let (n!/n, (n-1)!);
                                n!
(%o2)  -- --> (n - 1)!
                                n
(%i3) letrat: false$
(%i4) letsimp (a!/a);
                                a!
(%o4)  --
                                a
(%i5) letrat: true$
(%i6) letsimp (a!/a);
(%o6)  (a - 1)!
```

**letrules** () Function  
**letrules** (*package\_name*) Function

Displays the rules in a rule package. `letrules ()` displays the rules in the current rule package. `letrules (package_name)` displays the rules in `package_name`.

The current rule package is named by `current_let_rule_package`. If not otherwise specified, `current_let_rule_package` defaults to `default_let_rule_package`.

**letsimp** (*expr*) Function  
**letsimp** (*expr*, *package\_name*) Function

Applies the substitution rules previously defined by the function `let` until no further change is made to *expr*.

`letsimp (expr)` uses the rules from `current_let_rule_package`.

`letsimp (expr, package_name)` uses the rules from `package_name` without changing `current_let_rule_package`.

**let\_rule\_packages** Variable

default: [default\_let\_rule\_package]

`let_rule_packages` is a list of all the user-defined let rule packages plus the special package `default_let_rule_package`, which is the rule package used when a rule package is not otherwise specified.

**matchdeclare** (*a-1*, *pred-1*, ..., *a-n*, *pred-n*) Function

Associates a predicate *pred-k* with a variable or list of variables *a-k* so that *a-k* matches expressions for which the predicate is not `false`. Pattern matches are evaluated by one of the functions described below.

If the match succeeds then the variable is set to the matched expression. The predicate (in this case `freeof`) is written without the last argument which should be the one against which the pattern variable is to be tested. Note that the variable and the arguments to the predicate are evaluated at the time the match is evaluated.





```
(%i5) matchfix ("foo", "oof");
(%o5)                "foo"
(%i6) foo a, b, c oof;
(%o6)                fooa, b, coof
(%i7) >> w + foo x, y oof + z << / @ p, q ~;
                >>z + foox, yoof + w<<
(%o7)                -----
                        @p, q~
```

- Matchfix operators are ordinary user-defined functions.

```
(%i1) matchfix ("!-", "-!");
(%o1)                "!-"
(%i2) !- x, y -! := x/y - y/x;
(%o2)                !-x, y-! :=  $\frac{x}{y} - \frac{y}{x}$ 
(%i3) define (!-x, y-!, x/y - y/x);
(%o3)                !-x, y-! :=  $\frac{x}{y} - \frac{y}{x}$ 
(%i4) define ("!-" (x, y), x/y - y/x);
(%o4)                !-x, y-! :=  $\frac{x}{y} - \frac{y}{x}$ 
(%i5) dispfun ("!-");
(%t5)                !-x, y-! :=  $\frac{x}{y} - \frac{y}{x}$ 
(%o5)                done
(%i6) !-3, 5-!;
(%o6)                - --
                    16
(%i7) "!-" (3, 5);
(%o7)                - --
                    15
```

**remlet** (*prod*, *name*)

Function

deletes the substitution rule, *prod*  $\rightarrow$  *repl*, most recently defined by the **let** function. If *name* is supplied the rule is deleted from the rule package *name*. **REMLET**() and **REMLET(ALL)** delete all substitution rules from the current rule package. If the name of a rule package is supplied, e.g. **REMLET(ALL,name)**, the rule package, *name*, is also deleted. If a substitution is to be changed using the same product, **remlet** need not be called, just redefine the substitution using the same product (literally) with the **let** function and the new replacement and/or predicate name. Should **REMLET**(*product*) now be called the original substitution rule will be revived.

**remrule** (*function, rulename*) Function  
 will remove a rule with the name *rulename* from the function which was placed there by `defrule`, `defmatch`, `tellsimp`, or `tellsimpafter`. If *rule-name* is `all`, then all rules will be removed.

**tellsimp** (*pattern, replacement*) Function  
 is similar to `tellsimpafter` but places new information before old so that it is applied before the built-in simplification rules.

`tellsimp` is used when it is important to modify the expression before the simplifier works on it, for instance if the simplifier "knows" something about the expression, but what it returns is not to your liking. If the simplifier "knows" something about the main operator of the expression, but is simply not doing enough for you, you probably want to use `tellsimpafter`.

The pattern may not be a sum, product, single variable, or number.

*rules* is a list of names having simplification rules added to them by `defrule`, `defmatch`, `tellsimp`, or `tellsimpafter`.

Examples:

```
(%i1) matchdeclare (x, freeof (%i));
(%o1) done
(%i2) %iargs: false$
(%i3) tellsimp (sin(%i*x), %i*sinh(x));
(%o3) [sinrule1, simp-%sin]
(%i4) trigexpand (sin (%i*y + x));
(%o4) sin(x) cos(%i y) + %i cos(x) sinh(y)
(%i5) %iargs:true$
(%i6) errcatch(0^0);
0
0 has been generated
(%o6) []
(%i7) ev (tellsimp (0^0, 1), simp: false);
(%o7) [^rule1, simpexpt]
(%i8) 0^0;
(%o8) 1
(%i9) remrule ("^", %th(2)[1]);
(%o9) ^
(%i10) tellsimp (sin(x)^2, 1 - cos(x)^2);
(%o10) [^rule2, simpexpt]
(%i11) (1 + sin(x))^2;
(%o11) (sin(x) + 1)^2
(%i12) expand (%);
(%o12) 2 sin(x) - cos (x) + 2
(%i13) sin(x)^2;
(%o13) 1 - cos (x)
(%i14) kill (rules);
```

```
(%o14) done
(%i15) matchdeclare (a, true);
(%o15) done
(%i16) tellsimp (sin(a)^2, 1 - cos(a)^2);
(%o16) [^rule3, simpexpt]
(%i17) sin(y)^2;
(%o17) 1 - cos (y)2
```

**tellsimpafter** (*pattern, replacement*) Function  
defines a replacement for *pattern* which the Maxima simplifier uses after it applies the built-in simplification rules. The *pattern* may be anything but a single variable or a number.



## 39 Lists

### 39.1 Introduction to Lists

Lists are the basic building block for Maxima and Lisp. All data types other than arrays, hash tables, numbers are represented as Lisp lists, These Lisp lists have the form

```
((MPLUS) $A 2)
```

to indicate an expression  $a+2$ . At Maxima level one would see the infix notation  $a+2$ . Maxima also has lists which are printed as

```
[1, 2, 7, x+y]
```

for a list with 4 elements. Internally this corresponds to a Lisp list of the form

```
((MLIST) 1 2 7 ((MPLUS) $X $Y ))
```

The flag which denotes the type field of the Maxima expression is a list itself, since after it has been through the simplifier the list would become

```
((MLIST SIMP) 1 2 7 ((MPLUS SIMP) $X $Y))
```

### 39.2 Definitions for Lists

**append** (*list\_1, ..., list\_n*) Function

Returns a single list of the elements of *list\_1* followed by the elements of *list\_2*, ... **append** also works on general expressions, e.g. **append** ( $f(a,b)$ ,  $f(c,d,e)$ ); yields  $f(a,b,c,d,e)$ .

Do **example(append)**; for an example.

**assoc** (*key, list, default*) Function

**assoc** (*key, list*) Function

This function searches for the *key* in the left hand side of the input *list* of the form  $[x,y,z,...]$  where each of the *list* elements is an expression of a binary operand and 2 elements. For example  $x=1$ ,  $2^3$ ,  $[a,b]$  etc. The *key* is checked against the first operand. **assoc** returns the second operand if the *key* is found. If the *key* is not found it either returns the *default* value. *default* is optional and defaults to **false**.

**atom** (*expr*) Function

Returns **true** if *expr* is atomic (i.e. a number, name or string) else **false**. Thus **atom**(5) is **true** while **atom**( $a[1]$ ) and **atom**( $\sin(x)$ ) are **false** (assuming  $a[1]$  and  $x$  are unbound).

**cons** (*expr, list*) Function

Returns a new list constructed of the element *expr* as its first element, followed by the elements of *list*. **cons** also works on other expressions, e.g. **cons**( $x$ ,  $f(a,b,c)$ );  $\rightarrow f(x,a,b,c)$ .

**copylist** (*list*) Function

Returns a copy of the list *list*.

**delete** (*expr\_1*, *expr\_2*) Function

**delete** (*expr\_1*, *expr\_2*, *n*) Function

Removes all occurrences of *expr\_1* from *expr\_2*. *expr\_1* may be a term of *expr\_2* (if it is a sum) or a factor of *expr\_2* (if it is a product).

```
(%i1) delete(sin(x), x+sin(x)+y);
(%o1)                                     y + x
```

**delete**(*expr\_1*, *expr\_2*, *n*) removes the first *n* occurrences of *expr\_1* from *expr\_2*. If there are fewer than *n* occurrences of *expr\_1* in *expr\_2* then all occurrences will be deleted.

```
(%i1) delete(a, f(a,b,c,d,a));
(%o1)                                     f(b, c, d)
(%i2) delete(a, f(a,b,a,c,d,a), 2);
(%o2)                                     f(b, c, d, a)
```

**eighth** (*expr*) Function

Returns the 8'th item of expression or list *expr*. See **first** for more details.

**endcons** (*expr*, *list*) Function

Returns a new list consisting of the elements of *list* followed by *expr*. **endcons** also works on general expressions, e.g. **endcons**(*x*, *f*(*a*,*b*,*c*)); -> *f*(*a*,*b*,*c*,*x*).

**every** (*expr*) Function

This function takes a list, or a positive number of arguments and returns **true** if all its arguments are not **false**.

**fifth** (*expr*) Function

Returns the 5'th item of expression or list *expr*. See **first** for more details.

**first** (*expr*) Function

Returns the first part of *expr* which may result in the first element of a list, the first row of a matrix, the first term of a sum, etc. Note that **first** and its related functions, **rest** and **last**, work on the form of *expr* which is displayed not the form which is typed on input. If the variable **inflag** is set to **true** however, these functions will look at the internal form of *expr*. Note that the simplifier re-orders expressions. Thus **first**(*x*+*y*) will be *x* if **inflag** is **true** and *y* if **inflag** is **false** (**first**(*y*+*x*) gives the same results). The functions **second** .. **tenth** yield the second through the tenth part of their input argument.

**flatten** (*expr*) Function

Takes a list of the form [[1,2],[3,4]] and returns [1,2,3,4].

**fourth** (*expr*) Function

Returns the 4'th item of expression or list *expr*. See **first** for more details.

**get** (*a*, *i*) Function  
 Retrieves the user property indicated by *i* associated with atom *a* or returns **false** if *a* doesn't have property *i*.

```
(%i1) put (%e, 'transcendental, 'type);
(%o1)      transcendental
(%i2) put (%pi, 'transcendental, 'type)$
(%i3) put (%i, 'algebraic, 'type)$
(%i4) typeof (expr) := block ([q],
    if numberp (expr)
    then return ('algebraic),
    if not atom (expr)
    then return (maplist ('typeof, expr)),
    q: get (expr, 'type),
    if q=false
    then errcatch (error(expr,"is not numeric.)) else q)$
(%i5) typeof (2*%e + x*%pi);
x is not numeric.
(%o5)  [[transcendental, []], [algebraic, transcendental]]
(%i6) typeof (2*%e + %pi);
(%o6)  [transcendental, [algebraic, transcendental]]
```

**last** (*expr*) Function  
 Returns the last part (term, row, element, etc.) of the *expr*.

**length** (*expr*) Function  
 Returns (by default) the number of parts in the external (displayed) form of *expr*. For lists this is the number of elements, for matrices it is the number of rows, and for sums it is the number of terms (see `dispform`).

The `length` command is affected by the `inflag` switch. So, e.g. `length(a/(b*c))`; gives 2 if `inflag` is `false` (Assuming `exptdispflag` is `true`), but 3 if `inflag` is `true` (the internal representation is essentially  $a*b^{-1}*c^{-1}$ ).

**listarith** option variable  
 default value: `true` - if `false` causes any arithmetic operations with lists to be suppressed; when `true`, list-matrix operations are contagious causing lists to be converted to matrices yielding a result which is always a matrix. However, list-list operations should return lists.

**listp** (*expr*) Function  
 Returns `true` if *expr* is a list else `false`.

**makelist** (*expr*, *i*, *i.0*, *i.1*) Function

**makelist** (*expr*, *x*, *list*) Function

Constructs and returns a list, each element of which is generated from *expr*.

`makelist (expr, i, i.0, i.1)` returns a list, the *j*'th element of which is equal to `ev (expr, i=j)` for *j* equal to *i.0* through *i.1*.



`makelist (expr, x, list)` returns a list, the  $j$ 'th element of which is equal to `ev (expr, x=list[j])` for  $j$  equal to 1 through `length (list)`.

Examples:

```
(%i1) makelist(concat(x,i),i,1,6);
(%o1)          [x1, x2, x3, x4, x5, x6]
(%i2) makelist(x=y,y,[a,b,c]);
(%o2)          [x = a, x = b, x = c]
```

**member** (*expr*, *list*) Function

Returns `true` if *expr* occurs as a member of *list* (not within a member). Otherwise `false` is returned. `member` also works on non-list expressions, e.g. `member(b,f(a,b,c)); -> true`.

**ninth** (*expr*) Function

Returns the 9'th item of expression or list *expr*. See `first` for more details.

**rest** (*expr*, *n*) Function

**rest** (*expr*) Function

Returns *expr* with its first *n* elements removed if *n* is positive and its last - *n* elements removed if *n* is negative. If *n* is 1 it may be omitted. *expr* may be a list, matrix, or other expression.

**reverse** (*list*) Function

Reverses the order of the members of the *list* (not the members themselves). `reverse` also works on general expressions, e.g. `reverse(a=b);` gives `b=a`.

**second** (*expr*) Function

Returns the 2'nd item of expression or list *expr*. See `first` for more details.

**seventh** (*expr*) Function

Returns the 7'th item of expression or list *expr*. See `first` for more details.

**sixth** (*expr*) Function

Returns the 6'th item of expression or list *expr*. See `first` for more details.

**tenth** (*expr*) Function

Returns the 10'th item of expression or list *expr*. See `first` for more details.

**third** (*expr*) Function

Returns the 3'rd item of expression or list *expr*. See `first` for more details.

## 40 Function Definition

### 40.1 Introduction to Function Definition

### 40.2 Function

To define a function in Maxima you use the `:=` operator. E.g.

```
f(x) := sin(x)
```

defines a function `f`. Anonymous functions may also be created using `lambda`. For example

```
lambda ([i, j], ...)
```

can be used instead of `f` where

```
f(i,j) := block ([], ...);
map (lambda ([i], i+1), 1)
```

would return a list with 1 added to each term.

You may also define a function with a variable number of arguments, by having a final argument which is assigned to a list of the extra arguments:

```
(%i1) f ([u]) := u;
(%o1)          f([u]) := u
(%i2) f (1, 2, 3, 4);
(%o2)          [1, 2, 3, 4]
(%i3) f (a, b, [u]) := [a, b, u];
(%o3)          f(a, b, [u]) := [a, b, u]
(%i4) f (1, 2, 3, 4, 5, 6);
(%o4)          [1, 2, [3, 4, 5, 6]]
```

The right hand side of a function is an expression. Thus if you want a sequence of expressions, you do

```
f(x) := (expr1, expr2, ..., exprn);
```

and the value of `exprn` is what is returned by the function.

If you wish to make a `return` from some expression inside the function then you must use `block` and `return`.

```
block ([], expr1, ..., if (a > 10) then return(a), ..., exprn)
```

is itself an expression, and so could take the place of the right hand side of a function definition. Here it may happen that the return happens earlier than the last expression.

The first `[]` in the block, may contain a list of variables and variable assignments, such as `[a: 3, b, c: []]`, which would cause the three variables `a`, `b`, and `c` to not refer to their global values, but rather have these special values for as long as the code executes inside the `block`, or inside functions called from inside the `block`. This is called *dynamic* binding, since the variables last from the start of the block to the time it exits. Once you return from the `block`, or throw out of it, the old values (if any) of the variables will be restored. It is certainly a good idea to protect your variables in this way. Note that the assignments in the block variables, are done in parallel. This means, that if you had used `c: a` in the above, the value of `c` would have been the value of `a` at the time you just entered the block, but before `a` was bound. Thus doing something like

```
block ([a: a], expr1, ... a: a+3, ..., exprn)
```

will protect the external value of `a` from being altered, but would let you access what that value was. Thus the right hand side of the assignments, is evaluated in the entering context, before any binding occurs. Using just `block ([x], ...` would cause the `x` to have itself as value, just as if it would have if you entered a fresh **Maxima** session.

The actual arguments to a function are treated in exactly same way as the variables in a block. Thus in

```
f(x) := (expr1, ..., exprn);
and
f(1);
```

we would have a similar context for evaluation of the expressions as if we had done

```
block ([x: 1], expr1, ..., exprn)
```

Inside functions, when the right hand side of a definition, may be computed at runtime, it is useful to use `define` and possibly `buildq`.

## 40.3 Macros

**buildq** (*variables*, *expr*)

Function

*expr* is any single Maxima expression and *variables* is a list of elements of the form `<atom>` or `<atom>: <value>`.

### 40.3.1 Semantics

The elements of the list *variables* are evaluated left to right (the syntax *atom* is equivalent to *atom: atom*). then these values are substituted into `<expression>` in parallel. If any *atom* appears as a single argument to the special form `splice` (i.e. `splice (atom)`) inside *expr*, then the value associated with that *atom* must be a Maxima list, and it is spliced into *expr* instead of substituted.

### 40.3.2 Simplification

The arguments to `buildq` need to be protected from simplification until the substitutions have been carried out. This code should effect that by using `'`.

`buildq` can be useful for building functions on the fly. One of the powerful things about **Maxima** is that you can have your functions define other functions to help solve the problem. Further below we discuss building a recursive function, for a series solution. This defining of functions inside functions usually uses `define`, which evaluates its arguments. A number of examples are included under `splice`.

**splice** (*atom*)

Function

This is used with `buildq` to construct a list. This is handy for making argument lists, in conjunction with `buildq`.

```
mprint ([x]) ::= buildq ([u : x],
  if (debuglevel > 3) print (splice (u)));
```

Including a call like

```
mprint ("matrix is", mat, "with length", length(mat))
```

is equivalent to putting in the line

```
if (debuglevel > 3) print ("matrix is", mat, "with length", length(mat));
```

A more non trivial example would try to display the variable values and their names.

```
mshow (a, b, c)
```

should become

```
print ('a, "=", a, ",", 'b, "=", b, ", and", 'c, "=", c)
```

so that if it occurs as a line in a program we can print values.

```
(%i1) foo (x,y,z) := mshow (x, y, z);
(%i2) foo (1, 2, 3);
x = 1 , y = 2 , and z = 3
```

The actual definition of mshow is the following. Note how buildq lets you build "quoted" structure, so that the 'u lets you get the variable name. Note that in macros, the result is a piece of code which will then be substituted for the macro and evaluated.

```
mshow ([l]) ::= block ([ans:[], n:length(l)],
  for i:1 thru n do
    (ans: append (ans, buildq ([u: l[i]], ['u, "=", u])),
    if i < n then
      ans: append (ans, if i < n-1 then [","] else [", and"])),
  buildq ([u:ans], print (splice(u))));
```

The splice also works to put arguments into algebraic operations:

```
(%i1) buildq ([a: '[b, c, d]], +splice(a));
(%o1)          d + c + b
```

Note how the simplification only occurs *after* the substitution, The operation applying to the splice in the first case is the + while in the second it is the \*, yet logically you might think splice(a)+splice(a) could be replaced by 2\*splice(a). No simplification takes place with the buildq. To understand what splice is doing with the algebra you must understand that for Maxima, a formula an operation like a+b+c is really internally similar to +(a,b,c), and similarly for multiplication. Thus \*(2,b,c,d) is 2\*b\*c\*d.

```
(%i1) buildq ([a: '[b,c,d]], +splice(a));
(%o1)          d + c + b
(%i2) buildq ([a: '[b,c,d]], splice(a)+splice(a));
(%o2)          2 d + 2 c + 2 b
```

but

```
(%i3) buildq ([a: '[b,c,d]], 2*splice(a));
(%o3)          2 b c d
```

Finally buildq can be invaluable for building recursive functions. Suppose your program is solving a differential equation using the series method, and has determined that it needs to build a recursion relation

$$f[n] := -((n^2 - 2*n + 1)*f[n-1] + f[n-2] + f[n-3])/(n^2-n)$$

and it must do this on the fly inside your function. Now you would really like to add expand.

```
f[n] := expand (-((n^2 - 2*n + 1)*f[n-1] + f[n-2] + f[n-3])/(n^2-n))
```

but how do you build this code. You want the `expand` to happen each time the function runs, *not* before it.

```
(%i1) val: -((n^2 - 2*n + 1)*f[n-1] + f[n-2] + f[n-3])/(n^2-n)$
(%i2) define (f[n], buildq ([u: val], expand(u)))$
```

does the job. This might be useful, since when you do (with `expand`)

```
(%i3) f[0]: aa0$
(%i4) f[1]: aa1$
(%i5) f[2]: aa2$
(%i6) f[6];
```

```
(%o6)
          3 aa2   aa1   7 aa0
        ----- + --- + -----
          10     40     90
```

where as without it is kept unsimplified, and even after 6 terms it becomes:

```
(%i7) define (g[n], buildq ([u: val], u))$
(%i8) g[0]: bb0$
(%i9) g[1]: bb1$
(%i10) g[2]: bb2$
(%i11) g[6];
```

```
          aa2          7 aa2   aa1   11 aa0   aa1   aa0
        --- - 25 (- ----- - --- - -----) + --- + ---
          4           20     40     120     8     24
(%o11) -----
                               30
```

```
(%i12) expand (%);
```

```
(%o12)
          3 aa2   aa1   7 aa0
        ----- + --- + -----
          10     40     90
```

The expression quickly becomes complicated if not simplified at each stage, so the simplification must be part of the definition. Hence the `buildq` is useful for building the form.

## 40.4 Definitions for Function Definition

**apply** ( $f$ , [ $x_1$ , ...,  $x_n$ ])

Function

Returns the result of applying the function  $f$  to the list of arguments  $x_1, \dots, x_n$ .  $f$  is the name of a function or a lambda expression.

This is useful when it is desired to compute the arguments to a function before applying that function. For example, if `l` is the list `[1, 5, -10.2, 4, 3]`, then `apply (min, l)` gives `-10.2`. `apply` is also useful when calling functions which do not have their arguments evaluated if it is desired to cause evaluation of them. For example, if `filespec` is a variable bound to the list `[test, case]` then `apply (closefile, filespec)` is equivalent to `closefile (test, case)`. In general the first argument to `apply` should be preceded by a `'` to make it evaluate to itself. Since some atomic variables have the same name as certain functions the values of the variable would be

used rather than the function because `apply` has its first argument evaluated as well as its second.

**block** ( $[v_1, \dots, v_m], expr_1, \dots, expr_n$ ) Function

**block** ( $expr_1, \dots, expr_n$ ) Function

`block` evaluates  $expr_1, \dots, expr_n$  in sequence and returns the value of the last expression evaluated. The sequence can be modified by the `go`, `throw`, and `return` functions. The last expression is  $expr_n$  unless `return` or an expression containing `throw` is evaluated. Some variables  $v_1, \dots, v_m$  can be declared local to the block; these are distinguished from global variables of the same names. If no variables are declared local then the list may be omitted. Within the block, any variable other than  $v_1, \dots, v_m$  is a global variable.

`block` saves the current values of the variables  $v_1, \dots, v_m$  (if any) upon entry to the block, then unbinds the variables so that they evaluate to themselves. The local variables may be bound to arbitrary values within the block but when the block is exited the saved values are restored, and the values assigned within the block are lost.

`block` may appear within another `block`. Local variables are established each time a new `block` is evaluated. Local variables appear to be global to any enclosed blocks. If a variable is non-local in a block, its value is the value most recently assigned by an enclosing block, if any, otherwise, it is the value of the variable in the global environment. This policy may coincide with the usual understanding of "dynamic scope".

If it is desired to save and restore other local properties besides `value`, for example `array` (except for complete arrays), `function`, `dependencies`, `atvalue`, `matchdeclare`, `atomgrad`, `constant`, and `nonscalar` then the function `local` should be used inside of the block with arguments being the names of the variables.

The value of the block is the value of the last statement or the value of the argument to the function `return` which may be used to exit explicitly from the block. The function `go` may be used to transfer control to the statement of the block that is tagged with the argument to `go`. To tag a statement, precede it by an atomic argument as another statement in the block. For example: `block ([x], x:1, loop, x: x+1, ..., go(loop), ...)`. The argument to `go` must be the name of a tag appearing within the block. One cannot use `go` to transfer to a tag in a block other than the one containing the `go`.

Blocks typically appear on the right side of a function definition but can be used in other places as well.

**break** ( $expr_1, \dots, expr_n$ ) Function

Evaluates and prints  $expr_1, \dots, expr_n$  and then causes a Maxima break at which point the user can examine and change his environment. Upon typing `exit`; the computation resumes.

**catch** ( $expr_1, \dots, expr_n$ ) Function

Evaluates  $expr_1, \dots, expr_n$  one by one; if any leads to the evaluation of an expression of the form `throw (arg)`, then the value of the `catch` is the value of `throw (arg)`,

and no further expressions are evaluated. This "non-local return" thus goes through any depth of nesting to the nearest enclosing `catch`. If there is no `catch` enclosing a `throw`, an error message is printed.

If the evaluation of the arguments does not lead to the evaluation of any `throw` then the value of `catch` is the value of `expr_n`.

```
(%i1) lambda ([x], if x < 0 then throw(x) else f(x))$
(%i2) g(l) := catch (map ('%, l))$
(%i3) g ([1, 2, 3, 7]);
(%o3)          [f(1), f(2), f(3), f(7)]
(%i4) g ([1, 2, -3, 7]);
(%o4)          - 3
```

The function `g` returns a list of `f` of each element of `l` if `l` consists only of non-negative numbers; otherwise, `g` "catches" the first negative element of `l` and "throws" it up.

**compile** (*filename*, *f\_1*, ..., *f\_n*) Function

Translates Maxima functions *f\_1*, ..., *f\_n* into Lisp and writes the translated code into the file *filename*.

The Lisp translations are not evaluated, nor is the output file processed by the Lisp compiler. `translate` creates and evaluates Lisp translations. `compile_file` translates Maxima into Lisp, and then executes the Lisp compiler.

See also `translate`, `translate_file`, and `compile_file`.

**compile** (*f\_1*, ..., *f\_n*) Function

**compile** (*functions*) Function

**compile** (*all*) Function

Translates Maxima functions *f\_1*, ..., *f\_n* into Lisp, evaluates the Lisp translations, and calls the Lisp function `COMPILE` on each translated function. `compile` returns a list of the names of the compiled functions.

`compile (all)` or `compile (functions)` compiles all user-defined functions.

`compile` quotes its arguments; the double-single-quotes operator `''` defeats quotation.

**define** (*f(x\_1, ..., x\_n)*, *expr*) Function

Defines a function named *f* with arguments *x\_1*, ..., *x\_n* and function body *expr*.

`define` is similar to the function definition operator `:=`, but when `define` appears inside a function, the definition is created using the value of `expr` at execution time rather than at the time of definition of the function which contains it.

All function definitions appear in the same namespace; defining a function `f` within another function `g` does not limit the scope of `f` to `g`.

Examples:

```
(%i1) foo: 2^bar; bar (%o1) 2 (%i2) g(x) := (f_1 (y) := foo*x*y, f_2 (y) := "foo*x*y,
define (f_3 (y), foo*x*y), define (f_4 (y), "foo*x*y)); bar (%o2) g(x) := (f_1(y) := foo
x y, f_2(y) := 2 x y, bar define(f_3(y), foo x y), define(f_4(y), 2 x y)) (%i3) functions;
(%o3) [g(x)] (%i4) g(a); bar (%o4) f_4(y) := a 2 y (%i5) functions; (%o5) [g(x), f_1(y),
f_2(y), f_3(y), f_4(y)] (%i6) dispfun (f_1, f_2, f_3, f_4); (%t6) f_1(y) := foo x y
```

```

bar (%t7) f.2(y) := 2 x y
bar (%t8) f.3(y) := a 2 y
bar (%t9) f.4(y) := a 2 y
(%o9) done

```

**define\_variable** (*name, default\_value, mode, documentation*) Function

Introduces a global variable into the Maxima environment. This is for user-written packages, which are often translated or compiled. Thus

```
define_variable (foo, true, boolean);
```

does the following:

- (1) `mode_declare (foo, boolean)` sets it up for the translator.
- (2) If the variable is unbound, it sets it: `foo: true`.
- (3) `declare (foo, special)` declares it special.
- (4) Sets up an assign property for it to make sure that it never gets set to a value of the wrong mode. E.g. `foo: 44` would be an error once `foo` is defined `boolean`.

See `mode_declare` for a list of the possible modes.

The optional 4th argument is a documentation string. When `translate_file` is used on a package which includes documentation strings, a second file is output in addition to the Lisp file which will contain the documentation strings, formatted suitably for use in manuals, usage files, or (for instance) `describe`.

With any variable which has been `define_variable`'d with mode other than `any`, you can give a `value_check` property, which is a function of one argument called on the value the user is trying to set the variable to.

```
put ('g5, lambda([u], if u # 'g5 then error("Don't set g5")), 'value_check);
```

Use `define_variable (g5, 'g5, any_check, "this ain't supposed to be set by anyone but me.")` `any_check` is a mode which means the same as `any`, but which keeps `define_variable` from optimizing away the assign property.

**dispfun** (*f.1, ..., f.n*) Function

**dispfun** (*all*) Function

Displays the definition of the user-defined functions *f.1, ..., f.n*. Each argument may be the name of a macro (defined with `:=`), an ordinary function (defined with `:=` or `define`), an array function (defined with `:=` or `define`, but enclosing arguments in square brackets [ ]), a subscripted function, (defined with `:=` or `define`, but enclosing some arguments in square brackets and others in parentheses ( )) one of a family of subscripted functions selected by a particular subscript value, or a subscripted function defined with a constant subscript.

`dispfun (all)` displays all user-defined functions as given by the `functions`, `arrays`, and `macros` lists, omitting subscripted functions defined with constant subscripts.

`dispfun` creates an intermediate expression label (`%t1, %t2`, etc.) for each displayed function, and assigns the function definition to the label. In contrast, `fundef` returns the function definition.



`dispfun` quotes its arguments; the double-single-quote operator `''` defeats quotation. `dispfun` always returns `done`.

Examples:

```
(%i1) m(x, y) ::= x^(-y)$
(%i2) f(x, y) := x^(-y)$
(%i3) g[x, y] := x^(-y)$
(%i4) h[x](y) := x^(-y)$
(%i5) i[8](y) := 8^(-y)$
(%i6) dispfun (m, f, g, h, h[5], h[10], i[8])$
```

```
(%t6)          m(x, y) ::= x-y
```

```
(%t7)          f(x, y) := x-y
```

```
(%t8)          gx, y := x-y
```

```
(%t9)          hx(y) := x-y
```

```
(%t10)         h5(y) :=  $\frac{1}{y^5}$ 
```

```
(%t11)         h10(y) :=  $\frac{1}{y^{10}}$ 
```

```
(%t12)         i8(y) := 8-y
```

## functions

Variable

Default value: []

`functions` is the list of user-defined Maxima functions in the current session. A user-defined function is a function constructed by `define` or `:=`. A function may be defined at the Maxima prompt or in a Maxima file loaded by `load` or `batch`. Lisp functions, however, are not added to `functions`.

## fundef (f)

Function

Returns the definition of the function *f*.

The argument may be the name of a macro (defined with `:=`), an ordinary function (defined with `:=` or `define`), an array function (defined with `:=` or `define`, but

enclosing arguments in square brackets [ ]), a subscripted function, (defined with := or **define**, but enclosing some arguments in square brackets and others in parentheses ( )) one of a family of subscripted functions selected by a particular subscript value, or a subscripted function defined with a constant subscript.

**fundef** quotes its argument; the double-single-quote operator '' defeats quotation.

**fundef** (*f*) returns the definition of *f*. In contrast, **dispfun** (*f*) creates an intermediate expression label and assigns the definition to the label.

**funmake** (*name*, [*arg\_1*, ..., *arg\_n*]) Function

Returns an expression *name* (*arg\_1*, ..., *arg\_n*). The return value is simplified, but not evaluated, so the function is not called.

**funmake** evaluates its arguments.

Examples:

- **funmake** evaluates its arguments, but not the return value.

```
(%i1) det(a,b,c) := b^2 -4*a*c$
(%i2) x: 8$
(%i3) y: 10$
(%i4) z: 12$
(%i5) f: det$
(%i6) funmake (f, [x, y, z]);
(%o6)          det(8, 10, 12)
(%i7) ''%;
(%o7)          - 284
```

- Maxima simplifies **funmake**'s return value.

```
(%i1) funmake (sin, [%pi/2]);
(%o1)          1
```

**lambda** [*x\_1*, ..., *x\_m*], *expr\_1*, ..., *expr\_n*) Function

Defines and returns a lambda expression (that is, an anonymous function) with arguments *x\_1*, ..., *x\_m* and return value *expr\_n*. A lambda expression can be assigned to a variable and evaluated like an ordinary function. A lambda expression may appear in contexts in which a function evaluation (but not a function name) is expected.

When the function is evaluated, unbound local variables *x\_1*, ..., *x\_m* are created. **lambda** may appear within **block** or another **lambda**; local variables are established each time another **block** or **lambda** is evaluated. Local variables appear to be global to any enclosed **block** or **lambda**. If a variable is not local, its value is the value most recently assigned in an enclosing **block** or **lambda**, if any, otherwise, it is the value of the variable in the global environment. This policy may coincide with the usual understanding of "dynamic scope".

After local variables are established, *expr\_1* through *expr\_n* are evaluated in turn. The special variable %, representing the value of the preceding expression, is recognized. **throw** and **catch** may also appear in the list of expressions.

**return** cannot appear in a lambda expression unless enclosed by **block**, in which case **return** defines the return value of the block and not of the lambda expression, unless

the block happens to be *expr.n*. Likewise, *go* cannot appear in a lambda expression unless enclosed by *block*.

*lambda* quotes its arguments; the double-single-quote operator `''` defeats quotation.

Examples:

- A lambda expression can be assigned to a variable and evaluated like an ordinary function.

```
(%i1) f: lambda ([x], x^2);
(%o1)          lambda([x], x2)
(%i2) f(a);
(%o2)          a2
```

- A lambda expression may appear in contexts in which a function evaluation is expected.

```
(%i3) lambda ([x], x^2) (a);
(%o3)          a2
(%i4) apply (lambda ([x], x^2), [a]);
(%o4)          a2
(%i5) map (lambda ([x], x^2), [a, b, c, d, e]);
(%o5)          [a2, b2, c2, d2, e2]
```

- Argument variables are local variables. Other variables appear to be global variables. Global variables are evaluated at the time the lambda expression is evaluated, unless some special evaluation is forced by some means, such as `''`.

```
(%i6) a: %pi$
(%i7) b: %e$
(%i8) g: lambda ([a], a*b);
(%o8)          lambda([a], a b)
(%i9) b: %gamma$
(%i10) g(1/2);
(%o10)          %gamma
                -----
                2
(%i11) g2: lambda ([a], a*''b);
(%o11)          lambda([a], a %gamma)
(%i12) b: %e$
(%i13) g2(1/2);
(%o13)          %gamma
                -----
                2
```

- Lambda expressions may be nested. Local variables within the outer lambda expression appear to be global to the inner expression unless masked by local variables of the same names.

```
(%i14) h: lambda ([a, b], h2: lambda ([a], a*b), h2(1/2));
```

```
(%o14)      lambda([a, b], h2 : lambda([a], a b), h2(-))
                                                    1
                                                    2
```

```
(%i15) h(%pi, %gamma);
```

```
(%o15)
           %gamma
      -----
           2
```

- Since `lambda` quotes its arguments, `lambda` expression `i` below does not define a "multiply by `a`" function. Such a function can be defined via `buildq`, as in `lambda` expression `i2` below.

```
(%i16) i: lambda ([a], lambda ([x], a*x));
```

```
(%o16)      lambda([a], lambda([x], a x))
```

```
(%i17) i(1/2);
```

```
(%o17)
           lambda([x], a x)
```

```
(%i18) i2: lambda([a], buildq([a: a], lambda([x], a*x)));
```

```
(%o18)      lambda([a], buildq([a : a], lambda([x], a x)))
```

```
(%i19) i2(1/2);
```

```
(%o19)
           x
      lambda([x], -)
           2
```

```
(%i20) i2(1/2)(%pi);
```

```
(%o20)
           %pi
      ----
           2
```

### **local** ( $v_1, \dots, v_n$ )

Function

Declares the variables  $v_1, \dots, v_n$  to be local with respect to all the properties in the statement in which this function is used.

`local` may only be used in `block`, in the body of function definitions or `lambda` expressions, or in the `ev` function, and only one occurrence is permitted in each.

`local` is independent of `context`.

### **macroexpansion**

Variable

Default value: `false`

`macroexpansion` controls advanced features which affect the efficiency of macros.

Possible settings:

- `false` – Macros expand normally each time they are called.
- `expand` – The first time a particular call is evaluated, the expansion is remembered internally, so that it doesn't have to be recomputed on subsequent calls making subsequent calls faster. The macro call still calls `grind` and `display` normally. However, extra memory is required to remember all of the expansions.
- `displace` – The first time a particular call is evaluated, the expansion is substituted for the call. This requires slightly less storage than when `macroexpansion` is set to `expand` and is just as fast, but has the disadvantage that the original macro call is no longer remembered and hence the expansion will be seen if `display` or `grind` is called. See documentation for `translate` and `macros` for more details.

**mode\_checkp** Variable

Default value: `true`

When `mode_checkp` is `true`, `mode_declare` checks the modes of bound variables.

**mode\_check\_errorp** Variable

Default value: `false`

When `mode_check_errorp` is `true`, `mode_declare` calls error.

**mode\_check\_warnp** Variable

Default value: `true`

When `mode_check_warnp` is `true`, mode errors are described.

**mode\_declare** (*y<sub>1</sub>, mode<sub>1</sub>, ..., y<sub>n</sub>, mode<sub>n</sub>*) Function

`mode_declare` is used to declare the modes of variables and functions for subsequent translation or compilation of functions. `mode_declare` is typically placed at the beginning of a function definition, at the beginning of a Maxima script, or executed at the interactive prompt.

The arguments of `mode_declare` are pairs consisting of a variable and a mode which is one of `boolean`, `fixnum`, `number`, `rational`, or `float`. Each variable may also be a list of variables all of which are declared to have the same mode.

If a variable is an array, and if every element of the array which is referenced has a value then `array (yi, complete, dim1, dim2, ...)` rather than

```
array(yi, dim1, dim2, ...)
```

should be used when first declaring the bounds of the array. If all the elements of the array are of mode `fixnum` (`float`), use `fixnum` (`float`) instead of `complete`. Also if every element of the array is of the same mode, say `m`, then

```
mode_declare (completearray (yi), m)
```

should be used for efficient translation.

Numeric code using arrays might run faster by declaring the expected size of the array, as in:

```
mode_declare (completearray (a [10, 10]), float)
```

for a floating point number array which is 10 x 10.

One may declare the mode of the result of a function by using `function (f1, f2, ...)` as an argument; here `f1, f2, ...` are the names of functions. For example the expression,

```
mode_declare ([function (f1, f2, ...)], fixnum)
```

declares that the values returned by `f1, f2, ...` are single-word integers.

`modedeclare` is a synonym for `mode_declare`.

**mode\_identity** (*arg<sub>1</sub>, arg<sub>2</sub>*) Function

A special form used with `mode_declare` and `macros` to declare, e.g., a list of lists of flonums, or other compound data object. The first argument to `mode_identity` is a primitive value mode name as given to `mode_declare` (i.e., one of `float`, `fixnum`,

number, list, or any), and the second argument is an expression which is evaluated and returned as the value of `mode_identity`. However, if the return value is not allowed by the mode declared in the first argument, an error or warning is signalled. The important thing is that the mode of the expression as determined by the Maxima to Lisp translator, will be that given as the first argument, independent of anything that goes on in the second argument. E.g., `x: 3.3; mode_identity (fixnum, x);` yields an error. `mode_identity (flonum, x)` returns 3.3. This has a number of uses, e.g., if you knew that `first (l)` returned a number then you might write `mode_identity (number, first (l))`. However, a more efficient way to do it would be to define a new primitive,

```
firstnumb (x) ::= buildq ([x], mode_identity (number, x));
```

and use `firstnumb` every time you take the first of a list of numbers.

### **transcompile**

Variable

Default value: `true`

When `transcompile` is `true`, `translate` and `translate_file` generate declarations to make the translated code more suitable for compilation.

`compile` sets `transcompile: true` for the duration.

### **translate** (*f<sub>1</sub>, ..., f<sub>n</sub>*)

Function

### **translate** (*functions*)

Function

### **translate** (*all*)

Function

Translates the user-defined functions *f<sub>1</sub>, ..., f<sub>n</sub>* from the Maxima language into Lisp and evaluates the Lisp translations. Typically the translated functions run faster than the originals.

`translate (all)` or `translate (functions)` translates all user-defined functions.

Functions to be translated should include a call to `mode_declare` at the beginning when possible in order to produce more efficient code. For example:

```
f (x_1, x_2, ...) := block ([v_1, v_2, ...],
    mode_declare (v_1, mode_1, v_2, mode_2, ...), ...)
```

where the *x<sub>1</sub>, x<sub>2</sub>, ...* are the parameters to the function and the *v<sub>1</sub>, v<sub>2</sub>, ...* are the local variables.

The names of translated functions are removed from the `functions` list if `savedef` is `false` (see below) and are added to the `props` lists.

Functions should not be translated unless they are fully debugged.

Expressions are assumed simplified; if they are not, correct but non-optimal code gets generated. Thus, the user should not set the `simp` switch to `false` which inhibits simplification of the expressions to be translated.

The switch `translate`, if `true`, causes automatic translation of a user's function to Lisp.

Note that translated functions may not run identically to the way they did before translation as certain incompatibilities may exist between the Lisp and Maxima versions. Principally, the `rat` function with more than one argument and the `ratvars`

function should not be used if any variables are `mode_declare`'d canonical rational expressions (CRE). Also the `prederror: false` setting will not translate.

`savedef` - if `true` will cause the Maxima version of a user function to remain when the function is `translate`'d. This permits the definition to be displayed by `dispfun` and allows the function to be edited.

`transrun` - if `false` will cause the interpreted version of all functions to be run (provided they are still around) rather than the translated version.

The result returned by `translate` is a list of the names of the functions translated.

**translate\_file** (*maxima\_filename*) Function

**translate\_file** (*maxima\_filename, lisp\_filename*) Function

Translates a file of Maxima code into a file of Lisp code. `translate_file` returns a list of three filenames: the name of the Maxima file, the name of the Lisp file, and the name of file containing additional information about the translation. `translate_file` evaluates its arguments.

`translate_file ("foo.mac"); load("foo.LISP")` is the same as `batch ("foo.mac")` except for certain restrictions, the use of `'` and `%`, for example.

`translate_file (maxima_filename)` translates a Maxima file *maxima\_filename* into a similarly-named Lisp file. For example, `foo.mac` is translated into `foo.LISP`. The Maxima filename may include a directory name or names, in which case the Lisp output file is written to the same directory from which the Maxima input comes.

`translate_file (maxima_filename, lisp_filename)` translates a Maxima file *maxima\_filename* into a Lisp file *lisp\_filename*. `translate_file` ignores the filename extension, if any, of *lisp\_filename*; the filename extension of the Lisp output file is always `LISP`. The Lisp filename may include a directory name or names, in which case the Lisp output file is written to the specified directory.

`translate_file` also writes a file of translator warning messages of various degrees of severity. The filename extension of this file is `UNLISP`. This file may contain valuable information, though possibly obscure, for tracking down bugs in translated code. The `UNLISP` file is always written to the same directory from which the Maxima input comes.

`translate_file` emits Lisp code which causes some declarations and definitions to take effect as soon as the Lisp code is compiled. See `compile_file` for more on this topic.

See also `tr_array_as_ref`, `tr_bound_function_apply`, `tr_exponent`, `tr_file_tty_messagesp`, `tr_float_can_branch_complex`, `tr_function_call_default`, `tr_numer`, `tr_optimize_max_loop`, `tr_semicompile`, `tr_state_vars`, `tr_warnings_get`, `tr_warn_bad_function_calls`, `tr_warn_fexpr`, `tr_warn_meval`, `tr_warn_mode`, `tr_warn_undeclared`, `tr_warn_undefined_variable`, and `tr_windy`.

**transrun** Variable

Default value: `true`

When `transrun` is `false` will cause the interpreted version of all functions to be run (provided they are still around) rather than the translated version.

**tr\_array\_as\_ref** Variable

Default value: `true`

If `translate_fast_arrays` is `false`, array references in Lisp code emitted by `translate_file` are affected by `tr_array_as_ref`. When `tr_array_as_ref` is `true`, array names are evaluated, otherwise array names appear as literal symbols in translated code.

`tr_array_as_ref` has no effect if `translate_fast_arrays` is `true`.

**tr\_bound\_function\_applyp** Variable

Default value: `true`

When `tr_bound_function_applyp` is `true`, Maxima gives a warning if a bound variable (such as a function argument) is found being used as a function. `tr_bound_function_applyp` does not affect the code generated in such cases.

For example, an expression such as `g(f, x) := f(x+1)` will trigger the warning message.

**tr\_file\_tty\_messagesp** Variable

Default value: `false`

When `tr_file_tty_messagesp` is `true`, messages generated by `translate_file` during translation of a file are displayed on the console and inserted into the UNLISP file. When `false`, messages about translation of the file are only inserted into the UNLISP file.

**tr\_float\_can\_branch\_complex** Variable

Default value: `true`

Tells the Maxima-to-Lisp translator to assume that the functions `acos`, `asin`, `asec`, and `acsc` can return complex results.

The ostensible effect of `tr_float_can_branch_complex` is the following. However, it appears that this flag has no effect on the translator output.

When it is `true` then `acos(x)` is of mode `any` even if `x` is of mode `float` (as set by `mode_declare`). When `false` then `acos(x)` is of mode `float` if and only if `x` is of mode `float`.

**tr\_function\_call\_default** Variable

Default value: `general`

`false` means give up and call `meval`, `expr` means assume Lisp fixed arg function. `general`, the default gives code good for `mexprs` and `mlexprs` but not `macros`. `general` assures variable bindings are correct in compiled code. In `general` mode, when translating `F(X)`, if `F` is a bound variable, then it assumes that `apply(f, [x])` is meant, and translates a `such`, with appropriate warning. There is no need to turn this off. With the default settings, no warning messages implies full compatibility of translated and compiled code with the Maxima interpreter.



- tr\_numer** Variable  
 Default value: `false`  
 When `tr_numer` is `true` numer properties are used for atoms which have them, e.g. `%pi`.
- tr\_optimize\_max\_loop** Variable  
 Default value: `100`  
`tr_optimize_max_loop` is the maximum number of times the macro-expansion and optimization pass of the translator will loop in considering a form. This is to catch macro expansion errors, and non-terminating optimization properties.
- tr\_semicompile** Variable  
 Default value: `false`  
 When `tr_semicompile` is `true`, `translate_file` and `compfile` output forms which will be macroexpanded but not compiled into machine code by the Lisp compiler.
- tr\_state\_vars** Variable  
 Default value:  

```
[transcompile, tr_semicompile, tr_warn_undeclared, tr_warn_meval,
tr_warn_fexpr, tr_warn_mode, tr_warn_undefined_variable,
tr_function_call_default, tr_array_as_ref, tr_numer]
```

 The list of the switches that affect the form of the translated output. This information is useful to system people when trying to debug the translator. By comparing the translated product to what should have been produced for a given state, it is possible to track down bugs.
- tr\_warnings\_get ()** Function  
 Prints a list of warnings which have been given by the translator during the current translation.
- tr\_warn\_bad\_function\_calls** Variable  
 Default value: `true`  
 - Gives a warning when when function calls are being made which may not be correct due to improper declarations that were made at translate time.
- tr\_warn\_fexpr** Variable  
 Default value: `compfile`  
 - Gives a warning if any FEXPRs are encountered. FEXPRs should not normally be output in translated code, all legitimate special program forms are translated.
- tr\_warn\_meval** Variable  
 Default value: `compfile`  
 - Gives a warning if the function `meval` gets called. If `meval` is called that indicates problems in the translation.

- tr\_warn\_mode** Variable  
 Default value: `all`  
 - Gives a warning when variables are assigned values inappropriate for their mode.
- tr\_warn\_undeclared** Variable  
 Default value: `compile`  
 - Determines when to send warnings about undeclared variables to the TTY.
- tr\_warn\_undefined\_variable** Variable  
 Default value: `all`  
 - Gives a warning when undefined global variables are seen.
- tr\_windy** Variable  
 Default value: `true`  
 - Generate "helpfull" comments and programming hints.
- compile\_file** (*filename*) Function  
**compile\_file** (*filename*, *compiled\_filename*) Function  
**compile\_file** (*filename*, *compiled\_filename*, *lisp\_filename*) Function
- Translates the Maxima file *filename* into Lisp, executes the Lisp compiler, and, if the translation and compilation succeed, loads the compiled code into Maxima.
- `compile_file` returns a list of the names of four files: the original Maxima file, the Lisp translation, notes on translation, and the compiled code. If the compilation fails, the fourth item is `false`.
- Some declarations and definitions take effect as soon as the Lisp code is compiled (without loading the compiled code). These include functions defined with the `:=` operator, macros define with the `::=` operator, `alias`, `declare`, `define_variable`, `mode_declare`, and `infix`, `matchfix`, `nofix`, `postfix`, `prefix`, and `compile`.
- Assignments and function calls are not evaluated until the compiled code is loaded. In particular, within the Maxima file, assignments to the translation flags (`tr_numer`, etc.) have no effect on the translation.
- filename* may not contain `:lisp` statements.
- `compile_file` evaluates its arguments.
- declare\_translated** (*f.1*, *f.2*, ...) Function
- When translating a file of Maxima code to Lisp, it is important for the translator to know which functions it sees in the file are to be called as translated or compiled functions, and which ones are just Maxima functions or undefined. Putting this declaration at the top of the file, lets it know that although a symbol does which does not yet have a Lisp function value, will have one at call time. (`MFUNCTION-CALL fn arg1 arg2 ...`) is generated when the translator does not know `fn` is going to be a Lisp function.



## 41 Program Flow

### 41.1 Introduction to Program Flow

Maxima provides a `do` loop for iteration, as well as more primitive constructs such as `go`.

### 41.2 Definitions for Program Flow

**backtrace** () Function  
**backtrace** (*n*) Function

Prints the call stack, that is, the list of functions which called the currently active function.

`backtrace()` prints the entire call stack.

`backtrace(n)` prints the *n* most recent functions, including the currently active function.

`backtrace` can be called from a script, a function, or the interactive prompt (not only in a debugging context).

- `backtrace()` prints the entire call stack.

```
(%i1) h(x) := g(x/7)$
(%i2) g(x) := f(x-11)$
(%i3) f(x) := e(x^2)$
(%i4) e(x) := (backtrace(), 2*x + 13)$
(%i5) h(10);
#0: e(x=4489/49)
#1: f(x=-67/7)
#2: g(x=10/7)
#3: h(x=10)
                                     9615
(%o5)                                ----
                                     49
```

- `backtrace(n)` prints the *n* most recent functions, including the currently active function.

```
(%i1) h(x) := (backtrace(1), g(x/7))$
(%i2) g(x) := (backtrace(1), f(x-11))$
(%i3) f(x) := (backtrace(1), e(x^2))$
(%i4) e(x) := (backtrace(1), 2*x + 13)$
(%i5) h(10);
#0: h(x=10)
#0: g(x=10/7)
#0: f(x=-67/7)
#0: e(x=4489/49)
                                     9615
(%o5)                                ----
                                     49
```

**do**

special operator

The **do** statement is used for performing iteration. Due to its great generality the **do** statement will be described in two parts. First the usual form will be given which is analogous to that used in several other programming languages (Fortran, Algol, PL/I, etc.); then the other features will be mentioned.

There are three variants of this form that differ only in their terminating conditions. They are:

- **for** *variable*: *initial\_value* **step** *increment* **thru** *limit* **do** *body*
- **for** *variable*: *initial\_value* **step** *increment* **while** *condition* **do** *body*
- **for** *variable*: *initial\_value* **step** *increment* **unless** *condition* **do** *body*

(Alternatively, the **step** may be given after the termination condition or limit.)

*initial\_value*, *increment*, *limit*, and *body* can be any expressions. If the increment is 1 then "**step 1**" may be omitted.

The execution of the **do** statement proceeds by first assigning the *initial\_value* to the variable (henceforth called the control-variable). Then: (1) If the control-variable has exceeded the limit of a **thru** specification, or if the condition of the **unless** is **true**, or if the condition of the **while** is **false** then the **do** terminates. (2) The body is evaluated. (3) The increment is added to the control-variable. The process from (1) to (3) is performed repeatedly until the termination condition is satisfied. One may also give several termination conditions in which case the **do** terminates when any of them is satisfied.

In general the **thru** test is satisfied when the control-variable is greater than the limit if the increment was non-negative, or when the control-variable is less than the limit if the increment was negative. The increment and limit may be non-numeric expressions as long as this inequality can be determined. However, unless the increment is syntactically negative (e.g. is a negative number) at the time the **do** statement is input, Maxima assumes it will be positive when the **do** is executed. If it is not positive, then the **do** may not terminate properly.

Note that the limit, increment, and termination condition are evaluated each time through the loop. Thus if any of these involve much computation, and yield a result that does not change during all the executions of the body, then it is more efficient to set a variable to their value prior to the **do** and use this variable in the **do** form.

The value normally returned by a **do** statement is the atom **done**. However, the function **return** may be used inside the body to exit the **do** prematurely and give it any desired value. Note however that a **return** within a **do** that occurs in a **block** will exit only the **do** and not the **block**. Note also that the **go** function may not be used to exit from a **do** into a surrounding **block**.

The control-variable is always local to the **do** and thus any variable may be used without affecting the value of a variable with the same name outside of the **do**. The control-variable is unbound after the **do** terminates.

```
(%i1) for a:-3 thru 26 step 7 do display(a)$
      a = - 3

      a = 4
```

```

a = 11
a = 18
a = 25
(%i1) s: 0$
(%i2) for i: 1 while i <= 10 do s: s+i;
(%o2) done
(%i3) s;
(%o3) 55

```

Note that the condition `while i <= 10` is equivalent to `unless i > 10` and also `thru 10`.

```

(%i1) series: 1$
(%i2) term: exp (sin (x))$
(%i3) for p: 1 unless p > 7 do
      (term: diff (term, x)/p,
      series: series + subst (x=0, term)*x^p)$
(%i4) series;

```

$$\frac{x^7}{90} - \frac{x^6}{240} - \frac{x^5}{15} - \frac{x^4}{8} + \frac{x^2}{2} + x + 1$$

which gives 8 terms of the Taylor series for  $e^{\sin(x)}$ .

```

(%i1) poly: 0$
(%i2) for i: 1 thru 5 do
      for j: i step -1 thru 1 do
        poly: poly + i*x^j$
(%i3) poly;

```

$$5x^5 + 9x^4 + 12x^3 + 14x^2 + 15x$$

```

(%o3) 5 x + 9 x + 12 x + 14 x + 15 x
(%i4) guess: -3.0$
(%i5) for i: 1 thru 10 do
      (guess: subst (guess, x, 0.5*(x + 10/x)),
      if abs (guess^2 - 10) < 0.00005 then return (guess));
(%o5) - 3.162280701754386

```

This example computes the negative square root of 10 using the Newton- Raphson iteration a maximum of 10 times. Had the convergence criterion not been met the value returned would have been `done`. Additional Forms of the `do` Statement

Instead of always adding a quantity to the control-variable one may sometimes wish to change it in some other way for each iteration. In this case one may use `next expression` instead of `step increment`. This will cause the control-variable to be set to the result of evaluating expression each time through the loop.

```

(%i6) for count: 2 next 3*count thru 20 do display (count)$
      count = 2
      count = 6

```

```
count = 18
```

As an alternative to `for variable: value ...do...` the syntax `for variable from value ...do...` may be used. This permits the `from value` to be placed after the step or next value or after the termination condition. If `from value` is omitted then 1 is used as the initial value.

Sometimes one may be interested in performing an iteration where the control-variable is never actually used. It is thus permissible to give only the termination conditions omitting the initialization and updating information as in the following example to compute the square-root of 5 using a poor initial guess.

```
(%i1) x: 1000$
(%i2) thru 20 do x: 0.5*(x + 5.0/x)$
(%i3) x;
(%o3)                2.23606797749979
(%i4) sqrt(5), numer;
(%o4)                2.23606797749979
```

If it is desired one may even omit the termination conditions entirely and just give `do body` which will continue to evaluate the body indefinitely. In this case the function `return` should be used to terminate execution of the `do`.

```
(%i1) newton (f, x):= ([y, df, dfx], df: diff (f ('x), 'x),
do (y: ev(df), x: x - f(x)/y,
if abs (f (x)) < 5e-6 then return (x)))$
(%i2) sqr (x) := x^2 - 5.0$
(%i3) newton (sqr, 1000);
(%o3)                2.236068027062195
```

(Note that `return`, when executed, causes the current value of `x` to be returned as the value of the `do`. The `block` is exited and this value of the `do` is returned as the value of the `block` because the `do` is the last statement in the block.)

One other form of the `do` is available in Maxima. The syntax is:

```
for variable in list end_tests do body
```

The elements of `list` are any expressions which will successively be assigned to the variable on each iteration of the body. The optional termination tests `end_tests` can be used to terminate execution of the `do`; otherwise it will terminate when the list is exhausted or when a `return` is executed in the body. (In fact, list may be any non-atomic expression, and successive parts are taken.)

```
(%i1) for f in [log, rho, atan] do ldisp(f(1))$
(%t1)                0
(%t2)                rho(1)
                    %pi
(%t3)                ---
                    4
(%i4) ev(%t3,numer);
(%o4)                0.78539816
```

**errcatch** (*expr\_1*, ..., *expr\_n*) Function

Evaluates *expr\_1*, ..., *expr\_n* one by one and returns [*expr\_n*] (a list) if no error occurs. If an error occurs in the evaluation of any argument, **errcatch** prevents the error from propagating and returns the empty list [] without evaluating any more arguments.

**errcatch** is useful in **batch** files where one suspects an error might occur which would terminate the **batch** if the error weren't caught.

**error** (*expr\_1*, ..., *expr\_n*) Function  
**error** Variable

Evaluates and prints *expr\_1*, ..., *expr\_n*, and then causes an error return to top level Maxima or to the nearest enclosing **errcatch**.

The variable **error** is set to a list describing the error. The first element of **error** is a format string, which merges all the strings among the arguments *expr\_1*, ..., *expr\_n*, and the remaining elements are the values of any non-string arguments.

**errormsg()** formats and prints **error**. This is effectively reprinting the most recent error message.

**errormsg** () Function

Reprints the most recent error message. The variable **error** holds the message, and **errormsg** formats and prints it.

**for** special operator

Used in iterations. See **do** for a description of Maxima's iteration facilities.

**go** (*tag*) Function

is used within a **block** to transfer control to the statement of the block which is tagged with the argument to **go**. To tag a statement, precede it by an atomic argument as another statement in the **block**. For example:

```
block ([x], x:1, loop, x+1, ..., go(loop), ...)
```

The argument to **go** must be the name of a tag appearing in the same **block**. One cannot use **go** to transfer to tag in a **block** other than the one containing the **go**.

**if** special operator

The **if** statement is used for conditional execution. The syntax is:

```
if <condition> then <expr_1> else <expr_2>
```

The result of an **if** statement is *expr\_1* if condition is **true** and *expr\_2* otherwise. *expr\_1* and *expr\_2* are any Maxima expressions (including nested **if** statements), and *condition* is an expression which evaluates to **true** or **false** and is composed of relational and logical operators which are as follows:

Operation	Symbol	Type
less than	<	relational infix
less than or equal to	<=	relational infix



syntactic equality	=	relational infix
equivalence	equal	relational function
not equal to	#	relational infix
greater than	>=	
or equal to		relational infix
greater than	>	relational infix
and	and	logical infix
or	or	logical infix
not	not	logical prefix

**map** (*f*, *expr\_1*, ..., *expr\_n*) Function

Returns an expression whose leading operator is the same as that of the expressions *expr\_1*, ..., *expr\_n* but whose subparts are the results of applying *f* to the corresponding subparts of the expressions. *f* is either the name of a function of *n* arguments or is a `lambda` form of *n* arguments.

`maperror` - if `false` will cause all of the mapping functions to (1) stop when they finish going down the shortest `expi` if not all of the `expi` are of the same length and (2) apply `fn` to [`exp1`, `exp2`,...] if the `expi` are not all the same type of object. If `maperror` is `true` then an error message will be given in the above two instances.

One of the uses of this function is to `map` a function (e.g. `partfrac`) onto each term of a very large expression where it ordinarily wouldn't be possible to use the function on the entire expression due to an exhaustion of list storage space in the course of the computation.

```
(%i1) map(f,x+a*y+b*z);
(%o1)          f(b z) + f(a y) + f(x)
(%i2) map(lambda([u],partfrac(u,x)),x+1/(x^3+4*x^2+5*x+2));
(%o2)          1      1      1
          ----- - ----- + ----- + x
          x + 2   x + 1          2
                                (x + 1)
(%i3) map(ratsimp, x/(x^2+x)+(y^2+y)/y);
(%o3)          1
          y + ----- + 1
          x + 1
(%i4) map("=", [a,b], [-0.5,3]);
(%o4)          [a = - 0.5, b = 3]
```

**mapatom** (*expr*) Function

Returns `true` if and only if *expr* is treated by the mapping routines as an atom. "Mapatoms" are atoms, numbers (including rational numbers), and subscripted variables.

**maperror** Variable

Default value: `true`

When `maperror` is `false`, causes all of the mapping functions, for example

```
map (f, expr_1, expr_2, ...)
```

to (1) stop when they finish going down the shortest `expr_i` if not all of the `expr_i` are of the same length and (2) apply `f` to `[expr_1, expr_2, ...]` if the `expr_i` are not all the same type of object.

If `maperror` is `true` then an error message is displayed in the above two instances.

**maplist** (*f*, *expr\_1*, ..., *expr\_n*)

Function

Returns a list of the applications of *f* to the parts of the expressions *expr\_1*, ..., *expr\_n*. *f* is the name of a function, or a lambda expression.

`maplist` differs from `map (f, expr_1, ..., expr_n)` which returns an expression with the same main operator as *expr\_i* has (except for simplifications and the case where `map` does an `apply`).

**prederror**

Variable

Default value: `true`

When `prederror` is `true`, an error message is displayed whenever the predicate of an `if` statement or an `is` function fails to evaluate to either `true` or `false`.

If `false`, `unknown` is returned instead in this case. The `prederror:false` mode is not supported in translated code.

**return** (*value*)

Function

May be used to exit explicitly from a block, bringing its argument. See `block` for more information.

**scanmap** (*f*, *expr*)

Function

**scanmap** (*f*, *expr*, *bottomup*)

Function

Recursively applies *f* to *expr*, in a top down manner. This is most useful when complete factorization is desired, for example:

```
(%i1) exp:(a^2+2*a+1)*y + x^2$
(%i2) scanmap(factor,exp);

(%o2)                2      2
                (a + 1) y + x
```

Note the way in which `scanmap` applies the given function `factor` to the constituent subexpressions of *expr*; if another form of *expr* is presented to `scanmap` then the result may be different. Thus, `%o2` is not recovered when `scanmap` is applied to the expanded form of *exp*:

```
(%i3) scanmap(factor,expand(exp));

(%o3)                2      2
                a y + 2 a y + y + x
```

Here is another example of the way in which `scanmap` recursively applies a given function to all subexpressions, including exponents:

```
(%i4) expr : u*v^(a*x+b) + c$
(%i5) scanmap('f, expr);
                f(f(f(a) f(x)) + f(b))
(%o5) f(f(f(u) f(f(v)                )) + f(c))
```

`scanmap (f, expr, bottomup)` applies  $f$  to  $expr$  in a bottom-up manner. E.g., for undefined  $f$ ,

```
scanmap(f,a*x+b) ->
  f(a*x+b) -> f(f(a*x)+f(b)) -> f(f(f(a)*f(x))+f(b))
scanmap(f,a*x+b,bottomup) -> f(a)*f(x)+f(b)
  -> f(f(a)*f(x))+f(b) ->
  f(f(f(a)*f(x))+f(b))
```

In this case, you get the same answer both ways.

### **throw** (*expr*)

Function

Evaluates  $expr$  and throws the value back to the most recent `catch`. `throw` is used with `catch` as a nonlocal return mechanism.

## 42 Debugging

### 42.1 Source Level Debugging

Maxima has a built-in source level debugger. The user can set a breakpoint at a function, and then step line by line from there. The call stack may be examined, together with the variables bound at that level.

The command `:help` or `:h` shows the list of debugger commands. (In general, commands may be abbreviated if the abbreviation is unique. If not unique, the alternatives will be listed.) Within the debugger, the user can also use any ordinary Maxima functions to examine, define, and manipulate variables and expressions.

A breakpoint is set by the `:br` command at the Maxima prompt. Within the debugger, the user can advance one line at a time using the `:n` (“next”) command. The `:bt` (“back-trace”) command shows a list of stack frames. The `:r` (“resume”) command exits the debugger and continues with execution. These commands are demonstrated in the example below.

```
(%i1) load ("/tmp/foobar.mac");

(%o1)                                     /tmp/foobar.mac

(%i2) :br foo
Turning on debugging debugmode(true)
Bkpt 0 for foo (in /tmp/foobar.mac line 1)

(%i2) bar (2,3);
Bkpt 0:(foobar.mac 1)
/tmp/foobar.mac:1::

(dbm:1) :bt                                <-- :bt typed here gives a backtrace
#0: foo(y=5)(foobar.mac line 1)
#1: bar(x=2,y=3)(foobar.mac line 9)

(dbm:1) :n                                <-- Here type :n to advance line
(foobar.mac 2)
/tmp/foobar.mac:2::

(dbm:1) :n                                <-- Here type :n to advance line
(foobar.mac 3)
/tmp/foobar.mac:3::

(dbm:1) u;                                <-- Investigate value of u
28

(dbm:1) u: 33;                             <-- Change u to be 33
33

(dbm:1) :r                                <-- Type :r to resume the computation
```

```
(%o2) 1094
```

The file `/tmp/foobar.mac` is the following:

```
foo(y) := block ([u:y^2],
  u: u+3,
  u: u^2,
  u);

bar(x,y) := (
  x: x+2,
  y: y+2,
  x: foo(y),
  x+y);
```

### USE OF THE DEBUGGER THROUGH EMACS

If the user is running the code under GNU emacs in a shell window (dbl shell), or is running the graphical interface version, `xmaxima`, then if he stops at a break point, he will see his current position in the source file which will be displayed in the other half of the window, either highlighted in red, or with a little arrow pointing at the right line. He can advance single lines at a time by typing M-n (Alt-n).

Under Emacs you should run in a `dbl` shell, which requires the `dbl.el` file in the `elisp` directory. Make sure you install the `elisp` files or add the Maxima `elisp` directory to your path: e.g., add the following to your `.emacs` file or the `site-init.el`

```
(setq load-path (cons "/usr/share/maxima/5.9.1/emacs" load-path))
(autoload 'dbl "dbl")
```

then in emacs

```
M-x dbl
```

should start a shell window in which you can run programs, for example Maxima, `gcl`, `gdb` etc. This shell window also knows about source level debugging, and display of source code in the other window.

The user may set a break point at a certain line of the file by typing C-x space. This figures out which function the cursor is in, and then it sees which line of that function the cursor is on. If the cursor is on, say, line 2 of `foo`, then it will insert in the other window the command, `":br foo 2"`, to break `foo` at its second line. To have this enabled, the user must have `maxima-mode.el` turned on in the window in which the file `foobar.mac` is visiting. There are additional commands available in that file window, such as evaluating the function into the Maxima, by typing Alt-Control-x.

## 42.2 Keyword Commands

Keyword commands are special keywords which are not interpreted as Maxima expressions. A keyword command can be entered at the Maxima prompt or the debugger prompt, although not at the break prompt. Keyword commands start with a colon, `'`. For example, to evaluate a Lisp form you may type `:lisp` followed by the form to be evaluated.

```
(%i1) :lisp (+ 2 3)
5
```

The number of arguments taken depends on the particular command. Also, you need not type the whole command, just enough to be unique among the break keywords. Thus `:br` would suffice for `:break`.

The keyword commands are listed below.

<code>:break F n</code>	Set a breakpoint in function <code>F</code> at line offset <code>n</code> from the beginning of the function. If <code>F</code> is given as a string, then it is assumed to be a file, and <code>n</code> is the offset from the beginning of the file. The offset is optional. If not given, it is assumed to be zero (first line of the function or file).
<code>:bt</code>	Print a backtrace of the stack frames
<code>:continue</code>	Continue the computation
<code>:delete</code>	Delete the specified breakpoints, or all if none are specified
<code>:disable</code>	Disable the specified breakpoints, or all if none are specified
<code>:enable</code>	Enable the specified breakpoints, or all if none are specified
<code>:frame n</code>	Print stack frame <code>n</code> , or the current frame if none is specified
<code>:help</code>	Print help on a debugger command, or all commands if none is specified
<code>:info</code>	Print information about item
<code>:lisp some-form</code>	Evaluate <code>some-form</code> as a Lisp form
<code>:lisp-quiet some-form</code>	Evaluate Lisp form <code>some-form</code> without any output
<code>:next</code>	Like <code>:step</code> , except <code>:next</code> steps over function calls
<code>:quit</code>	Quit the current debugger level without completing the computation
<code>:resume</code>	Continue the computation
<code>:step</code>	Continue the computation until it reaches a new source line
<code>:top</code>	Return to the Maxima prompt (from any debugger level) without completing the computation

### 42.3 Definitions for Debugging

#### `refcheck`

Variable

Default value: `false`

When `refcheck` is `true`, Maxima prints a message each time a bound variable is used for the first time in a computation.

**setcheck**

Variable

Default value: `false`

If `setcheck` is set to a list of variables (which can be subscripted), Maxima prints a message whenever the variables, or subscripted occurrences of them, are bound with the ordinary assignment operator `:`, the `::` assignment operator, or function argument binding, but not the function assignment `:=` nor the macro assignment `::=` operators. The message comprises the name of the variable and the value it is bound to.

`setcheck` may be set to `all` or `true` thereby including all variables.

Each new assignment of `setcheck` establishes a new list of variables to check, and any variables previously assigned to `setcheck` are forgotten.

The names assigned to `setcheck` must be quoted if they would otherwise evaluate to something other than themselves. For example, if `x`, `y`, and `z` are already bound, then enter

```
setcheck: ['x, 'y, 'z]$
```

to put them on the list of variables to check.

No printout is generated when a variable on the `setcheck` list is assigned to itself, e.g., `X: 'X`.

**setcheckbreak**

Variable

Default value: `false`

When `setcheckbreak` is `true`, Maxima will present a break prompt whenever a variable on the `setcheck` list is assigned a new value. The break occurs before the assignment is carried out. At this point, `setval` holds the value to which the variable is about to be assigned. Hence, one may assign a different value by assigning to `setval`.

See also `setcheck` and `setval`.

**setval**

Variable

Holds the value to which a variable is about to be set when a `setcheckbreak` occurs. Hence, one may assign a different value by assigning to `setval`.

See also `setcheck` and `setcheckbreak`.

**timer** (*f*<sub>1</sub>, ..., *f*<sub>*n*</sub>)

Function

**timer** ()

Function

Given functions *f*<sub>1</sub>, ..., *f*<sub>*n*</sub>, `timer` puts each one on the list of functions for which timing statistics are collected. `timer(f)$ timer(g)$` puts `f` and then `g` onto the list; the list accumulates from one call to the next.

With no arguments, `timer` returns the list of timed functions.

Maxima records how much time is spent executing each function on the list of timed functions. `timer_info` returns the timing statistics, including the average time elapsed per function call, the number of calls, and the total time elapsed. `untimer` removes functions from the list of timed functions.

`timer` quotes its arguments. `f(x) := x^2$ g:f$ timer(g)$` does not put `f` on the timer list.

If `trace(f)` is in effect, then `timer(f)` has no effect; `trace` and `timer` cannot both be in effect at the same time.

See also `timer_devalue`.

**untimer** (*f<sub>1</sub>*, ..., *f<sub>n</sub>*) Function  
**untimer** () Function

Given functions *f<sub>1</sub>*, ..., *f<sub>n</sub>*, `untimer` removes each function from the timer list.

With no arguments, `untimer` removes all functions currently on the timer list.

After `untimer (f)` is executed, `timer_info (f)` still returns previously collected timing statistics, although `timer_info()` (with no arguments) does not return information about any function not currently on the timer list. `timer (f)` resets all timing statistics to zero and puts `f` on the timer list again.

**timer\_devalue** Variable

Default value: `false`

When `timer_devalue` is `true`, Maxima subtracts from each timed function the time spent in other timed functions. Otherwise, the time reported for each function includes the time spent in other functions. Note that time spent in untimed functions is not subtracted from the total time.

See also `timer` and `timer_info`.

**timer\_info** (*f<sub>1</sub>*, ..., *f<sub>n</sub>*) Function  
**timer\_info** () Function

Given functions *f<sub>1</sub>*, ..., *f<sub>n</sub>*, `timer_info` returns a matrix containing timing information for each function. With no arguments, `timer_info` returns timing information for all functions currently on the timer list.

The matrix returned by `timer_info` contains the function name, time per function call, number of function calls, total time, and `gctime`, which meant "garbage collection time" in the original Macsyma but is now always zero.

The data from which `timer_info` constructs its return value can also be obtained by the `get` function:

```
get(f, 'calls); get(f, 'runtime); get(f, 'gctime);
```

See also `timer`.

**trace** (*f<sub>1</sub>*, ..., *f<sub>n</sub>*) Function  
**trace** () Function

Given functions *f<sub>1</sub>*, ..., *f<sub>n</sub>*, `trace` instructs Maxima to print out debugging information whenever those functions are called. `trace(f)$ trace(g)$` puts `f` and then `g` onto the list of functions to be traced; the list accumulates from one call to the next.

With no arguments, `trace` returns a list of all the functions currently being traced.

The `untrace` function disables tracing. See also `trace_options`.



`trace` quotes its arguments. Thus, `f(x) := x^2$ g:f$ trace(g)$` does not put `f` on the trace list.

When a function is redefined, it is removed from the timer list. Thus after `timer(f)$ f(x) := x^2$`, function `f` is no longer on the timer list.

If `timer(f)` is in effect, then `trace(f)` has no effect; `trace` and `timer` can't both be in effect for the same function.

**trace\_options** (*f*, *option\_1*, ..., *option\_n*) Function  
**trace\_options** (*f*) Function

Sets the trace options for function *f*. Any previous options are superseded. `trace_options(f, ...)` has no effect unless `trace(f)` is also called (either before or after `trace_options`).

`trace_options(f)` resets all options to their default values.

The option keywords are:

- `noprint` Do not print a message at function entry and exit.
- `break` Put a breakpoint before the function is entered, and after the function is exited. See `break`.
- `lisp_print` Display arguments and return values as Lisp objects.
- `info` Print `-> true` at function entry and exit.
- `errorcatch` Catch errors, giving the option to signal an error, retry the function call, or specify a return value.

Trace options are specified in two forms. The presence of the option keyword alone puts the option into effect unconditionally. (Note that option *foo* is not put into effect by specifying `foo: true` or a similar form; note also that keywords need not be quoted.) Specifying the option keyword with a predicate function makes the option conditional on the predicate.

The argument list to the predicate function is always [`level`, `direction`, `function`, `item`] where `level` is the recursion level for the function, `direction` is either `enter` or `exit`, `function` is the name of the function, and `item` is the argument list (on entering) or the return value (on exiting).

Here is an example of unconditional trace options:

```
(%i1) ff(n) := if equal(n, 0) then 1 else n * ff(n - 1)$
(%i2) trace (ff)$
(%i3) trace_options (ff, lisp_print, break)$
(%i4) ff(3);
```

Here is the same function, with the `break` option conditional on a predicate:

```
(%i5) trace_options (ff, break(pp))$
(%i6) pp (level, direction, function, item) := block (print (item),
  return (function = 'ff and level = 3 and direction = exit))$
(%i7) ff(6);
```

**untrace** (*f\_1*, ..., *f\_n*)

Function

**untrace** ()

Function

Given functions *f\_1*, ..., *f\_n*, **untrace** disables tracing enabled by the **trace** function.

With no arguments, **untrace** disables tracing for all functions.

**untrace** returns a list of the functions for which it disabled tracing.



## 43 Indices



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