TRIP is a general computer algebra system dedicated to celestial mechanics.

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This software includes a library developed by the NetBSD foundation and its contributors.
This software includes the libraries LAPACK and 'SCSCP C Library'.
This software is dynamically linked to the libraries LTDL, GMP, MPFR et MPC. The source of these libraries could be downloaded from the web site specified in the Chapter References (see Chapter 19 [References], page 209).

TRIP version 1.4.120

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1 Introduction

In order to execute TRIP, you must
− Under Unix or MacOs X, enter trip in a shell window.
− Under Windows, click on trip icon for the console version.
− Under Windows, click on tripstudio icon for the graphic version.

You will see on screen :

```
**** BIENVENUE SUR TRIP v1.0.0 ****
Taper 'help;' ou '?;' pour obtenir de l’aide
>
```

TRIP is ready to receive your commands. TRIP has a case sensitive interpreter.

In order to compute \( S = (X + Y)^2 \), you must just enter :

```
**** BIENVENUE SUR TRIP v1.0.0 ****
Taper 'help;' ou '?;' pour obtenir de l’aide
>
S=(X+Y)^2;
S(X,Y) = 1*Y**2 + 2*Y*X + 1*X**2
```

Use the up-arrow key, you could bring back the previous command. Command-line input can be edited using arrow keys.

To exit TRIP, you enter `quit` or `exit`.

The reserved keywords for TRIP language are specified in the Annexes. TRIP contains several keyword categories :
− TRIP global variables, TRIP constants and mathematic symbol.
− Functions : command returns a value
− Procedures : command doesn’t return a value.

When TRIP starts or the commands reset or @@ are executed, it reads these files in the following order :

- ‘\$TRIPDIR/etc/trip.ini’ where \$TRIPDIR is the installation directory of TRIP.
- ‘\$HOME/.trip’ where \$HOME is the user home directory.
- ‘trip.ini’ in the current directory.

These file may contain any valid instructions.

Recommendations :
− The following environment variables should be set and exported:
  ● TMPDIR : pathname of a directory in which temporary files may be written.
  ● LC_MESSAGES : language used to display messages.
2 Semantic

2.1 expression
An expression can contain all defined operators on object identifiers and can call functions or commands. An expression must be always ended with ; or $. The character ; implies the execution of the expression and displays the result. The character $ implies only the execution of the expression. Several instructions separated with the character ; or $ could follow. They will be executed sequentially.

A comment on one or more lines is defined as /* ... */, like in the C language. A single-line comment starts with the characters //, like in the C++ language.

Example:
> // compute 1+t and display the result
> s=1+t;
> s(t) =  
>  1 + 1*t

> // compute 1+t but does not display the result
> s1=(1+x+y)^2$
> // display the content of s1
> s1;
> s1(x,y) =  
>  1 + 2*y + 1*y**2 + 2*x + 2*x*y + 1*x**2


2.2 object identifier
An object identifier must begin with a letter and must contain only letters (lower-case or upper-case) and/or number (09) and the characters _ or '.

By default, an object identifier is global, so it will be viewed by any instruction. An object identifier could be local to a macro, so it will be only viewed from this macro and it will be destroyed each time the execution of this macro finished. A TRIP reserved keyword can’t be used as the name of an object identifier.

The object identifier can have the following type:
- variable
- serie
- constant
- array of series
- array of variables
- numerical vector
- numerical matrix
- string of characters
structure

Example :
> ch="file1"
> s=1+x
> dim t[1:2];
> z=1+2*i;
> bilan;
  ch CHAINE
  s SERIE
  t TAB
  x VAR

The definition of the function, subroutines and variables use the following types :

<object identifier> object identifier
<integer> natural integer number or an operation which returns an integer
<real> real floating-point number or an operation which returns a real floating-point number
<complex> complex number or an operation which returns a complex number
<variable> variable
<constant> an operation which returns an integer, a real number or a complex number
<serie> serie
<operation> an operation which returns a number (constant) or a serie
<array> array of series
<array of variables> array of variables
<array of) variables> variable or array of variables
<num. vec.> numerical vector
<real vec.> numerical vector of real numbers
<complex vec.> numerical vector of complex numbers
<array of real vec.> array of numerical vectors of real numbers
<array of complex vec.> array of numerical vectors of complex numbers
<array of num. vec.> array of numerical vectors
<(array of) real vec.> (array of) numerical vectors of real numbers
<(array of) num. vec.> (array of) numerical vectors
<constant or num. vec.> a constant or a numerical vector
<constant or matrix> a constant or a numerical matrix
<real or real vec.> real floating-point number or numerical vector of real numbers

matrixtype numerical matrix
matrixrtype real matrix
matrixctype complex matrix
<filename> file
<file> macro
<macro> <dimension of an array> two integers separated with the character :
<list of dimension> list of <dimension of an array> separated with a comma
<two dimensions of a matrix> list of two dimensions separated with a comma. Each dimension consists of two integers separated with the character :
Chapter 2: Semantic

<condition> comparison between two operations
<string> an operation which returns a string of characters
<(array of) string> an operation which returns a string of characters or an array of string of characters
<List of variables> list of variables or an array of variables
<List of parameters> list of parameters of a macro
<body> list of trip expression
<scscp client> connection to a SCSCP server
<remote object> object stored on a remote SCSCP server
<Maple client> connection to a local Maple session
<Mathematica client> connection to a local Mathematica session

2.2.1 serie
A serie is a polynomial of degree n (n is an integer). It depends on one or more variables.

Example:
> s=1+x;
s(x) = 1 + 1*x
> s2=(1+x+y);
s2(x,y) = 1 + 1*y + 1*x
> bilan;
s SERIE
s2 SERIE
x VAR
y VAR

2.2.2 constant
A constant is an integer, a real, a complex, a rational number or an interval. In rational numerical mode (\_modenum = NUMRAT or \_modenum = NUMRATMP), the input of 0.5 creates a floating-point number and not the rational number 1/2.

Example:
> x = 2;
2
> y = 2.25;
9/4
> z = 1+2*i;
(1+i*2)
> bilan;
x CONST
y CONST
z CONST
> \_modenum=NUMDBLINT;
\_modenum = NUMDBLINT
> s = <1|2>;
s = [+1.00000000000000E+00,+2.00000000000000E+00]

2.2.3 string
It’s a list of characters surrounded with " ". It’s used to specify the _path, to define the name of the files, to display messages, ... . There is no limit on the size. But, for the global variable _path, it’s length is truncated to 256 (this value depends on the system). To produce a double-quotes (" ) in a string, two double-quotes must be used.

Example:
> ch = "file" + "1";
> ch = "file1"
> sch = "../" + ch;
> sch1 = sch + "." + str(12);
> ch3="nomsuivide""fin";
> ch3 = "nomsuivide"fin"

2.2.4 real

\(<\text{nom}> = <\text{real}> ;

A real floating-point number is always displayed or inputed in base 10. The symbol \(d\), \(e\), \(E\) or \(D\) is the exponentiation symbol.

Example :
> a=2.125E6;
a = 2125000
> q=3D2;
q = 300
> r=0.123554545;
r = 0.123554545

2.2.5 complex

\(<\text{nom}> = <\text{real}> + i * <\text{real}> ;

A complex number is allowed. The lower-case \(i\) or upper-case \(I\) specifies the imaginary part of the complex number.

Example :
> x=2+i*3;
x = (2+i*3)
> y=-5+i*4;i;
y = (-5+i*4)

2.2.6 filename

It’s a string of characters. This string must be surrounded with " " if the string contains the character space " [ ] { } ( ) or more than one points. A file name could be the content of an object identifier of type string. These file names are relative to the variable .path, excepted for the file names build with the function file_fullname (see Section 18.20 [file_fullname], page 205).

Example :
> read("fichier.1.dat",T);
> s="ell."+str(10)+".asc";
s = "ell.10.asc"
> read(s,T);

2.2.7 file

It’s an object identifier which specifies a file, which is opened in reading or writing mode.

Example :
> f = file_open("fichier.1.dat",read);
> file_close(f);

2.3 Visibility
2.3.1 private

**private**

private <object identifier> x , ...

The specified object identifiers will be local to the macro, to the loop (for, while, sum) or to a file. The local object identifiers are destroyed at the end of the execution of the macro or at the end of each iteration of loop (for, while, sum).

If a local object identifier is local to a file, this object is visible only to the macros of this file or during the execution of the file.

If a local object identifier has the same name as a global object identifier, then the local object identifier will be used. In this case, the local object identifier hides the global object identifier during the execution of the macro.

**private _ALL**

private _ALL;

All object identifiers used in a macro and a file will be local to the macro or to a file.

The behavior is similar to the previous definition.

The command **public** (see Section 2.3.2 [public], page 8) allow to access to a global object.

Example :

> // S, SY, SXY, SXX, DEL are private objects
> macro least_ab[TX,TY,a,b] {
  private S, SY, SXY, SXX, DEL;
  S=size(TX)$
  SX=sum(TX)$
  SY=sum(TY)$
  SXY=sum(TX*TY)$
  SXX=sum(TX*TX)$
  DEL=S*SXX-SX*SX$
  a = (S*SXY-SX*SY)/DEL$
  b = (SXX*SY-SX*SXY)/DEL$
}
>
> tx=1,10;
  tx double precision real vector : number of elements =10
> ty=3*tx;
  ty double precision real vector : number of elements =10
> %least_ab[tx,ty,[a],[b]];
> bilan;
  SX CONST
  a CONST
  b CONST
  tx VNUMR
  ty VNUMR
>
> // T2 is a private object
> macro detval[T] {
  private _ALL[T];
  dim T2[1:3,1:3]$ T2[:,::]=0$
  T2[1,1]=T**2$
  T2[2,2]=1-T$
}
T2[3,3]=T$
  \text{return } \det(T2);
};
  > \%\text{detval}[5];
    -500
  > \text{bilan};
  SX \text{CONST}
a \text{CONST}
b \text{CONST}
tx \text{VNUMR}
ty \text{VNUMR}
>

2.3.2 public

public <object identifier> x ,...;  
[Commande]

This command must be only used after the command \texttt{private _\texttt{ALL};}. It specifies that the 
identifiers of this list will be global whereas the other objects will be local to the macro or to 
the file.

\textbf{Example :}
> // w is local and z is global
> \texttt{macro myfunc}[T] 
  \{ 
  \texttt{private _\texttt{ALL};}
  \texttt{public z;}
  \texttt{w = cos(T);}
  \texttt{z = T;}
  \};
> \%\text{myfunc}[5];
  w = 0.2836621854632262
  z = \text{5}
  > \text{bilan};
  z \text{CONST}
  >
3 Global variables

The content of a global variable is performed by entering its name followed with ;. The content of all global variables is displayed with the command vartrip (see Section 18.12 [vartrip], page 201). All global variables, containing a real number, are stored in a double-precision real number.

3.1 _affc

integer _affc

_set = {1, 2, 3, 4, 5, 6, 7, 8, 9};

Set the format to display the numerical coefficients of series or constants.

Default value = 4.

1 : %G (default short format)
2 : %.8G (8 digits after the decimal points)
3 : %.10G (10 digits after the decimal points)
4 : %.20.15G (15 digits or more after the decimal points)
5 : rational approximation (double-precision real numbers are converted to rationals number using the continued fraction, else %.8G)
6 : same as 5 but %.15G (15 digits after the decimal points)
7 : rational approximation (fixed length format)
8 : %15.6E (6 digits after the decimal points)
9 : same as 8 but displays the diameter of intervals

Example:
> _affc=1;
  _affc = 1
> 1/3;
  0.333333
> _affc=2;
  _affc = 2
> 1/3;
  0.33333333
> _affc=3;
  _affc = 3
> 1/3;
  0.3333333333
> _affc=4;
  _affc = 4
> 1/3;
  0.333333333333333
> _affc=5;
  _affc = 5
> 1/3;
  1/3
> _affc=6;
  _affc = 6
> 1/3;
  1/3
> _affc=7;
.affc = 7
> 1/3;
    1/3
> .affc=8;
    .affc = 8
> 1/3;
    3.333333E-01
> .affc=9; _modenum=NUMDBLINT;
    .affc = 9
    _modenum = NUMDBLINT
> 1/3;
    [+3.333333333333333E-1,+3.333333333333334E-1 ( 5.551115E-17)]

3.2 _affdist

integer _affdist

_affdist = {0, 1, 2, 3, 4, 5, 6, 7, 8};

Set the format to display the series

Default value = 2.

= 0 : recursive.
= 1 : distributed.
= 2 : distributed with one term by line.
= 3 : distributed with (co)sinus .
= 4 : distributed with (co)sinus and one term by line.
= 5 : distributed and aligned with (co)sinus and one term by line.
= 6 : distributed and factorized for variables specified in _affvar.
= 7 : distributed and factorized for variables specified in _affvar, with one term by line
    for the two parts.
= 8 : distributed and factorized for variables specified in _affvar, with one term by line
    for the distributed part and the factorized part is on the same line.

Remarks : if _affvar is empty, then the format of _affdist = 6, respectively 7 or 8, is the
same as _affdist=1, respectively 2.

Example :
> _affdist = 2;
> x = 1+y;
  x(y) =
  1
  + 1*y

3.3 _comment

boolean _comment

_comment {on,off};

Enable or disable the display of comments.

Default value = off.

Example :
> _comment;
Chapter 3: Global variables

3.4 _cpu

integer _cpu

Specify the number of processors which will be used for later computations.

The specified value can’t be greater than the number of available processors or cores on the computer.

Default value: number of active processors or cores on the computer.

Example:
> _mode=POLP;
> _cpu=1;
> time_s; s=(1+x+y+z+t+u)^30$ time_t(usertime, realtime);
03.321s - (99.44% CPU)
> msg("the real time is %g\n", realtime);
the real time is 3.33976
> _cpu=4;
> time_s; s=(1+x+y+z+t+u)^30$ time_t(usertime, realtime);
03.339s - (322.89% CPU)
> msg("the real time is %g\n", realtime);
the real time is 1.03414

3.5 _echo

boolean _echo

Enable or disable the display of the echo of the executed commands.

Default value = off.

= 1: Display the echo of the commands (same as on).
= 0: Don’t Display the echo of the commands (same as off)

Example:
>_echo = 1;
>msg "exemple"
exemple
cde>>
msg "exemple"
### 3.6 `_ endian`

**name** `_ endian` [Variable]

```
_endian = {little, big};
```

Specify the byte order (big-endian or little-endian) in binary files.

Default value = endianness in the computer.

- `big`: big-endian
- `little`: little-endian

Example:
```
> _endian = big;
> _endian = big
> vnumR ieps;
> readbin(file1.dat, "%u", ieps);
```

### 3.7 `_ graph`

**name** `_ graph` [Variable]

```
_graph = gnuplot;
```

Specify if the plot, replot, plotps, plotf commands will use the grace or gnuplot (see Chapter 12 [Graphiques], page 111) software to display graphics.

Default value = gnuplot.

- `gnuplot`: gnuplot will display the graphics.
- `grace`: grace will display the graphics.

Example:
```
> _graph = grace;
> _graph = gnuplot
```

### 3.8 `_ hist`

**boolean** `_ hist` [Variable]

```
_hist {on, off};
```

Specify if the file `history.trip` is opened or not.

If the file is opened, all commands are stored in this file.

Default value = on.

Example:
```
>_hist;
_hist = ON
```

### 3.9 `_ history`

**string** `_ history` [Variable]

```
_history = <string>;
```

Specify the folder where the file `history.trip` will be written.

Default value = "".
Example :
> _history="/users/toto/";
    _history = /users/toto/

3.10 _info

boolean _info [Variable]

    _info {on, off};

Enable or disable the display of the message information.
Default value = on.

Example :
> _modenum=NUMQUAD;
    _modenum = NUMQUAD
> x1=0,2*pi,2*pi/59$
> x2=0,3,0.5$
> _info off;
    _info off
> sl=interpol(LINEAR,x1,cos(x1),x2)$
> _info on;
    _info on
> sl=interpol(LINEAR,x1,cos(x1),x2)$
    Information : interpol perform computations in double precision.

3.11 _integnum

boolean _integnum [Variable]

    _integnum = {DOPRI8, ODEX1, ADAMS};

Defines the numerical integrator used by the numerical integration performed by integnum (see Section 17.4 [integnum], page 183) and integnumfcn (see Section 17.5 [integnumfcn], page 188).
Default value = DOPRI8.


Example :
> _integnum=ODEX1;
    _integnum = ODEX1
> _integnum=DOPRI8;
    _integnum = DOPRI8

3.12 _language

_language [Variable]

    _language = {fr, en};

Select the language to display messages.
Default value =
- On Unix, Linux and MacOS X, depends on the environment variable LC_MESSAGES.
  If it doesn’t exist, the french language is used.
- On Windows, depends on the system preferences.

Example :
> _language = en ;
  _language = en
> Exp(x, y);
TRIP[ 3 ] : This parameter must be an integer
  Command : ' Exp(x, y);'

> _language = fr;
  _language = fr
> Exp(x, y);
TRIP[ 3 ] : Cet argument doit etre un entier
  Commande : 'Exp(x, y);'

3.13 _mode

_mode [Variable]

_mode = {POLY, POLP, POLYV, POLPV};

Specify the internal representation of the series for their storage in the memory and during the computations.

The internal representation could be changed during the computations.

Default value = POLY.
- POLY : series are stored using a recursive list (recursive sparse representation). This is the default mode. The leaf nodes use generic containers. This mode is efficient for the very degree.
- POLP : series are stored using a recursive vector (recursive dense representation). The leaf nodes use generic containers. It should be used if the degree of series are low.
- POLYV : series are stored using a recursive list (recursive leaf sparse representation). The leaf nodes use optimized containers. This mode is efficient for the very degree.
- POLPV : series are stored using a recursive vector (recursive leaf dense representation). The leaf nodes use optimized containers. It should be used if the degree of series are low.

Example :
> _mode=POLY;
  _mode = POLY
> s1=(1+x)^100$
> s2=1+x^100$
> bilan mem;

<table>
<thead>
<tr>
<th>Nom</th>
<th>Memoire (octets)</th>
<th>Nb de termes</th>
</tr>
</thead>
<tbody>
<tr>
<td>s1</td>
<td>4848</td>
<td>101</td>
</tr>
<tr>
<td>s2</td>
<td>96</td>
<td>2</td>
</tr>
<tr>
<td></td>
<td>4944</td>
<td></td>
</tr>
</tbody>
</table>

Memoire physique utilisee = 7118848 octets
Memoire maximale utilisee = 2783817728 octets
> _mode=POLP;
  _mode = POLP
> s1=(1+x)^100$
> s2=1+x^100$
> bilan mem;

<table>
<thead>
<tr>
<th>Nom</th>
<th>Memoire (octets)</th>
<th>Nb de termes</th>
</tr>
</thead>
<tbody>
<tr>
<td>s1</td>
<td>3232</td>
<td>101</td>
</tr>
<tr>
<td>s2</td>
<td>3232</td>
<td>2</td>
</tr>
</tbody>
</table>

Memoire physique utilisee = 7208960 octets
Memoire maximale utilisee = 2784210944 octets

> _mode=POLY$ time_s; s1=(1+x)^100$ time_t;
00.240s - ( 47.74% CPU)
> _mode=POLP$ time_s; s1=(1+x)^100$ time_t;
00.013s - ( 40.57% CPU)

### 3.14 _modenum

_modenum

 modenum = {NUMDBL, NUMRAT, NUMRATMP, NUMQUAD, NUMFPMP};

Specify the representation of the numerical coefficients in the series and the precision of the floating point numbers in the numerical vectors or matrices.

This mode could be changed during the session.

Default value = NUMDBL.

The numerical coefficients are encoded as followed:

- NUMDBL : double precision floating-point. (default)
- NUMQUAD : quadruple precision floating-point. This representation isn’t supported on Windows operating system.
- NUMRAT : integer or rational number, otherwise double precision floating-point.

The absolute value for the numerator and denominator is between 0 et 2**62 on most computers. If the numerator or denominator becomes too large, rational number is converted in a double precision floating-point.

The real numbers are stored in double precision format. For example, the coefficient "2.0" will be stored as a double precision floating-point while the coefficient "2" is stored as a rational number.

- NUMRATMP : multiprecision rational, otherwise double precision real.

The absolute value for the numerator and denominator don’t have any limits.

The real numbers are stored in double precision format. For example, the coefficient "2.0" will be stored as a double precision floating-point while the coefficient "2" is stored as a rational number.

This numerical representation uses the GMP library (type mpz_t or mpq_t). On the 64 bit operating systems, integers smaller than $2^{-63}-1$ (in absolute value) use the hardware instructions instead of the GMP library.
NUMFPMP : multiprecision floating-point.

The number of number of digits of precision is specified by the global variable \_modenum\_prec (see Section 3.15 \[\_modenum\_prec\], page 16). This numerical representation uses the MPFR and the MPC library.

In the mode NUMRAT and NUMRATMP, numerical vectors or matrices contain always double precision floating-points (real or complex numbers).

Example :
> \_affc=4;
  \_affc = 4
> \_modenum=NUMDBL;
  \_modenum = NUMDBL
> s=1/11;
  s = 0.0909090909090909
> \_modenum=NUMRAT;
  \_modenum = NUMRAT
> s=1/11;
  s = 1/11
> \_modenum=NUMFPMP;
  \_modenum = NUMFPMP
> pi;
  3.1415926535897932384626433832795028841971693993751058209749445924

3.15 \_modenum\_prec

integer \_modenum\_prec [Variable]

\_modenum\_prec = \<integer\>

Variable used if \_modenum = NUMFPMP (see Section 3.14 \[\_modenum\], page 15).

Specify the number of significant decimal digits for the multiprecision real numbers.

Default value = 64.

Example :
> \_modenum=NUMFPMP;
  \_modenum = NUMFPMP
> \_modenum\_prec= 20;
  \_modenum\_prec = 20
> pi;
  3.14159265358979323846
> \_modenum\_prec= 40;
  \_modenum\_prec = 40
> pi;
  3.1415926535897932384626433832795028841975

3.16 \_naf\_dtour

real \_naf\_dtour [Variable]

\_naf\_dtour = \<real\>

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).

Specify the "length of the dial turn ".

Default value = 2*\pi.

Example :
> \_naf\_dtour=360;
3.17 _naf_icplx

integer _naf_icplx
_naf_icplx = {0, 1};

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify if the function is real or complex.
Default value = 1.
The available values are:
= 0 : real function
= 1 : complex function

Example :
> _naf_icplx=0;

3.18 _naf_iprt

integer _naf_iprt
_naf_iprt = {-1, 0, 1, 2};

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify the verbosity to display messages in the commands naf and naftab. Intermediate
results are stored in a file.
Default value = -1.
The available values are:
= -1 : No display.
= 0 : Few messages are displayed.
= 1 : More messages are displayed.
= 2 : Many messages are displayed.

Example :
> _naf_iprt = 2;

3.19 _naf_isec

integer _naf_isec
_naf_isec = {0, 1};

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify if the algorithm "secantes" is used or not in the commands naf and naftab.
Default value = 1.
The available values are:
= 0 : algorithm "secantes" is not used.
= 1 : algorithm "secantes" is used.

Example :
> _naf_isec=1;
3.20 _naf_iw

integer _naf_iw [Variable]

_naf_iw = <integer> ;

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify if a window filter is set.
Default value = 1.
The available values are :

= -1 : exponential window \( PHI(T) = 1/CE \times \exp\left(-1/(1-T^2)\right) \)
   avec \( CE=0.22199690808403971891E0 \)
= 0 : no window
= N > 0 : \( PHI(T) = C_N \times (1 + \cos(\pi T))^N \) avec \( C_N = 2^N \times (N!)^2/(2N)! \)

Example :
> _naf_iw=-1;

3.21 _naf_nulin

integer _naf_nulin [Variable]

_naf_nulin = <integer> ;

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify the number of lines to be ignored at the beginning of the data file.
Default value = 1.

Example :
> _naf_nulin=0;

3.22 _naf_tol

real _naf_tol [Variable]

_naf_tol = <real> ;

Variable used by naf (see Section 17.1 [Analyse en frequence], page 177).
Specify the tolerance to check if two frequencues are equal.
Default value = 1E-10.

Example :
> _naf_tol=1E-4;

3.23 _path

string _path [Variable]

_path = <string> ;

_path = "path directory";

Specify the folder used to save or load files. All TRIP commands or functions writing or
reading files will prepend the filename with this path. loading trip programs (include),
creating psotscript files(plotps) and plotting files (plotf) will use this path.
Default value = "" (current folder).
Remarkes :
− Don’t miss the character "/" at the end of the path on UNIX or MACOS X.
− Don’t miss the character "\" at the end of the path on WINDOWS.
3.24 _quiet

boolean _quiet {on, off};
Remove, respectively enable, the future display produced by the commands followed by ;
only in the current macro. Default value = off.

Example:
> macro silence{ s=1; _quiet on; s=pi; }
> %silence;
> stat(s);
constant has name s and has the value 3.141592653589793

3.25 _read

string _read

_example = ();
_example = ("skipbrokenlines", "skipemptylines");
_example is a list of no, one or more strings which specify the behavior of the functions read
and file_read when an empty or partial line is encountered.
_example = () specifies that these functions will generate an error when such lines are encoun-
tered.
Les valeurs possibles sont:

- "skipbrokenlines": the partial lines are ignored and a warning message is displayed with
the indices of the ignored lines.
- "skipemptylines": the empty lines are ignored and a warning message is displayed with
the indices of the ignored lines.

Default value = ()

Example:
> t1=1,3;
t1 double precision real vector : number of elements = 3
> t2=11,13;
t2 double precision real vector : number of elements = 3
> f=file_open(temp001, write);
f = file "temp001" opened in write mode
> file_writemsg(f,"# header line\n");
> file_write(f,t1);
> file_writemsg(f,"err...\n");
> file_write(f,t2);
> file_close(f);
> _read= ("skipbrokenlines");
_example = ("skipbrokenlines")
> vnumR Q;
> read(temp001, Q);
Information : The line 1 is ignored : the line is incomplete or invalid
Information: The line 5 is ignored: the line is incomplete or invalid
> writes(Q);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
+1.1000000000000000E+01
+1.2000000000000000E+01
+1.3000000000000000E+01

3.26 _read_history

string _read_history [Variable]
_read_history = <string>;

_read_history specifies the name of the file in which the skipped errors by _read are recorded. This file is emptied when this variable is assigned. Each line of this ascii file contains two columns: the line number and the name of the file where the error is produced.

_read_history = "" specifies that the skipped errors by _read are not recorded in a file.

Default value = "".

Example:
> t1=1,3;
  t1 double precision real vector: number of elements =3
> t2=11,13;
  t2 double precision real vector: number of elements =3
> f=file_open(temp001, write);
  f = file "temp001" opened in write mode
> file_writemsg(f,"# header line\n");
> file_write(f,t1);
> file_writemsg(f,"err...\n");
> file_write(f,t2);
> file_close(f);
> _read=( "skipbrokenlines" );
  _read = ( "skipbrokenlines" )
> _read_history="errors.dat"
  _read_history = "errors.dat"
> vnumR Q;
> read(temp001, Q);
  Information: The line 1 is ignored: the line is incomplete or invalid
  Information: The line 5 is ignored: the line is incomplete or invalid
> vnumR lerr;
> fmt="%g %s"
  fmt = "%g %s"
> read("errors.dat", fmt, lerr, ferr);
> writes("%g\n",lerr);
1
5
> afftab(ferr);
  ferr[1] = "temp001"
  ferr[2] = "temp001"
### 3.27 _time

**boolean _time**

```plaintext
_time {on, off};
```

enable or disable display of partial time after each executed command.

When the command `_time on` is performed, an implicit call to `time_s` and the time `0.00s` is always displayed after it.

Default value = `off`.

**Example:**
```
> macro temps {
  s=(1+x+y+z)**20$
  _time on;
  s=(1+x+y+z)**20$
  q=s$
  _time off;
};

> %temps;
```

### 3.28 _userlibrary_path

**string _userlibrary_path**

```plaintext
_userlibrary_path = <string> ;
```

```
_userlibrary_path = "path to the user functions";
```

When this variable is assigned, TRIP reads recursively the specified directory all the file with the extension .t. It loads all the macros which are preceded by a comment starting by `//!trip extern_function` and renders them as functions to the user.

The format of this comment:
```
//!trip extern_function name_of_the_function attribute="parameter_attribute" "parameter_inout"
```

The string `parameter_attribute` may contain the following keywords separated by a comma.

- if this string is empty, then the the function is visible by the user.
- `private` : the function is private and is not visible by the user.
- `parametres_inout` must contain the same number of elements as the number of parameter of the function. Each element are separated with a comma. An element could have the following values:
  - `in` : the parameter is an input only.
  - `out` : the parameter is an output only.
  - `inout` : the parameter is an input and output.

Several macros may have the same value `name_of_the_function` if they have a different number of parameters.

Default value = "".
Example:

_userlibrary_path="./myuserlibrary/";
userfunc1(3);
userfunc2(3,y);

The directory myuserlibrary contains the following file:

Example:

```cpp
//!trip extern_function userfunc1 "in"
macro userfunc1[x]
{
    return 2*x+1;
};

//!trip extern_function userfunc2 "in,out"
macro userfunc2[x,y]
{
    y=2*x+1;
};
```

3.29 I

complex i [constant]
complex I [constant]

\[ i^2 = -1 \]

Example:

\[ z = x + y*i; \]
\[ z(x,y) = (0+i*1)*y+1*x \]

3.30 PI

real pi [constant]
real PI [constant]

The mathematical symbol \( \pi \)

Example:

\[ z = \pi; \]
\[ 3.14159265358979 \]
4 Variables

4.1 crevar

crevar

Function

crevar (<string or name> radical, <integer> indice1, <integer> indice2,...);
Create and return a variable with the name radical +"_" + indice1 +...+ "_" + indicen.
crevar (<string or name> radical);
Create and return a variable with the name radical. This function allow to create variables with a non-standard name.

Example :
> dimvar t[1:3];
> t[1]:=crevar("t",1);
  t[1] = t_1 = 1*t_1
> t[3]:=crevar("tab",3,2,1);
  t[3] = tab_3_2_1 = 1*tab_3_2_1

> afftab(t);
  t[1] = t_1 = 1*t_1
  t[2] =
  t[3] = tab_3_2_1 = 1*tab_3_2_1
> dimvar u[1:2];
> u[1]:=crevar("\bar{X}");
  u[1] = \bar{X} = 1*\bar{X}
> u[1];
  u[1] = \bar{X} = 1*\bar{X}

4.2 Operators

:=

Operator

<object identifier> [...] := <variable> ;
Create a reference to an existing variable.

Example :
> // w is a reference to x
> s=(1+x+y)**3$
> w:=x$
> stat(w);
Variable x type : 2 ordres : 2 2 2 2
dependances :
variables dependant de celle-ci :
> deriv(s,w);
3
+ 6*y
+ 3*y**2
+ 6*x
+ 6*x*y
+ 3*x**2
>

5 Series

5.1 Operators

\(+\ -\ *\ /\)

TRIP performs the addition (or unary plus), the subtraction (or unary minus), multiplication, or division between series, variables, numbers and expressions.

To perform the addition, multiplication, subtraction and division of the arrays, the operators are \(+\), \(*\), \(-\), and \(/\).

The operators \(+\), \(*\), \(-\), \(/\) could be applied on numerical vectors.

The unary \(+\) or \(-\) could be applied to any type.

Remark : The division by a serie isn’t available. You should use the function \texttt{div} to perform the euclidian division.

Example :
> s=1\times-1-y+z+t/2;
s(x,y,z,t) = 1 + 1/2*t - 1*y*z + 1*x
> \_modenum=NUMRATMP;
\_modenum = NUMRATMP
> p=5/3;
p = 5/3
> \_modenum=NUMQUAD;
\_modenum = NUMQUAD
> p=5/3;
p = 1.6666666666666666666666666666667

** ^

\langle\text{serie}\rangle ^{\langle\text{real}\rangle}

It performs the exponentiation.

It supports integer powers.

The exponentiation \(*\)-1 on an array, with 2 same dimensions, performs the matrix inversion.

Remarks : The exponentiation, term by term, on arrays isn’t available.

Example :
> u=2^{3};
8
> z=(1+i)^{3};
(-2+i*2)

> s=(1+x+y)**2;
s(y,x) = 1 + 2*x + 1*x**2 + 2*y + 2*y*x + 1*y**2

5.2 Standard functions

\texttt{size}

\texttt{size(<serie> \ s \ )}

It returns the number of terms in the serie \( s \).

Example :
> s=1+x\times y+z\times t-i*p;
s(x,y,z,t,p) = 1 + (-0-i*1)*p + 1*t*z + 1*y*x
> size(s);
4
5.3 Differentiation and integration

5.3.1 deriv

derv <serie>, <variable>

It computes the total derivative of the serie with respect to one variable.

Example:
> s = (1 + x + y)**2;
> deriv(s, x);
    2 + 2*y + 2*x

5.3.2 integ

integ <serie>, <variable>

It integrates the serie with respect to one variable.

Example:
> s = (1 + x + y)**2;
> integ(s, x);
    1*x + 2*x*y + 1*x*y**2 + 1*x**2 + 1*x**2*y + 1/3*x**3

5.4 Euclidian division

div <serie> f, <serie> g, <object identifier> q, <object identifier> r

It computes the quotient and the remainder of f divided by g such that f = q * g + r and degree(r) < degree(g). The quotient is stored in q and the remainder in r.

Example:
> div(x^3 + x + 1, x^2 + x + 1, q, r);
> q(x) =
    - 1
    + 1*x

> r(x) =
    2
    + 1*x

5.5 Selection

5.5.1 coef_ext

coef_ext <serie>, (<variable> , <integer> ),...)

Return the coefficient of the variable raised to the specified value from the serie.
coef_ext(S, (X, n), (Y, m)) returns the coefficient of X^n*Y^m in the serie S.
Example:
> S= (1+x+y+z)**4 $
> \text{coef\_ext}(S,(x,1),(y,2));
12 + 12*z

5.6 Evaluation

5.6.1 coef\_num

coe\_num  [Function]

coe\_num(<serie> , (<variable> ,<constant> ),...)
Substitute in a serie one (or more) variable(s) by one (or more) numerical value(s).

Example:
> S= (1+x+y+z)**4 $
> \text{coef\_num}(S,(x,0.1),(y,2));
923521/10000 + 29791/250*z + 2883/50*z**2 + 62/5*z**3 + 1*z**4

5.6.2 eval\_num

eval\_num  [Function]
eval\_num(<serie> , {REAL/COMPLEX} , (<variable> ,<num. vec.> ),...)
Evaluates the serie with substituting all variables by their value in the associated numerical vectors.

The function returns a numerical vector of real numbers if REAL is specified or else a numerical vector of complex numbers COMPLEX is specified.

All numerical vectors must have the same size.

Example:
> serie=sin(x+y)-2*y;
serie(y,_EXy1,_EXx1) =
( -0+i* 0.5)*_EXy1**-1*_EXx1**-1
+ ( 0-1* 0.5)*_EXy1*_EXx1
- 2*y

> TABX=0,pi,pi/6;
TABX double precision real vector : number of elements =7
> TABY=-pi,0,pi/6;
TABY double precision real vector : number of elements =7
> TABRES=eval\_num(serie,REAL,(x,TABX),(y,TABY));
TABRES double precision real vector : number of elements =7
> writes(TABRES);
+6.2831853071795862E+00
+4.3699623521985504E+00
+3.227648010019526E+00
+3.1415926535897931E+00
+2.960604205061776341E+00
+1.913229549810371E+00
+1.2246467991473532E+00
> writes(TABX);
+0.0000000000000000E+00
+5.23598775559829882E-01
+1.0471975511965976E+00
+1.5707963267948966E+00
+2.0943951023931953E+00
+2.6179938779914940E+00
+3.1415926535897931E+00

> TABRES=evalnum(serie,COMPLEX,(x,TABX),(y,TABY));

TABRES Double precision complex vector : number of elements =7

> writes(TABRES);
+6.2831853071795862E+00 +1.4997597826618576E-32
+4.3699623521985504E+00 +0.0000000000000000E+00
+3.3227648010019526E+00 +0.0000000000000000E+00
+3.1415926535897931E+00 +0.0000000000000000E+00
+2.9604205061776341E+00 -5.5511151231257827E-17
+1.913229549810371E+00 +5.5511151231257827E-17
+1.2246467991473532E-16 +0.0000000000000000E+00

>
6 Constants

6.1 Standard functions
Most of the functions are described in (see Chapter 10 [Vecteurs numeriques], page 55).

6.1.1 factorial

fac

    fac( <integer> n);

    Return $n!$ (factorial function).

Example :

    > fac(3);
    6

    > n=5;
    n = 5

    > fac(n+1);
    720

6.2 Input/output on real numbers
These functions perform the sequential writing and reading of a text file containing only real values.

ecriture

    ecriture(<filename>);

    Open a file in writing mode in the folder specified by _path. If the file doesn’t exist, it will be automatically created.

Example :

    > ecriture("fichier1.dat");

lecture

    lecture(<filename>);

    Open a file in reading mode in the folder specified by _path.

Example :

    > lecture("fichier1.dat");

print

    print(<real>);

    Write a double-precision floating-point number in the file opened (with the command ecriture).

Example :

    > ecriture("fichier1.dat");

    > print(atan(1)); /* on écrit pi/4 dans fichier1.dat */

read

    read;

    Read a floating-point number in the file opened (with the command lecture) and return a double-precision floating-point number.
Example:

```plaintext
> ecriture("fichier1.dat");
> print(atan(1));
> close;
> lecture("fichier1.dat");
> s=read;
> s = 0.785398163397448
> close;
```

**close**

[Procedure]

Close the file of real numbers if it has been opened (for reading or writing).

Example:

```plaintext
> ecriture("fichier1.dat");
> print(atan(1));
> close;
```
7 Strings

7.1 Declaration and assignment

Character string’s declaration is implicit. It will be realized when assignment are performed. To produce double-quote (") in a string, you must double it.

\[ \text{Example:} \]
\[ /* L'exemple suivant declare les chaines ch et ch2. */ \]
\[ > \text{ch}="file"; \]
\[ \text{ch} = "file" \]
\[ > \text{ch2} = \text{ch}".txt"; \]
\[ \text{ch2} = "file.txt" \]

7.2 Concatenation

\[ \text{Example:} \]
\[ > \text{ch} = "file" + "1"; \]
\[ \text{ch} = "file1" \]
\[ > \text{sch} = "./" + \text{ch}; \]
\[ \text{sch} = "./file1" \]
\[ > \text{sch1} = \text{sch} + "." + \text{str(12)}; \]
\[ \text{sch1} = "./file1.12" \]

7.3 Repetition

\[ \text{Example:} \]
\[ > \text{s}=" %g"; \]
\[ \text{s} = " %g" \]
\[ > \text{format}=4*\text{s}+"\n"; \]
\[ \text{format} = " %g %g %g %g\n" \]
\[ > \text{t}=1,10$ \]
\[ > \text{writes}()\text{format}, \text{t},\text{t}**2, \text{t}**3, \text{t}**4); \]

7.4 Extraction

\[ \text{Example:} \]
\[ > \text{<string> [ <integer> j ]} \]
\[ \text{Return the value of the element specified by the index j. The index starts at 1.} \]
\[ > \text{<string> [ <integer> binf : <integer> bsup ]}; \]
\[ \text{Return an array containing only the elements located between the lower and uper bounds with the specified step.} \]
If the lower bound is omitted then its value is assumed to be 1. If the upper bound is omitted then its value is assumed to be the length of the string.

Remarks: any combination of omission are permitted.

Example:
> s="bonjour";
s = "bonjour"
> s1=s[4];
s1 = "j"
> s2=s[:3];
s2 = "bon"
> s3=s[4:5];
s3 = "jo"

7.5 Comparaison

== [Operator]
<string> == <string>

The Operator == returns true if the strings of characters are equal else it returns false.

Example:
> s="monchemin";
s = "monchemin"
> if (s=="monchemin") then { msg "true"; } else { msg "false"; };
true
>

!= [Operator]
<string> != <string>

The Operator != returns false if the strings of characters are equal else it returns true.

Example:
> s="monchemin";
s = "monchemin"
> if (s=="MON") then { msg "true"; } else { msg "false"; };
false
>

< [Operator]
<string> s1 < <string> s2

The Operator < returns true if s1 is less than s2 using the lexicographic order else it returns false.

Example:
> s="abc";
s = "abc"
> if (s<"abd") then { msg "true"; } else { msg "false"; };
true
>

<= [Operator]
<string> s1 <= <string> s2

The Operator <= returns true if s1 is less than or equal to s2 using the lexicographic order else it returns false.
Example:
> s="abc";
s = "abc"
> if (s<="abe") then { msg "true"; } else { msg "false"; };
true
>
> [Operator]
<string> s1 > <string> s2
The Operator > returns true if s1 is greater than s2 using the lexicographic order else it returns false.

Example :
> s="abc";
s = "abc"
> if (s>"abd") then { msg "true"; } else { msg "false"; };
false
>

>= [Operator]
<string> s1 >= <string> s2
The Operator >= returns true if s1 is greater than or equal to s2 using the lexicographic order else it returns false.

Example :
> s="abc";
s = "abc"
> if (s=>"abe") then { msg "true"; } else { msg "false"; };
false
>

7.6 Conversion from an integer, a real or a serie to a string

str [Function]
str( <integer> );
str( <real> );
str( <serie> );
str( <string> format, <real> );

Convert an integer, a real number or an object in a character string. If format is specified, then the integer or real number is converted to this format. The format specification is similar to that for the printf function of the C-language.

The conversion specifiers are
- g,G The number is converted as a real (style [-]dddd.dddd or scientific style [-]d.dddE+-dd if exponent is too large).
- e,E The number is converted as a real (scientific style [-]d.dddE+-dd).
- f,F The number is converted as a real (style [-]d.dddE+-dd).
- d,i The number is converted as an integer. If the argument is a real number, only its integer part is converted. If the numerical mode NUMRAT or NUMRATMP is selected, then the rational numbers as written as "numerator/denominator".
The length modifier aren’t supported. For example, the format "%lg" will be rejected and produce an error.

If the numerical mode NUMRAT or NUMRATMP is selected, then the rational numbers as written as "numerator/denominator" if no format are specified.

Example :
> ch = str(1235);
ch = "1235"
> ch = str(int(2*pi));
ch = "6"
> s = str(2E3);
s = "2000"
> s = str("%.4g",pi);
s = "3.142"
>_modenum = NUMRATMP;
_modenum = NUMRATMP
> s = str("%d", 3/2);
s = "3/2"
> s = str("%g", 3/2);
s = "1.5"
> s = str(1+x**2);
s = " 1 + 1*x**2"

7.7 Conversion from a list of integers or reals to a string

msg
msg(<string> textformat, <real> x, ... );
Create a string using the format with the real constants.
The format is the same as the command printf in language C (see Section 7.6 [str], page 33, for the valid conversion specifiers). This format must be a string and could be on several lines.
To create a double-quotes, two double-quotes must be used.

Example :
> ch = msg("pi=%.8E",pi,pi);
ch = "pi=3.14159265E+00"
>_modenum=NUMRATMP;
_modenum = NUMRATMP
> s=msg("%d %g", 1/11, 1/11);
s = "1/11 0.0909091"

7.8 Conversion from a string to an integer or a real

dto_num
coef_num
coef_num( <string> );
It converts a character string to an integer or a real number.

Example :
Chapter 7: Strings

7.9 Length of a string

```plaintext
size( <string> );
```

it computes the length of the string, that is to compute the number of characters.

Example:
```plaintext
> ch = "1235";
ch = "1235"
> size(ch);
4
```
8 Arrays

TRIP has 4 types of arrays:
- arrays of series and constants.
- array of variables.
- numerical vectors.
- numerical matrices.

Matrix are arrays of series and constants. For the numerical vectors or matrices, see the appropriate chapter.

8.1 Declaration of array of series

```plaintext
dim

dim <name> [ <list of dimension> ], ...;
Define one or more array of series with the specified dimensions.
Each dimension is separated with a comma. A dimension is defined with two integers, separated by the character : which specify the index of the first element, the index of the last element, and the step for this dimension. If the third integer is missing, the step is 1. The dimensions may stored in an object identifier.
This type of array can contain series, constants, truncatures, strings of characters... but no variables.

Example :
> // It defines the array tl of series or numbers
> // with two dimensions
> dim tl [1:22,-2:6];
> // it defines 3 arrays t1, t2 , t3 with different bounds
> bounds1 = 1:4;
> bounds1 = bounds [ 1:4 ]
> dim t1[bounds1], t2[5:6], t3[-1:3];
> // it defines an array with a step not equal to 1
> dim t5[1:5:2];
> t5[::]=7;
> afftab(t5);
t5[1] = 7
t5[3] = 7
t5[5] = 7
```

8.2 Initialization of an array of series

```plaintext
<name> = [ <serie> , ... : <serie> , ...];
<name> = dim[<serie> , ... : <serie> , ...];
Create and initialize an array of series with the specified series and constants.
The symbol : specifies a new line and the symbol , separates the columns. It must have the same number of columns on each line.

Example :
> // define a 2-dimensional array which contains polynomials
> tab=[1,2+2*x:(1+x)**2,(2+2*x)**2];
```
tab [1:2, 1:2 ] number of elements = 4

> stat(tab);
  Array of series
  tab [ 1:2 , 1:2 ]
  list of array's elements :
  tab [ 1 , 1 ] =
    constante has name tab and has the value  1
  tab [ 1 , 2 ] =
    serie tab ( x )
    number of variables : 1  size of the descriptor : 80 bytes
    number of terms : 2  size : 304 bytes
  tab [ 2 , 1 ] =
    serie tab ( x )
    number of variables : 1  size of the descriptor : 80 bytes
    number of terms : 3  size : 352 bytes
  tab [ 2 , 2 ] =
    serie tab ( x )
    number of variables : 1  size of the descriptor : 80 bytes
    number of terms : 3  size : 352 bytes

> // define a 2-dimensional array which contains numbers
> tab2=[1,2,3:4,5,6];

  tab2 [1:2, 1:3 ] number of elements = 6

>

8.3 Declaration of array of variables

\textbf{dimvar} \hspace{1cm} [\textbf{Procedure}]
\texttt{dimvar <name> [ <list of dimension> ], ...;}

Define one or more array of variables with the specified dimensions.

Each dimension is separated with a comma. A dimension is defined with two integers which specify the index of the first element and the index of the last element for this dimension.

To assign a existing variable to an array of variables, you should the symbol \texttt{:=} instead of \texttt{=}.

Example :
> // define the array of variables t2 with one dimension
> dimvar t2[1:5];
> t2[1] := x;
>  t2[1] = x = 1*x

> // we obtain the derivative of S by x.
> S=1+3*x$
> deriv(S,t2[1]);
> 3
>
> // we define 2 arrays of variables tv1 and tv2
> dimvar tv1[1:2], tv2[1:3];
>
8.4 Initialization of an array of variables

\[ \text{<name> = dimvar[<serie>, ... : <serie>, ...];} \]

Create and intialize an array of variables with the specified variables.

The symbol : specifies a new line and the symbol , separates the columns. It must have the same number of columns on each line.

Example :
> // define the array of variables t2 with the variables x, y, z and u
> i+x+y+z$
> t2 = dimvar[x:y:z:u];

\[ t2[1:4 \text{ ] number of elements = 4} \]

8.5 Generation of an array of variables

[tabvar] [Procedure]

\[ \text{tabvar(<array of variables> );} \]

Generate automatically the variables inside the specified array. The name of variables use the name of the array as radical and its indice.

Example :
> dimvar t[0:3];
> tabvar(t);
> afftab(t);
\[ t[0] = t_0 = 1 \times t_0 \]
\[ t[1] = t_1 = 1 \times t_1 \]
\[ t[2] = t_2 = 1 \times t_2 \]
\[ t[3] = t_3 = 1 \times t_3 \]

8.6 Assignment to an array

[Operator]

\[ \text{= <name> = <array> ;} \]

Assign the array (operation between arrays) to the specified object identifier.

\[ \text{<array> [<integer>, ...] = <operation> ;} \]

Assign a value to an element of an array. This operation can be a serie, a constant, a numerical vector, a truncature but not an array.

\[ \text{<array> [ <integer> binf : <integer> bsup : <integer> pas, ...] = <operation> ;} \]

Assign a value to a part of an array.

If the operation is a serie, a constant, a numerical vector, a truncature then the value is copied to each element of this part of array. If the operation is an array then each element of this array will be assigned to the corresponding element. In this case, it verifies that the number of elements and dimension is compatible.
If the lower bound is omitted then its value is assumed to be 1. If the upper bound is omitted then its value is assumed to be the size of the array. If the step is omitted then its value is assumed to be 1.

When the number of dimension is unknown or variable, the variadic statement ... could be provided as the last argument in order to specify an undefined number of ;;;;

Remarks : any combination of omission are permitted.

Example :
> _affc=1$
> dim t[1:4,7:25];
> t[1,7]=1+x;
t[1,7] = 1
+ 1*x

> r=t;

r [1:4, 7:25 ] number of elements = 76

> dim t[1:4,7:25];
> dim t2[1:4];
> t2[:]=5;
> t[::8]=t2;
> t[:,:,:2]=1+x;
> t[2,...]=3;

:=
<array> [...] := <variable> ;

Assign a variable to an array of variables.

Example :
> dimvar X[1:4];
> X[1]:=crevar("e",1,1);
X[1] = e_1_1 = 1*e_1_1
> deriv(1+2*e_1_1,X[1]);
2

8.7 Size of an array

size [Function]

size(<array> );
size(<array> ,<integer> );

Return the number of elements in the first dimension of the array or in the specified dimension by the integer.

If the specified dimension doesn’t exist in the array then the function returns -1.

Example :
> dim t[1:4,7:25];
> size(t,2);
19
> size(t);
4
>

**num_dim**

`num_dim(<array> t)`

It returns the number of dimensions of the array `t`.

Example:
> dim t1[1:5];
> size(t1);
5
> dim t2[1:22,-2:6];
> size(t2);
22
>

### 8.8 bounds of an array

**inf**

`inf(<array>,<integer>)`

Return the lower bound of the array of the specified dimension by the integer.

If the specified dimension doesn’t exist in the array then the function returns -1.

Example:
> dim t[1:2,0:3];
> inf(t,2);
0
> for k=inf(t,2) to sup(t,2) { t[:,k]= k$ };
> afftab(t);

```
 0  1  2  3  
```

**sup**

`sup(<array>,<integer>)`

Return the upper bound of the array of the specified dimension by the integer.

If the specified dimension doesn’t exist in the array then the function returns -1.

Example:
> dim t[1:2,0:3];
> sup(t,2);
3
> for k=inf(t,1) to sup(t,1) { t[k,]= k$ };
> afftab(t);

```
 1  1  
```
8.9 Display an array

afftab

afftab(<array>);

Display the content of the arrays of series or variables.

Example:
> _affc=1$
> dim t[1:4];
> t[1]=1+x$
> t[2]=2*i$
> t[3]=((x,y),2))$
> afftab(t);
> 
> t[1] = 1
+ 1*x

> t[2] = (0+i*2)
> t[3] = ((y, x), 2))
> t[4] =
>

8.10 Extract an element

[]

Return the value of the element specified by the indexes.

<num. vec.> [ <integer> binf : <integer> bsup : <integer> pas,... ];
<num. vec.> [ <integer> binf : <integer> bsup, ... ];

Return an array containing only the elements located between the lower and upper bounds with the specified step.

If the lower bound is omitted then its value is assumed to be the lower bound of that dimension of this array. If the upper bound is omitted then its value is assumed to be the upper bound of that dimension of this array. If the step is omitted then its value is assumed to be 1.

When the number of dimension is unknown or variable, the variadic statement ... could be provided as the last argument in order to specify an undefined number of ;;;;.

Remarks: any combination of omission are permitted.

Example:
> _affc=1$
> dim t[1:4];
> t[1]=1+x$
> s=t[1];
> s(x) =
Chapter 8: Arrays

1
+ 1*x

> t1=t[3:];

t1 [3:4 ] number of elements = 2

> dim t[1:4,5:10];
> v2=t[:,6:];

v2 [1:4, 6:10 ] number of elements = 20

> v3=t[4,...];

v3 [5:10 ] number of elements = 6

select

select ( <condition> , <array> );

Return an array which only contains the elements of the array where the condition is true.

If the condition is always false, the function returns the constant number 0. To test if the constant value 0 was returned, the following code could be used: if (size(result)==0) then { ... } else { ... }

The condition and the array must have the same number of elements and only one dimension size.

Example :

> // extract from tm the elements such that t1[j]=1
> _modenum=NUMRATMP;
>_modenum = NUMRATMP
> dim t1[1:5], tm[1:5]$;
> for j=1 to 5 { t1[j]=abs(3-j)$ tm[j]=(1+x)**j$ }
> q = select(t1==1, tm);

q [1:2 ] number of elements = 2

> if (size(q)!=0) then { afftab(q); }
q[1] = 1
+ 2*x
+ 1*x**2

q[2] = 1
+ 4*x
+ 6*x**2
+ 4*x**3
+ 1*x**4

>
8.11 Operations on array of series

Most of the functions are described in (see Chapter 10 [Vecteurs numeriques], page 55). The following functions could be applied on the arrays of series (as on the series) to avoid the loops for:

- `coef_ext` see Section 5.5.1 [coef_ext], page 26,
- `coef_num` see Section 5.6.1 [coef_num], page 27,
- `deriv` see Section 5.3.1 [deriv], page 26,
- `integ` see Section 5.3.2 [integ], page 26,

8.11.1 Matrix product

`&*` [Operator]

<array> &* <array>

Compute the matrix product of the arrays (with 1 or 2 dimensions).

Compute \( r = a &* b \) such that \( r[i,j] = \sum_{k=1}^{\text{size}(a,1)} a[i,k] \times b[k,j] \)

If the array has only one dimension then it’s considered as a column-vector.

Remark : the number of columns of the first array must be same as the number of lines of the second array.

Example:

\[
>_ \text{affc}=1$ \text{affdist}=0$
>_ t=[x,y;x-y,x+y]$
>_ t2=[x+y,x-y;x,y]$
>_ r=t&*t2;

\[
\begin{align*}
  r[1:2, 1:2] \text{ number of elements } &= 4 \\
  r[1,1] &= 2*y*x + x**2 \\
  r[1,2] &= y**2 - y*x + x**2 \\
  r[2,1] &= - y**2 + y*x + 2*x**2 \\
  r[2,2] &= 2*y**2 - y*x + x**2 \\
\end{align*}
\]

8.11.2 Matrix determinant

`det` [Function]

\[
\text{det( <array> M )}
\text{det( <array> M, <string> method )}
\]

Computes the determinant of a matrix. The array must have two dimensions.

- If \( M \) is a matrix with numerical entries, the LU algorithm is used in all numerical precision. When computations are performed in double precision, the Lapack Library is used.
- If \( M \) is a matrix with polynomial entries, un algorithme de Bareiss est utilisé par défaut.

If `method` is specified, it uses that algorithm to compute the determinant if the content of the matrix allows it. In other case, the determinant is computed using the default algorithm.

The available values for `method` are:

- "minor" : Determinant expansion by minors using large memory. In practice, it could be used only for the matrices’size less than 12.
Chapter 8: Arrays

- "bareiss": Gauss-Bareiss algorithm.
- "interpol": Interpolation algorithm. It could be only used on the matrices with polynomial entries. Their coefficients must be integer or rational numbers.
- "interpol_modular": Interpolation algorithm using the modular arithmetic and the Chinese remainder. It could be only used on the matrices with polynomial entries. Their coefficients must be integer numbers.

Example:
> \( t = \begin{bmatrix} 9, 0, 7 \\ 1, 2, 3 \\ 4, 5, 6 \end{bmatrix} \) number of elements = 9

> det(t); -48

8.11.3 Matrix inversion

**-1 [Function]

<array> **-1

Computes the inverse of a matrix. The array must have two dimensions.
Remarks: When computations are performed in double precision, the Lapack Library is used. The LU algorithm is used in all numerical precision.

Example:
> _modenum=NUMRATMP$
> t = \begin{bmatrix} 9, 0, 7 \\ 1, 2, 3 \\ 4, 5, 6 \end{bmatrix} \) number of elements = 9

> r = t ** -1;

r [1:3, 1:3 ] number of elements = 9

> afftab(r);
r[1,1] = 1/16 
r[1,2] = -35/48 
r[1,3] = 7/24 
r[2,1] = -1/8 
r[2,2] = -13/24 
r[2,3] = 5/12 
r[3,1] = 1/16 
r[3,2] = 15/16 
r[3,3] = -3/8

8.11.4 Eigenvalues

eigenvalues [Function]
eigenvalues(<array> )
Compute the eigenvalues of the 2-dimensional array (square matrix) using a QR algorithm (lapack library).

Example:
> t=[9,0,7
:1,2,3
:4,5,6];

t [1:3, 1:3 ] number of elements = 9

> l=eigenvalues(t);

l [1:3 ] number of elements = 3

> afftab(l);
1[1] = 13.76915041895731
1[2] = 4.084362034809442
1[3] = -0.8535124537667539

8.11.5 Eigenvectors

eigenvectors

eigenvectors(<array> MAT, <array> TVECT, <array> TVAL)
eigenvectors(<array> MAT, <array> TVECT)

Compute the eigenvectors of the 2-dimensional array MAT (square matrix) using a QR algorithm (lapack library). It stores the eigenvectors in the array TVECT and the eigenvalues in the array TVAL.

Example:
> t=[9,0,7
:1,2,3
:4,5,6];

t [1:3, 1:3 ] number of elements = 9

> eigenvectors(t,vectp,valp);
> afftab(vectp);
vectp[1,1] = -0.8081690381494434
vectp[1,2] = -0.7505769919464202
vectp[1,3] = -0.4846638559537327
vectp[2,1] = -0.209021306608322
vectp[2,2] = 0.3985224414326275
vectp[2,3] = -0.5474094124384613
vectp[3,1] = -0.550611386696964
vectp[3,2] = 0.5270806796287867
vectp[3,3] = 0.6822344772186747
> afftab(valp);
valp[1] = 13.76915041895731
valp[2] = 4.084362034809442
valp[3] = -0.8535124537667539
> // first vector
> p=vectp[:,1]$
> afftab(p);
8.11.6 Arithmetic

The arrays must have the same number of elements on each dimension and the number of
dimension must be the same.

+  
<array> + <array>  
Add term by term two arrays of series.

+  
<array> + <serie>  
<serie> + <array>  
Add a serie with each element of the array of series.

*  
<array> * <array>  
Multiply term by term two arrays of series.

*  
<array> * <serie>  
<serie> * <array>  
Multiply a serie with each element of the array of series.

-  
<array> - <array>  
Subtract term by term two arrays of series.

-  
<array> - <serie>  
<serie> - <array>  
Subtract a serie with each element of the array of series.

/  
<array> / <array>  
Divide term by term two arrays of series.

/  
<array> / <serie>  
<serie> / <array>  
Divide a serie with each element of the array of series.

Example :
> .affc=1$
> t2=x+y,x-y:x,y];

t2 [1:2, 1:2 ] number of elements = 4
> t=[x,y:x+y,x+y];
t [1:2, 1:2 ] number of elements = 4

> r=t*t2+i-t*x;

r [1:2, 1:2 ] number of elements = 4

> afftab(r);
r[1,1] = (0+i*1)*x*y + (-1+i*1)*x**2

r[1,2] = (-0-i*1)*y**2 + (-1+i*1)*x*y

r[2,1] = (1-i*1)*x*y + (-1+i*1)*x**2

r[2,2] = (0+i*1)*y**2 + (-1+i*1)*x*y - 1*x**2

>

8.12 Conversion

For the conversion to numerical vectors, (see Section 10.12 [Conversion], page 94).
9 Structures and OOP

The data structures contain other values indexed by their names. These fields may have a global or private visibility. The macros can be associated to these data structures. These macros may access to all the fields.

9.1 Type declaration

```
struct <structtype> { <list of parameters> ; private <list of parameters> ; };
struct <structtype> { <object identifier> = <value> ; ... private <object identifier> = <value> ; private ... };
```

It defines a new type of records whose fields of the first list can be accessed globally and the other list which follows the keyword `private` can only be accessed by an associated macro.

The fields may be initialized to a default value.

Example :

```c
> // declaration of the type POINT
> struct POINT {
    x, y;
};
>
> // declaration of the type MyData
> struct MyData {
    x = 0;
    y = 0;
    z = -1;
    vnumr v([1:3]);
    dim armap[1:5];
    private name = "NONAME";
    private file = "/tmp/myfile";
};
> MyData a;
> afftab(a);
  a@x = 0
  a@y = 0
  a@z = -1
  a@v = a double precision real vector : number of elements =3
  a@armap = a [1:5] number of elements = 5
  a@name = "NONAME"
  a@file = "/tmp/myfile"
```

9.2 Object identifier declaration

```
struct <structtype> <object identifier> p1, <object identifier> p2, ... ;
```

It creates the object identifiers whose the name are given in the list. These objects are data structure of the specified type.
Example:
> // declaration of the type POINT
> struct POINT {
>   x, y;
>};
> // declaration of the object of type POINT
> struct POINT p1, p2, p3;
> p1;
p1 = structure POINT
> afftab(p1);
p1@x = /* UNINITIALIZED */
p1@y = /* UNINITIALIZED */

9.3 Access to member attributes

@ [Operator]

<structure> @<name>

It allows to access to the value of a field of the data structure. If this field is private, this field can only be accessed from a macro associated to this structure.

Example:
> struct MyData {
>   x,y,z;
>   name, file;
>};
> struct MyData s;
> s@x = 3;
s@x = 3
> s@y = 1;
s@y = 1
> s@z = s@x + 2*s@y;
s@z = 5
> s@name = "temp001";
s@name = "temp001"
s@file=file_open(s@name,write);
s@file = file "temp001" opened in write mode

9.4 Display

afftab [Procedure]

afftab(<structure> x);

It displays the value of all the fields of the data structure x.

Example:
> struct MyData {
>   x,y,z;
>   name, file;
>};
> struct MyData s;
Chapter 9: Structures and OOP

9.5 Constructor macros

macro [Procedure]
macro <structtype> @ <structtype> { <body> }

It defines a constructor of the data structure with no parameter and a body of trip code. This macro is implicitly called when an object is initialized.

The macro allows to access to all the private or public fields of the data structure. A direct access to these fields is allowed and does not require to be prefixed by @.

The object self is already defined during the execution of the macro. It refers to the value of the data structure which responsible of the call. This object is useful for a global assignment or to call another macro with this value.

Example:
> struct MyData {
  t;
  vnumR v([1:3])$
};
> macro MyData@MyData {
  private _ALL;
  t = log(2)
  v[:]=4$
};
> MyData a;
> afftab(a);
a@t = 0.6931471805599453
a@v = a double precision real vector : number of elements =3
>

9.6 Member macros

9.6.1 Declaration

macro [Procedure]
macro <structtype> @ <name> [ <list of parameters> ] { <body> }
macro <structtype> @ <name> { <body> }
private macro <structure> @ <name> [ <list of parameters> ] { <body> }
private macro <structure> @ <name> { <body> };

It defines a member macro of the data structure with 0 or more parameter and a body of trip code.

The macro allows to access to all the private or public fields of the data structure. A direct access to these fields is allowed and does not require to be prefixed by @.

The object self is already defined during the execution of the macro. It refers to the value of the data structure which responsible of the call. This object is useful for a global assignment or to call another macro with this value.

If the macro is defined as private, then this macro can be called only other macros associated to the same data structure.

Example:
> struct MyData {
  x,y,z;
  private name, file;
};
>
> // Define Init as a macro of MyData
> macro MyData@Init[vx,vy,vz,vname] {
  x = vx$
  y = vy$
  z = vz$
  name = vname$
  %OpenFile$
  afftab(self);
};
>
> // Define OpenFile as a private macro of MyData
> private macro MyData@OpenFile {
  file = file_open(name,write)$
};
>

9.6.2 Execution

% [opérateur]

<structure> % <name> [ list of parameters ];
<structure> % <name> ;

It calls a member macro of the data structure and defines the object self as the value of the data structure during its execution.

Example:
> struct MyData {
  x,y,z;
  private name, file;
};
>
> macro MyData@Init[vx,vy,vz,vname] {
  x = vx$
  y = vy$
  z = vz$
name = vname$
%OpenFile$
afftab(self);
};
> macro MyData@Clear {
%CloseFile;
};
> private macro MyData@OpenFile {
  file = file_open(name,write)$
};
> private macro MyData@CloseFile {
  file_close(file);
};
> struct MyData s;
> // execute the macro Init
> s%Init[1,2,3,"temp001"];
s@x = 1
s@y = 2
s@z = 3
s@name = "temp001"
s@file = file "temp001" opened in write mode
>
> // execute the macro Clear
> s%Clear;
>
10 Numerical vectors

The numerical values stored in vectors are always real or complex double-precision, quadruple-precision or multiprecision numbers depending on the current numerical mode. A numerical vector is assumed to be a column vector.

In the numerical mode NUMRATMP, the rational numbers may be stored in the numerical vector of rational numbers.

10.1 Declaration

Explicit declaration for numerical vectors are only required before call to read, readbin (see Section 10.8 [Entree/SortieTabNum], page 63) and resize. Explicit declaration for array of numerical vectors are always required.

10.1.1 vnumR

```
vnumR <name> , ... ;
vnumR <name> [ <dimension of an array> ] , ... ;
```

It declares a real vector or an array of real vectors.

The real vector’s size is dynamic. After the declaration, the real vectors are empty. To set their sizes, the command resize must be used.

The dimension specifies the number of elements in the array of real vectors.

```
vnumR <name> ([ 1: <integer> n ]), ... ;
vnumR <name> [ <dimension of an array> ] ([ 1: <integer> n ]), ... ;
```

It declares a vector of \( n \) reals or an array of vectors of \( n \) reals.

Example:
```
> vnumR A, C([1:5]);
> stat(A);
Numerical vector A contains 0 double precision reals.
> stat(C);
Numerical vector C contains 5 double precision reals.
Size of the array in bytes: 40
> bounds=1:2;
bounds = bounds [ 1:2 ]
> vnumR TE[bounds], T0[bounds]([1:6]);
> stat(TE);
Array of series
TE [ 1:2 ]
list of array’s elements :
TE [ 1 ] =
Numerical vector TE contains 0 double precision reals.
TE [ 2 ] =
Numerical vector TE contains 0 double precision reals.
> stat(T0);
Array of series
T0 [ 1:2 ]
list of array’s elements :
```
T0[1] =
Numerical vector T0 contains 6 double precision reals.
Size of the array in bytes: 48
T0[2] =
Numerical vector T0 contains 6 double precision reals.
Size of the array in bytes: 48

10.1.2 vnumC

vnumC
vnumC <name> , ... ;
vnumC <name> [ <dimension of an array> ] , ... ;
It declares a complex vector or an array of complex vectors.
The complex vector’s size is dynamic. After the declaration, the complex vectors are empty.
To set their sizes, the command resize must be used.
The dimension specifies the number of elements in the array of complex vectors.

Example:
> vnumC A, C([1:5]);
> stat(A);
Numerical vector A contains 0 double precision complex.
> stat(C);
Numerical vector C contains 5 double precision complex.
Size of the array in bytes: 80
> vnumC TE[1:2], T0[1:2]([1:6]);
> stat(TE);
Array of series
TE [1:2]
list of array’s elements:
TE[1] =
Numerical vector TE contains 0 double precision complex.
TE[2] =
Numerical vector TE contains 0 double precision complex.
> stat(T0);
Array of series
T0 [1:2]
list of array’s elements:
T0[1] =
Numerical vector T0 contains 6 double precision complex.
Size of the array in bytes: 96
T0[2] =
Numerical vector T0 contains 6 double precision complex.
Size of the array in bytes: 96
10.1.3 vnumQ

vnumQ [Procedure]

vnumQ <name> , ... ;

vnumQ <name> [ <dimension of an array> ] , ... ;

It declares a vector of rational numbers or an array of vector of rational numbers.

The rational vector’s size is dynamic. After the declaration, the rational vectors are empty.

To set their sizes, the command resize must be used.

The dimension specifies the number of elements in the array of rational vectors.

vnumQ [Procedure]

vnumQ <name> ([ 1: <integer> n ]), ... ;

vnumQ <name> [ <dimension of an array> ] ([ 1: <integer> n ]), ... ;

It declares a vector of n rational numbers or an array of vectors of n rational numbers.

After the declaration, the elements of the vector are initialized with the value 0.

Example:

> _modenum=NUMRATMP;

> vnumQ A, C([1:5]);

> stat(A);

Numerical vector A contains 0 rational numbers.

> stat(C);

Numerical vector C contains 5 rational numbers.

Size of the array in bytes: 160

> vnumQ TE[1:2], T0[1:2]([1:6]);

> stat(TE);

Array of series

TE [ 1:2 ]
list of array’s elements:

TE [ 1 ] =

Numerical vector TE contains 0 rational numbers.

TE [ 2 ] =

Numerical vector TE contains 0 rational numbers.

> stat(T0);

Array of series

T0 [ 1:2 ]
list of array’s elements:

T0 [ 1 ] =

Numerical vector T0 contains 6 rational numbers.

Size of the array in bytes: 192

T0 [ 2 ] =

Numerical vector T0 contains 6 rational numbers.

Size of the array in bytes: 192

>

10.2 Initialization

... [Procedure]

<name> = <real> binf , <real> bsup , <real> step ;
It declares and initializes a real vectors such that all elements are initialized by a loop (from $b_{inf}$ to $b_{sup}$ with the step $b_{step}$):

$$\text{for } (j=1, \text{valeur} = b_{inf}) \text{ to } \text{valeur} \leq b_{sup} \text{ step } (j=j+1, \text{valeur} = \text{valeur} + b_{step}) \text{ <name> } [j]=\text{valeur}$$

The step $b_{step}$ is optional; By default, its value is 1.

Example:
- $t = 0, 10$
  - $t$ double precision real vector : number of elements = 11
  - writes($t$);
    +0.0000000000000000E+00
    +1.0000000000000000E+00
    +2.0000000000000000E+00
    +3.0000000000000000E+00
    +4.0000000000000000E+00
    +5.0000000000000000E+00
    +6.0000000000000000E+00
    +7.0000000000000000E+00
    +8.0000000000000000E+00
    +9.0000000000000000E+00
    +1.0000000000000001E+01

- $tab = -\pi, 6, \pi/100$
  - $tab$ double precision real vector : number of elements = 291
  - writes([::30], $tab$);
    -3.1415926535897931E+00
    -2.1991148575128552E+00
    -1.2566370614359170E+00
    -3.1415926535897887E-01
    +6.2831853071795907E-01
    +1.5707963267948966E+00
    +2.5132741228718354E+00
    +3.4557519189487733E+00
    +4.3982297150257113E+00
    +5.3407075111026483E+00

Function $vnumR[]$:

$name$ = $vnumR$ [<real> or <real vec.> or <array of real vec.> , ... : <real> ,... ] ; It declares and initializes a real vector or an array of real vectors with the specified reals or real vectors.

The character : separates the lines and the character , separates the columns. All columns must have the same line counts.

Example:
- // declare an array of 3 vectors of 2 reals.
- $tab3$ = $vnumR[1, 2, 3:4, 5, 6]$;

$tab3$ [1:3 ] number of elements = 3

writes($tab3$);
+1.0000000000000000E+00 +2.0000000000000000E+00 +3.0000000000000000E+00
+4.0000000000000000E+00 +5.0000000000000000E+00 +6.0000000000000000E+00
$t = 7, 8$;
- $t$ double precision real vector : number of elements = 2
> tab4=vnumR[t,tab3];

    tab4 [1:4 ] number of elements = 4

> // declare a vector of 5 reals.
> t2=vnumR[1:2:4:8:9];
    t2 double precision real vector : number of elements =5
> writes(t2);
    +1.0000000000000000E+00
    +2.0000000000000000E+00
    +4.0000000000000000E+00
    +8.0000000000000000E+00
    +9.0000000000000000E+00

vnumC[] [Function]

<name> = vnumC[<complex> or <complex vec.> or <array of complex vec.>, ... : <complex> , ...
It declares and initializes a complex vector or an array of complex vectors with the specified
complexs or complex vectors.

The character : separates the lines and the character , separates the columns. All columns
must have the same line counts.

Example :
> // declare a vector of 5 complex numbers
> tab3=vnumC[1:1+i:2+2*i:3+3*i];
    tab3 Double precision complex vector : number of elements =4
> writes(tab3);
    +1.0000000000000000E+00 +0.0000000000000000E+00
    +1.0000000000000000E+00 +1.0000000000000000E+00
    +2.0000000000000000E+00 +2.0000000000000000E+00
    +3.0000000000000000E+00 +3.0000000000000000E+00
> vnumC ti;
> resize(ti,5,3-5*i);
> tab4=vnumC[tab3 : 5+7*i : ti];
    tab4 Double precision complex vector : number of elements =10
> // declare an array of 2 vectors of 2 complex numbers
> z4=vnumC[5,2+i:4,9+3*i];

    z4 [1:2 ] number of elements = 2

> writes("%g %g %g %g\n",z4[1],z4[2]);
      5 0 2 1
      4 0 9 3

10.3 Display

writes [Procedure]

    writes([<integer> : <integer> : <integer> ],<string>,<(array of) num. vec.>,...);
    writes(<string>,<(array of) num. vec.>,...);
    writes(<(array of) num. vec.>,...);
writes( [{\texttt{binf}}:{{\texttt{bsup}}}:\{\texttt{step}\}],  \{\texttt{format}\} <(array of) num. vec.> ,...).

It prints, to the screen, numerical vectors or the array of numerical vectors in columns form.

From the \texttt{binf}th element of each numerical vector (if \texttt{binf} isn’t specified, it starts from the first element).

To the \texttt{bsup}th element of each numerical vector (if \texttt{bsup} isn’t specified, it stops to the last element of the biggest vector).

Every \texttt{step} elements (if \texttt{step} isn’t specified, the step is 1).

\texttt{format} is optional. This format is similar to the C format (cf. printf) and must be between double quotes (\"). To display a double quote (\") as a character, you must double it:

For example, if the C format is " \%g \"titi\" \%g", you must write " \%g\"titi\"\" \%g"

A complex vector will be printed on two columns (the first one for the real part and the second one for the imaginary part).

Example :

Ecriture de tous les éléments de T et de X.
La première colonne correspondra à T.
La deuxième colonne correspondra à la partie réelle de X.
La troisième colonne correspondra à la partie imaginaire de X.

...>
stat(X);
vecteur numérique X de 6 complexes.
   taille en octets du tableau: 96
>stat(T);
vecteur numérique T de 6 réels.
   taille en octets du tableau: 48
>writes(T,X);
+9.999993149794888E-02 +1.00000000456180E-01 +1.095970178673141E-06
-2.000000940350956E-01 +9.999999679956E-03 +1.31200753238499E-07
+3.0000003832689856E-01 +1.00000314882390E-03 -1.403292661135361E-08
-4.000000970924361E-01 +9.999995669530993E-05 +1.695880029074994E-09
+4.99999805039361E-01 +1.000003117384769E-05 +3.216502329016007E-11
-5.999999830866213E-01 +9.999979187109419E-07 -1.796078221829677E-12
>

Ecriture du 2 au 4 éléments de T et de X
avec un format "%.1g\t(%.5g+i*%.5E)n".
> writes([2:4],"%.1g\t(%.5g+i*%.5E)n",T,X);
-0.2  (0.01+i*1.31201e-07)
 0.3  (0.001+i*-1.40329e-08)
-0.4  (0.0001+i*1.69588e-09)

Ecriture du 1 au 5 éléments de T et de X avec un pas de 2 sans format.
> writes([1:5:2],T,X);
+9.99993149794888E-02 +1.00000000456180E-01 +1.095970178673141E-06
+3.0000000832689856E-01 +1.00000314882390E-03 -1.403292661135361E-08
+4.99999805039361E-01 +1.000003117384769E-05 +3.216502329016007E-11

Ecriture des éléments de T et de X avec un pas de 5 sans format.
> writes([::5],T,X);
Chapter 10: Numerical vectors 61

+9.999993149794888E-02 +1.0000000000456180E-01 +1.095970178673141E-06
-5.999999830666213E-01 +9.999979187109419E-07 -1.796078221829677E-12

10.4 Size

size
size(<num. vec.>)

Return the number of elements in the numerical vector.

Example:
> t=1,10;
t double precision real vector : number of elements =10
> size(t);
10
> p=-pi,pi,pi/400;
p double precision real vector : number of elements =800
> size(p);
800
>

10.5 Resize

resize
resize(<(array of) num. vec.>, <integer>);
resize(<(array of) num. vec.>, <integer>, <constant>);

Change the size of the numerical vector or of the array of numerical vectors.

If a constant isn’t specified then all elements will be initialized to 0 else all elements will initialized with this constant.

Example:
> vnumR t;
> resize(t,3);
> vnumR t2;
> resize(t2,3,5);
> writes(t,t2);
+0.0000000000000000E+00 +5.0000000000000000E+00
+0.0000000000000000E+00 +5.0000000000000000E+00
+0.0000000000000000E+00 +5.0000000000000000E+00
> vnumC t[1:3];
> resize(t[2],2,1-5*i);
> writes(t[2]);
+1.0000000000000000E+00 -5.0000000000000000E+00
+1.0000000000000000E+00 -5.0000000000000000E+00
>

10.6 Bounds

inf
inf(<num. vec.>,<integer>);

Return the lower bound of the numerical vector. The supplied integer must be equal to 1.

If the specified dimension is not 1, then the function returns -1.
Example :
> t=3,10;
  t double precision real vector : number of elements =8
> inf(t,1);
  1
>
sup
    sup(<num. vec.>,<integer>);
    Return the upper bound of the numerical vector. The supplied integer must be equal to 1.
    If the specified dimension is not 1, then the function returns -1.
    Example :
    > t=3,10;
    t double precision real vector : number of elements =8
    > sup(t,1);
    8
>
10.7 Data retrieval

10.7.1 select

select
    select ( <condition>, <num. vec.> );
    Return a numerical vector which only contains the elements of the numerical vector where
    the condition is true.
    The condition and the numerical vector must have the same number of elements.
    Example :
    > // return the elements which are multiple of 3
    > t=1,10;
    t double precision real vector : number of elements =10
    > r=select((t mod 3)==0, t);
    r double precision real vector : number of elements =3
    > writes(r);
    +3.0000000000000000E+00
    +6.0000000000000000E+00
    +9.0000000000000000E+00
>
10.7.2 Data retrieval

[]
    <num. vec.> [ <real vec.> ];
    Return a numerical vector which only contains the elements of the numerical vector located
    at the indexes stored in the real vector.
    Remark : The real vector must be an object identifier and not be a operation result.
    Example :
    > r=vnumR[1:5:7:9];
    r double precision real vector : number of elements =4
    > t=20,30;
Chapter 10: Numerical vectors

```plaintext
t double precision real vector : number of elements =11
> b=t[r];
b double precision real vector : number of elements =4
> writes("%g\n",b);
20
24
26
28

[::]
[num. vec.] [ <integer> binf : <integer> bsup : <integer> pas ];
[num. vec.] [ <integer> binf : <integer> bsup ];

Return a numerical vector which contains only the elements located between the lower bound and the upper bound with the specified step.
If the lower bound isn’t specified, then the lower bound is assumed to be 1.
If the upper bound isn’t specified, then the upper bound is assumed to be the size of the numerical vector.
If the step isn’t specified, then the step is assumed to be 1.
Remark : All missing combinations are allowed.

Example :
> t=1,10;
t double precision real vector : number of elements =10
> r=t[::2];
r double precision real vector : number of elements =5
> v=t[7:9];
v double precision real vector : number of elements =3
> y=t[5:10:2];
y double precision real vector : number of elements =3
> writes("%g %g\n",v,y);
7 5
8 7
9 9

10.8 Input/Output

10.8.1 read

read

read(<filename> , [ <integer> : <integer> : <integer> ],

  (array of) num. vec. , ... ,
  (array of) num. vec. , <integer> ), ...
  (array of) num. vec. , <integer> , <integer> ), ...);

equivalent to
read(fichier.dat, [binf:bsup:step], T, (T1,n1), (T3,n2,n3));
read(fichier.dat, T, (T1,n1), (T3,n2,n3));

Read in a text file the specified columns and store them to numerical vectors.
starting at the the line binf (if binf isn’t specified, then it will start at the first line of the file)
```
to the line bsup (if bsup isn’t specified, then it will read to the end of file)
every bstep lines (if bstep isn’t specified, then the step is assumed to 1)

- If the index of the column for a numerical vector of real numbers is specified, then the
column for this vector will be the index of the last column +1.

- If the index of the column for the real part of a numerical vector of complex numbers is
specified, then the column for this vector will be the index of the last column +1.

- If the index of the column for the imaginary part of a numerical vector of complex
numbers isn’t specified and the column for the real part is specified, then the imaginary
part is initialized to 0.

If the file contains the strings NAN or NANQ, then it assumes to be the value "Not A Number". If
the file contains the strings INF, Inf ou Infinity, then it assumes to be the value "infinite".

When an empty or incomplete line is encountered, the behavior of this function depends on
the value of _read (see Section 3.25 [_read], page 19).

Example :
Lecture dans le fichier tessin.out de la ligne 2 à ligne 100
avec un pas de 3.
Le vecteur T contiendra la première colonne,
la partie réelle de X contiendra la deuxième colonne,
la partie imaginaire de X contiendra la troisième colonne,
TAB[1] contiendra la 4eme colonne,
TAB[2] contiendra la 5eme colonne,
> vnumR T;
vnumC X;
vnumR TAB[1:3];
read(tessin.out,[2:100:3],T,X,TAB);
stat(T);
stat(X);
stat(TAB);
T nb elements reels =0
>
X nb elements complexes =0
>
TAB [1:3] nb elements = 3

> > Tableau numerique T de 33 reels.
taille en octets du tableau: 264
> > Tableau numerique X de 33 complexes.
taille en octets du tableau: 528
> > Tableau de series
TAB [ 1:3 ]
liste des elements du tableau :
taille en octets du tableau: 264
taille en octets du tableau: 264
Lecture dans le fichier tessin.out de la ligne 2 à ligne 100.
Le vecteur T contiendra la première colonne,
la partie réelle de X contiendra la 4ème colonne,
la partie imaginaire de X sera nulle,
> read(tessin.out,[2:100],T,(X,4),(TAB[3],5));

Lecture dans le fichier tessin.out de l’ensemble des lignes.
Le vecteur T contiendra la 2ème colonne,
la partie réelle de X contiendra la 3ème colonne,
la partie imaginaire de X contiendra la 4ème colonne,
> read(tessin.out,(T,2),(X,3,4),(TAB[3],5));

Lecture dans le fichier tessin.out à partir de la ligne 1000.
Le vecteur T contiendra la 2ème colonne,
la partie réelle de X contiendra la 3ème colonne,
la partie imaginaire de X contiendra la 4ème colonne,
> read(tessin.out,[1000:],(T,2),(X,3,4),(TAB[3],5));

Lecture dans le fichier tessin.out avec un pas de 50 lignes.
Le vecteur T contiendra la 2ème colonne,
la partie réelle de X contiendra la 3ème colonne,
la partie imaginaire de X contiendra la 4ème colonne,
> read(tessin.out,[::50],(T,2),(X,3,4),TAB[3]);

This function is similar to the above command but allows to specify a format to read the
data. It is able to read columns of numbers or strings of characters. The type of the data is
specified by the format.
The allowed formats are:
- %E double-precision floating-point numbers.
- %e double-precision floating-point numbers.
- %g double-precision floating-point numbers.
- %s string of characters
For the format specifier %s, the object must not be declared before ; the returned object is an array of string.

For the format specifiers %g, %e, %E, the objects must be declared before this command as a (array of) numerical vector (vnumR). The returned object will be of the same type.

By default, it assumes that the columns are separated by spaces or tabulations. If a length is specified in the format (e.g., %10g or %10s for a column of 10 characters), this length specifies the number of characters of the column. In this case, there is no distinction between spaces, tabulations and other characters.

When an empty or incomplete line is encountered, the behavior of this function depends on the value of \_read (see Section 3.25 \_read, page 19).

Limitation : It is not possible to write the following command : \read(file,"....", T1, T2, ...);. It must be rewritten as : fmt=" ...$ \read(file,fmt, T1, T2, ...);

Example :
> // generates a file with 2 columns : asteroid number and name
> f=file_open(data.dat, write)$
> file_writemsg(f, " 1 Ceres \n 2 Pallas\n 3 Juno \n");
> file_close(f);
>
> // reads two columns like a numerical vector and an array of string
> // using a space separator
> vnumR id;
> fmt="%g %s";
fmt = "%g %s"
> read(data.dat, fmt, id, name);
> writes(id);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
> afftab(name);
name[1] = "Ceres"
name[2] = "Pallas"
name[3] = "Juno"
>
> // reads two columns like a numerical vector and an array of string
> // using a fixed size for each column
> vnumR id2;
> fmt="%5g %7s";
fmt = "%5g %7s"
> read(data.dat, fmt, id2, name2);
> writes(id2);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
> afftab(name2);
name2[1] = "Ceres"
name2[2] = "Pallas"
name2[3] = "Juno"

The following example show how to read the astorb data files using the format.

Example :
vnumR number, H, Slope, ColorIndex, v1, v2, v3, v4, v5, v6;
vnumR arc, observations, epochyy, epochymm, epochdd;
vnumR anomaly, perihelion, ascendingnode, Inclination, Eccentricity, Semimajoraxis, Date;

fmt="%7s%19s%16s%6s%5g%5s%6s%6s"+6*"%4g"+2*"%g"+"%4g"
+2*"%2g"+5*"%11g"+"%12g%9s";

read(astorb.dat, [:10000], fmt, number, name, computer, H, Slope, ColorIndex, diameter, Taxonomic, v1, v2, v3, v4, v5, v6, arc, observations, epochyy, epochymm, epochdd, anomaly, perihelion, ascendingnode, Inclination, Eccentricity, Semimajoraxis, Date);

10.8.2 readappend

readappend

readappend(filename, [integer : integer : integer],
(array of) num. vec., ...
(array of) num. vec., <integer>, ...
(array of) num. vec., <integer>, <integer>, ...
);
equivalent to
readappend(fichier.dat, [binf:bsup:step], T, (T1,n1), (T3,n2,n3));
readappend(fichier.dat, T, (T1,n1), (T3,n2,n3));

This function is similar to the function read (see Section 10.8.1 [read], page 63) but it stores read data to the end of the numerical vectors (with an automatic growing step) instead of overwriting their contents.

Example :
> t1=1:5;
t1 Tableau de reels double-precision : nb reels =5
> write(temp001, t1);
> write(temp002, t1**2);
> vnumR w;
> readappend(temp001,w);
> readappend(temp002,w);
> writes(w);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
+4.0000000000000000E+00
+5.0000000000000000E+00
+6.0000000000000000E+00
+7.0000000000000000E+00
+8.0000000000000000E+00
+9.0000000000000000E+00
+1.0000000000000000E+01
+1.6000000000000000E+01
+2.5000000000000000E+01

readappend

readappend(filename, [integer : integer : integer], textfmt,
(array of) num. vec., ...
(array of) num. vec., <integer>, ...
This function is similar to the function read (see Section 10.8.1 [read], page 63) using the format `textfmt`. But it stores read data to the end of the numerical vectors (with an automatic growing step) instead of overwriting their contents.

### 10.8.3 write

```
write( <filename> , [ <integer> : <integer> : <integer> ] , <string> , <(array of) num. vec.> , ...);
write( <filename> , <string> , <(array of) num. vec.> , ...);
write( <filename> , <(array of) num. vec.> , ...);
```

equivalent to

```
write( fichier.dat, [binf:bsup:step], "format", T, T1, T2).
write( fichier.dat, "format", T, T1, T2).
write( fichier.dat, T, T1, T2).
write( fichier.dat, [binf:bsup:step], T, T1, T2).
```

Write to a file the numerical vector or the arrays of numerical vectors as columns of numbers.

- from the element `binf` of each numerical vector (if `binf` isn’t specified, then it’s assumed to be the first one).
- to the element `bsup` of each numerical vector (if `bsup` isn’t specified, then it’s assumed to be the last one of the largest vector).
- every `step` element (if `step` isn’t specified, then it’s assumed to be 1).

The format is optional. The format is the format of C language (See `printf`) and surrounded with double-quotes (""`). To obtain a double-quotes, two double-quotes must be used:

For example: if the C format is "%g \"titi\" %E", it must be written as, "%g \"titi\"\" %E"

A numerical vector of complex numbers uses two columns (the first one for the real part and the second one for the imaginary part).

#### Example:

```c
Ecriture de tous les éléments de T et de X dans le fichier tx.out.
La première colonne correspondra à T.
la deuxième colonne correspondra à la partie réelle de X.
la troisième colonne correspondra à la partie imaginaire de X.
> vnumR T;
vnumC X;
...
stat(T);
stat(X);
```

```c
> Tableau numérique T de 33 réels.
   taille en octets du tableau: 264
> Tableau numérique X de 33 complexes.
   taille en octets du tableau: 528
> write(tx.out,T,X);
>
Ecriture de 10 au 20 éléments de T et de X dans le fichier tx.out avec
```
un format "\%g\t(\%8.4E+i*\%e)\n".
La première colonne correspondra à T.
la deuxième colonne correspondra à la partie réelle de X.
la troisième colonne correspondra à la partie imaginaire de X.
> write(tx.out, [10:20], "\%g\t(\%8.4E+i*\%e)\n", T, X);

Ecriture de 1 au 20 éléments de T et de X avec un pas de 2 dans le fichier tx.out sans format.
La première colonne correspondra à T.
la deuxième colonne correspondra à la partie réelle de X.
la troisième colonne correspondra à la partie imaginaire de X.
> write(tx.out, [1:20:2], T, X);

Ecriture des éléments de T et de X avec un pas de 5 dans le fichier tx.out sans format.
La première colonne correspondra à T.
la deuxième colonne correspondra à la partie réelle de X.
la troisième colonne correspondra à la partie imaginaire de X.
> write(tx.out, [:5], T, X);

10.8.4 readbin

readbin [Procedure]
readbin(<filename> , [ <integer> : <integer> : <integer> ] , <string> , <(array of) real vec.>, ...);
readbin(<filename> , <string> , <(array of) real vec.>, ...);
equivalent to
readbin(fichier.dat, [binf:bsup:step], format, T, T2, T3);
readbin(fichier.dat, format, T, T2, T3);

Read in a binary file the data with the specified format and store them to numerical vectors.
beginning at the record binf (if binf isn’t specified, then it’s assumed to be first record).
to the record bsup (if bsup isn’t specified, then it’s assumed to be the end of file)
every bstep records (if bstep isn’t specified, then it’s assumed to be 1)

The allowed format to define a record are :
- %2d signed integer on 2 bytes.
- %4d signed integer on 4 bytes.
- %8d signed integer on 8 bytes.
- %d signed integer on (default size of the system for an integer).
- %2u unsigned integer on 2 bytes.
- %4u unsigned integer on 4 bytes.
- %8u unsigned integer on 8 bytes.
- %u unsigned integer (default size of the system for an integer).
- %E double-precision floating-point number.
- %e double-precision floating-point number.
- %g double-precision floating-point number.
- %8g double-precision floating-point number on 8 bytes.
- %hE simple precision floating-point number.
The number of format must be the same as the number of vectors. If the provided object identifier is an array of real vectors, then it must have the same number of format as the number of elements in the array.

The data are previously converted if the global variable \_endian (see Chapter 3 \_endian, page 9) isn’t the same as its default value.

Example :
Lecture dans le fichier binaire test1.dat
d’entiers signés sur 4 octets
et stockés dans le vecteur de réels T1.

\begin{verbatim}
  vnumR T1;
  readbin(test1.dat,"%4d",T1);
\end{verbatim}

Lecture dans le fichier binaire test1.dat
des 30 premiers entiers signés sur 4 octets
et stockés dans le vecteur de réels T1.

\begin{verbatim}
  vnumR T1;
  readbin(test1.dat,[:30],"%4d",T1);
\end{verbatim}

Lecture dans le fichier binaire test2.dat
dont chaque enregistrement est composé de 2 réels
et stockés dans le vecteur de réels T1 et T2.

\begin{verbatim}
  vnumR T1,T2;
  readbin(test2.dat,[:30],"%g%g",T1,T2);
\end{verbatim}

Lecture dans le fichier binaire test3.dat
dont chaque enregistrement est composé de 4 réels
et stockés dans le tableau T3.

\begin{verbatim}
  vnumR T3[1:4];
  readbin(test3.dat,"%g%g%g%g",T3);
\end{verbatim}

\begin{verbatim}
est équivalent à

  readbin(test3.dat,"%g%g%g%g",T3[1],T3[2],T3[3],T3[4]);
\end{verbatim}

10.8.5 writebin

\begin{verbatim}
writebin(\texttt{<filename>},[ \texttt{<integer> : <integer> : <integer> }] , \texttt{<string>}, \texttt{<(array of) real vec.>}, ...);
writebin(\texttt{<filename>}, \texttt{<string>}, \texttt{<(array of) real vec.>}, ...);
equivalent to
writebin(fichier.dat, [binf:bsup:step], format, T, T2, T3);
writebin(fichier.dat, format, T, T2, T3);
\end{verbatim}

Write the numerical vectors to a binary file the data with the specified format.

beginning at the index binf (if binf isn’t specified, then it’s assumed to be first element of the numerical vectors).
to the index bsup (if bsup isn’t specified, then it’s assumed to be the size of the largest numerical vector)
every bstep elements (if bstep isn’t specified, then it’s assumed to be 1)

The allowed format to define a record are:
- %2d signed integer on 2 bytes.
- %4d signed integer on 4 bytes.
- %8d signed integer on 8 bytes.
- %d signed integer on (default size of the system for an integer).
- %2u unsigned integer on 2 bytes.
- %4u unsigned integer on 4 bytes.
- %8u unsigned integer on 8 bytes.
- %u unsigned integer (default size of the system for an integer).
- %E double-precision floating-point number.
- %e double-precision floating-point number.
- %g double-precision floating-point number.
- %8g double-precision floating-point number on 8 bytes.
- %hE simple precision floating-point number.
- %he simple precision floating-point number.
- %hg simple precision floating-point number.
- %4g simple precision floating-point number on 4 bytes.

The number of format must be the same as the number of vectors. If the provided object identifier is an array of real vectors, then it must have the same number of format as the number of elements in the array.

If the numerical vectors doesn’t have the same size, then the missing data are completed with 0.

The data are previously converted if the global variable _endian (see Chapter 3 [endian], page 9) isn’t the same as its default value.

Example:
Ecriture du vecteur de réels T1 dans le fichier binaire test1.dat sous la forme d’entiers signés sur 4 octets.
```plaintext
vnumR T1;
T1=1,10;
writebin(test1.dat,"%4d",T1);
```

Ecriture des 30 premiers éléments du vecteur de réels T1 dans le fichier binaire test1.dat sous la forme d’entiers signés sur 4 octets.
```plaintext
vnumR T1;
T1=1,100;
writebin(test1.dat,[:30],"%4d",T1);
```

Ecriture des 30 premiers éléments du vecteur de réels T1 et de T2 dans le fichier binaire test2.dat dont chaque enregistrement est composé de 2 réels.
```plaintext
vnumR T1,T2;
T1=1,110;
```
T2=-100,-1;
writebin(test2.dat,:30,"%g%g",T1,T2);

Ecriture des vecteurs de réels de T3 dans le fichier binaire test3.dat
dont chaque enregistrement est composé de 4 réels.
vnumR T3[1:4];
T3[:]=1,10;
writebin(test3.dat,"%g%g%g%g",T3);

est équivalent à

writebin(test3.dat,"%g%g%g%g",T3[1],T3[2],T3[3],T3[4]);

10.9 Input/Output low level

10.9.1 file_open

<file> file_open

file_open(<filename> , read);
file_open(<filename> , write);
file_open(<filename> , write, append);
It opens a file in reading or writing mode depending on the value of the second parameter.
It returns an object identifier of type file.
In writing mode, the existing file is kept and the new writing operations are performed at
the end of the file if append is given, otherwise, the file is overwritten.
Example :
> f=file_open("sim2007.dat",read);
f = fichier "sim2007.dat" ouvert en lecture
> vnumR t;
> file_read(f,t);
> file_close(f);

10.9.2 file_close

file_close()

It closes a file previously opened with the function file_open.
Example :
> f=file_open("sim2007.dat",read);
f = fichier "sim2007.dat" ouvert en lecture
> vnumR t;
> file_read(f,t);
> file_close(f);
> stat(f);
fichier f = "sim2007.dat" ferme
Aucune erreur en lecture/ecriture

10.9.3 file_write

file_write

vec.> , ...);
file_write( <file> , <string> , (array of) num. vec. , ... );
file_write( <file> , (array of) num. vec. , ... );" equivalent to
file_write( fichier, [binf:bsup:step], "format", T, T1, T2).
file_write( fichier, "format", T, T1, T2).
file_write( fichier, T, T1, T2).
file_write( fichier, [binf:bsup:step], T, T1, T2).
This function is similar as the function write (see Section 10.8.3 [write], page 68) but this function requires an object identifier of type file instead of a filename. The function writes data in the file from the current position.

Example :
> f=file_open(sim2007.dat, write);
> t1=1:10;
> t2=-t1;
> file_write(f,t1);
> file_write(f,t2);
> file_close(f);

10.9.4 file_read

file_read( [Procedure]
file_read( <file> , [ <integer> : <integer> : <integer> ] ,
          (array of) num. vec. , ... ,
          (array of) num. vec. , <integer> ) , ... );

This function is similar as the function read (see Section 10.8.1 [read], page 63) but this function requires an object identifier of type file instead of a filename. The function reads data from the file from the current position.

Example :
> f=file_open(sim2007.dat, read);
> t1=1:10;
> file_read(f,[:5],t);
> vnumR t;
> file_read(f,t2);
> writes(t);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
+4.0000000000000000E+00
+5.0000000000000000E+00
> writes(t2);
+6.0000000000000000E+00
+7.0000000000000000E+00
+8.0000000000000000E+00
+9.0000000000000000E+00
+1.0000000000000000E+01
-1.0000000000000000E+00
-2.0000000000000000E+00
-3.0000000000000000E+00
-4.0000000000000000E+00
-5.0000000000000000E+00
-6.0000000000000000E+00
-7.0000000000000000E+00
-8.0000000000000000E+00
-9.0000000000000000E+00
-1.0000000000000000E+01
> file_close(f);

10.9.5 file_readappend

file_readappend

file_readappend(<file>, [ <integer> : <integer> : <integer> ]),
< (array of) num. vec. > ,...,
((<array of) num. vec.>, <integer>), ...
((<array of) num. vec.>, <integer>, <integer>), ...

equivalent to
file_readappend(fichier, [binf:bsup:step], T, (T1,n1), (T3,n2,n3));
file_readappend(fichier, T, (T1,n1), (T3,n2,n3));

This function is similar as the function readappend (see Section 10.8.2 [readappend], page 67) but this function requires an object identifier of type file instead of a filename. The function reads data from the file from the current position.

Example :
> f=file_open(sim2007.dat, read);
f = fichier "sim2007.dat" ouvert en lecture
> vnumR t;
> file_readappend(f,[:5],t);
> file_readappend(f,[10:15],t);
> writes(t);
+1.0000000000000000E+00
+2.0000000000000000E+00
+3.0000000000000000E+00
+4.0000000000000000E+00
+5.0000000000000000E+00
-5.0000000000000000E+00
-6.0000000000000000E+00
-7.0000000000000000E+00
-8.0000000000000000E+00
-9.0000000000000000E+00
-1.0000000000000000E+01
> file_close(f);
10.9.6 file_writemsg

file_writemsg

file_writemsg(<file> , <string> textformat);

file_writemsg(<file> , <string> textformat, <real> x, ...);

Write formatted messages to a file with(out) real constants.

The real constants must be formatted. The format is the same as the command printf in language C (see Section 7.6 [str], page 33, for the valid conversion specifiers). This message must be a string or a text between double-quotes. The messages could be on several lines.

To display double-quotes, two double-quotes must be used.

Example:
> f=file_open("temp001", write);
f = fichier "temp001" ouvert en ecriture
> x=4;
x = 4
> file_writemsg(f," resultat=%g\n", x);
> file_writemsg(f," termine\n");
> file_close(f);
> !"cat temp001";
resultat=4

> f=file_open("data1.txt",write);
f = fichier "data1.txt" ouvert en ecriture
> file_writemsg(f,"%d\n",3/2);
> file_writemsg(f,"%4.2E\n",3/2);
> file_close(f);
> !"cat data1.txt";
3/2
1.50E+00

10.9.7 file_writebin

file_writebin

file_writebin( <file> , [ <integer> : <integer> : <integer> ], <string> , <(array of) num. vec.> , ...);

file_writebin( <file> , <string> , <(array of) num. vec.> , ...);

equivalent to

file_writebin(file, [binf:bsup:step], "format", T, T1, T2).

This function is similar as the function writebin (see Section 10.8.5 [writebin], page 70) but this function requires an object identifier of type file instead of a filename. The function writes data in the file from the current position.

Example:
> f=file_open(sim2007.dat, write);
f = file "sim2007.dat" opened in write mode
> t1=1,10;
t1 double precision real vector : number of elements =10
> t2=-t1;
t2 double precision real vector : number of elements =10
Reference manual TRIP

10.10 Standard functions

10.10.1 minimum and maximum

imin
imin( <real vec.> )
Return the index of the minimal value of the numerical vector of real numbers. If more than one elements have the minimal value, then this function returns the index of the first element. If the numerical vector is empty then it returns 0.

```
Example :
> t=-5,5;
> imin(t);
1
```

imax
imax( <real vec.> )
Return the index of the maximal value of the numerical vector of real numbers. If more than one elements have the maximal value, then this function returns the index of the first element. If the numerical vector is empty then it returns 0.

```
Example :
> t=-5,5;
> imax(t);
10
```

min
min( <real or real vec.> , ...);
Return the minimum value of the parameters which could be real constants or real vectors.

```
Example :
>min(1.5,2.5);
3/2
> x=1.5;
> y=2.5;
> min(x,y);
3/2
> t=-10,10;
> p=abs(t);
> min(-5,t,p,20);
-10
```

max
max( <real or real vec.> , ...);
Return the maximum value of the parameters which could be real constants or real vectors.

```
Example :
> max(1.5,2.5);
5/2
> x=1.5;
```
Chapter 10: Numerical vectors

```
y=2.5;
> max(x,y);
5/2
> t=-10:10;
> p=abs(t);
> max(p,t,8);
10

IMIN

IMIN( <array of real vec.> )

Return a numerical vector of real numbers such that its elements have the index of the column where is located the minimal value term by term of each numerical vector.

The numerical vectors must have the same size.

Example:
> A=vnumR[1,2,3;6,7,1;9,2,5;13,11,16];
> writes("%g %g %g\n", A);
1 2 3
6 7 1
9 2 5
13 11 16
> B=IMIN(A);
B double precision real vector: number of elements = 4
> writes("%g\n", B);
1
3
2
2
```

IMAX

IMAX( <array of real vec.> )

Return a numerical vector of real numbers such that its elements have the index of the column where is located the maximal value term by term of each numerical vector.

The numerical vectors must have the same size.

Example:
> A=vnumR[1,2,3;6,7,1;9,2,5;13,11,16];
> writes("%g %g %g\n", A);
1 2 3
6 7 1
9 2 5
13 11 16
> B=IMAX(A);
B double precision real vector: number of elements = 4
> writes("%g\n", B);
3
2
1
3
```

MIN

MIN( <real vec.>, ...)
Return a numerical vector of real numbers such that its elements have the minimal value
term by term of each numerical vector.

The numerical vectors must have the same size.

For an array of numerical vector of real numbers, the operation is performed on each element.

Example :
> t=0,pi,pi/6$
> a=\text{MIN}(\text{cos}(t),\text{sin}(t))$
> c=\text{MAX}(t,\text{cos}(t),\text{sin}(t))$
> \text{writes}(a,c)$
+0.0000000000000000E+00 +1.0000000000000000E+00
+4.9999999999999994E-01 +8.6602540378443871E-01
+5.0000000000000011E-01 +1.0471975511965976E+00
+6.12323957367660E-17 +1.5707963267948966E+00
-4.9999999999999978E-01 +2.0943951023931953E+00
-8.6602540378443849E-01 +2.6179938779914940E+00
-1.0000000000000000E+00 +3.1415926535897931E+00
> vnumR cs[1:2]$
> cs[1]=\text{cos}(t)$$
> cs[2]=\text{sin}(t)$$
> b=\text{MIN}(cs)$$
> d=\text{MAX}(t,cs)$$
> \text{writes}(b,d);$
Chapter 10: Numerical vectors

> cs[2]=sin(t)
> b=MIN(cs)
> d=MAX(t,cs)
> writes(b,d);
+0.0000000000000000E+00 +1.0000000000000000E+00
+4.9999999999999994E-01 +8.6602540378443871E-01
+5.0000000000000011E-01 +1.0471975511965976E+00
+6.1232339957367660E-17 +1.5707963267948966E+00
-4.9999999999999978E-01 +2.0943951023931953E+00
-8.6602540378443849E-01 +2.6179938779914940E+00
-1.0000000000000000E+00 +3.1415926535897931E+00

10.10.2 addition and product

sum [Function]

sum( <num. vec.> )
sum( <num. vec.>, "kahan" )
sum( <num. vec.>, "sort", "kahan" )

Return the sum of elements in the numerical vector.

The option "sort" sort the vector using the ascending norm before the summation is done.
If the option "kahan" is set, the function uses the Kahan method to perform the summation (compensated summation).


Example :
> r=0,100$
> sum(r);
5050
> p=-10000,10000$
> p=p*pi/10000$
> sum(p);
3.885780586188048E-13
> // compensated summation
> sum(p,"sort","kahan");
0

prod [Function]

prod( <num. vec.> )

Return the product of elements in the numerical vector.

Example :
> p=1,10;
p Tableau de reels : nb reels =10
> prod(p);
3628800
> c=p+2*i*p;
c vecteur de complexes : nb complexes =10
> prod(c);
(860025600-i*11307340800)
accum  [Function]

accum( <num. vec.> )

Return the partial sum of elements in the numerical vector.

\[ Y = accum(X) : Y[N] = \sum_{i=1}^{N} X[i] \]

Example:
> tx=1:1:10;
> ty=accum(tx);
> writes(accum(tx));
+1.000000000000000E+00
+3.000000000000000E+00
+6.000000000000000E+00
+1.000000000000000E+01
+1.500000000000000E+01
+2.000000000000000E+01
+2.500000000000000E+01
+3.000000000000000E+01
+3.500000000000000E+01
+4.000000000000000E+01
+4.500000000000000E+01
+5.000000000000000E+01

10.10.3 sort

intersectvnum  [Function]

intersectvnum(<num. vec.> A ,<num. vec.> B)

It returns the values common to both numerical vectors A and B. The values of returned vector are in sorted order using ascending order.

Example:
> A = vnumR[7:1:7:5:4:6];
> B = vnumR[0:7:4:-2:5:3];
> C = intersectvnum(A,B);
> writes(C);
+4.000000000000000E+00
+5.000000000000000E+00
+7.000000000000000E+00

unionvnum  [Function]

unionvnum(<num. vec.> A ,<num. vec.> B)

It returns a vector which contains the combined values from A and B with no repetitions. The values of returned vector are in sorted order using ascending order.

Example:
> A = vnumR[7:1:7:5:4:6];
> B = vnumR[0:7:4:-2:5:3];
> C = unionvnum(A,B);
> writes(C);
reversevnum
reversevnum (<num. vec.> )
Reverse the order of elements in the numerical vector.

Example :
> p=1,10;
p Tableau de reels : nb reels =10
> writes(p, reversevnum(p));
+1.000000000000000E+00 +1.000000000000000E+01
+2.000000000000000E+00 +9.000000000000000E+00
+3.000000000000000E+00 +8.000000000000000E+00
+4.000000000000000E+00 +7.000000000000000E+00
+5.000000000000000E+00 +6.000000000000000E+00
+6.000000000000000E+00 +5.000000000000000E+00
+7.000000000000000E+00 +4.000000000000000E+00
+8.000000000000000E+00 +3.000000000000000E+00
+9.000000000000000E+00 +2.000000000000000E+00
+1.000000000000000E+01 +1.000000000000000E+00

sort
sort ( <real vec.> )
Sort in ascending order the elements in the real vector.
sort ( <real vec.> , <(array of) num. vec.> , ...);
Sort the elements in the (array of) numerical vectors according to the ascending sort of the first real vector.
Remarks : If you call ‘sort(TX,TY,TZ);’, the vector TY et TZ are sorted but TX isn’t sorted.

Example :
> p=vnumR[5:2:-3:7:1:6];
p Tableau de reels : nb reels =6
> sort(p);
> writes(p);
-3.000000000000000E+00 +1.000000000000000E+00
+2.000000000000000E+00 +9.000000000000000E+00
+5.000000000000000E+00 +6.000000000000000E+00
+6.000000000000000E+00 +5.000000000000000E+00
+7.000000000000000E+00 +4.000000000000000E+00
+8.000000000000000E+00 +3.000000000000000E+00
+9.000000000000000E+00 +2.000000000000000E+00
+1.000000000000000E+01 +1.000000000000000E+00

T=vnumC[1+i:2+2*i:3+3*i:1:4+4*i:1:5+5*i:1:7+7*i];
T vecteur de complexes : nb complexes =6
> TAB[1]=abs(T);
> TAB[2]=-real(T);
> sort(-p,T,TAB);

### 10.10.4 transposevnum

**transposevnum**

*Function*

**transposevnum** ( <array of num. vec.> )

Transpose the array of numerical vectors.

The array must have only one dimension and its numerical vectors must have the same size.

**Example:**

```plaintext
t = 1,5;
> t
Tableau de reels : nb reels =5
> vnumR TSRC[1:2];
> TSRC[1]=t;
> TSRC[2]=reversevnum(t);
> writes("%g %g\n",TSRC);
1 5
2 4
3 3
4 2
5 1
> TRES=transposevnum(TSRC);

TRES [1:5] nb elements = 5

> writes("%g %g %g %g %g\n",TRES);
1 2 3 4 5
5 4 3 2 1
```

### 10.10.5 standard math function

**abs**

*Function*

**abs** ( <constant or num. vec.> )

Return the absolute value of the constant if the value is a constant.

Return a numerical vector containing the absolute value of each element if the value is a numerical vector.

**Example:**

```plaintext
> abs(-2.3);
2.3
> x = 2.3;
> abs(x);
2.3
> t=-pi,pi,pi/100;
> at=abs(t);
at Tableau de reels : nb reels =200
> abs(1+i);
1.4142135623731
```

**arg**

*Function*

**arg** ( <constant or num. vec.> )
Return the argument (also called phase angle) of the constant if the value is a constant.
Return a numerical vector containing the argument (also called phase angle) of each element if the value is a numerical vector.

Example:
> arg(i);
  1.5707963267949
> arg(1+i);
  0.785398163397448
> t=0,pi,pi/6;
t Tableau de reels : nb reels =7
> writes(arg(exp(i*t)));
+0.000000000000000E+00
+5.235987755982987E-01
+1.047197551196598E+00
+1.570796326794897E+00
+2.094395102393195E+00
+2.617993877991494E+00
+3.141592653589793E+00

real
real(<constant or num. vec.>)

Return the real part of the constant if the value is a constant.
Return a numerical vector containing the real part of each element if the value is a numerical vector.

Example:
> real(i);
  0
> real(2+3*i);
  2
> t=0,pi,pi/6;
t Tableau de reels : nb reels =7
> writes(real(exp(i*t)));
+1.000000000000000E+00
+8.660254037844387E-01
+5.000000000000000E-01
-0.000000000000000E+00
-4.999999999999999E-01
-8.660254037844385E-01
-1.000000000000000E+00
> real(4);
  4

imag
imag(<constant or num. vec.>)

Return the imaginary part of the constant if the value is a constant.
Return a numerical vector containing the imaginary part of each element if the value is a numerical vector.

Example:
> imag(i);
  1
> imag(2+3*i);
  3
> t=0,pi,pi/6;
> writes(real(exp(i*t)));
  +1.000000000000000E+00
  +8.660254037844387E-01
  +5.000000000000000E-01
  -0.000000000000000E+00
  -4.999999999999998E-01
  -8.660254037844385E-01
  -1.000000000000000E+00
> imag(4);
  0

conj

conj( <constant or num. vec.> )
Return the conjugate of the constant if the value is a constant.
Return a numerical vector containing the conjugate of each element if the value is a numerical vector.

Example:
> conj(1+i);
  (1-i*1)
> conj(3);  

z vecteur de complexes : nb complexes =3
> writes(conj(z));
  +1.000000000000000E+00 -1.000000000000000E+00
  +3.000000000000000E+00 +5.000000000000000E+00
  +0.000000000000000E+00 -1.000000000000000E+00

erf

erf( <real or real vec.> )
Return the error function of the constant if the value is a real.
Return a numerical vector containing the error function of each element if the value is a numerical vector of real numbers.

The error function is defined by:

\[ erf(x) = \frac{2}{\sqrt{\pi}} \int_0^x \exp(-t^2) dt \]

Example:
> erf(1);
  0.842700792949715
> erf(0.5);
  0.520499877813047
> t=-2,2;
  t Tableau de reels : nb reels =5
> r=erf(t);
  r Tableau de reels : nb reels =5
> writes(t,r);
  -2.000000000000000E+00 -9.9532226501895271E-01
Chapter 10: Numerical vectors

-1.0000000000000000E+00 -8.4270079294971478E-01
+0.0000000000000000E+00 +0.0000000000000000E+00
+1.0000000000000000E+00 +8.4270079294971478E-01
+2.0000000000000000E+00 +9.953226501895271E-01

erfc

erfc(<real or real vec.>)
Return the complementary error function of the constant if the value is a real.
Return a numerical vector containing the complementary error function of each element if the value is a numerical vector of real numbers.
The complementary error function is defined by:

\[
erfc(x) = 1 - \frac{2}{(\sqrt{\pi})} \int_0^x \exp(-t^2) dt
\]

Example:
> erf(1);
0.842700792949715
> erf(0.5);
0.520499877813047
> t=-2,2;
t Tableau de reels : nb reels =5
> r=erf(t);
r Tableau de reels : nb reels =5
> writes(t,r);
-2.0000000000000000E+00 -9.9533226501895271E-01
-1.0000000000000000E+00 -8.4270079294971478E-01
+0.0000000000000000E+00 +0.0000000000000000E+00
+1.0000000000000000E+00 +8.4270079294971478E-01
+2.0000000000000000E+00 +9.953226501895271E-01

exp

exp(<constant or num. vec.>)
Return the base-e exponential of the constant if the value is a constant.
Return a numerical vector containing the base-e exponential of each element if the value is a numerical vector.

Example:
> exp(1.2);
3.32011692273655
> x = 1.2;
> exp(x);
3.32011692273655
> exp(2*i);
(-0.416146836547142+i*0.909297426825682)
> t=-2,2;
t Tableau de reels : nb reels =5
> r=exp(t);
r Tableau de reels : nb reels =5
> writes(t,r);
-2.0000000000000000E+00 +1.353352832366127E-01
-1.0000000000000000E+00 +0.3678794411714423E-01
+0.0000000000000000E+00 +1.0000000000000000E+00
**fac**

`fac(<constant or num. vec.>)`

Returns the factorial of a non-negative integer of the constant if the value is a constant.

Return a numerical vector containing the factorial of each element if the value is a numerical vector.

Example:

```plaintext
> fac(10);
3628800
> t = 1,5;
t double precision real vector : number of elements = 5
> r = fac(t);
r double precision real vector : number of elements = 5
> writes(t,r);
+1.000000000000000E+00 +1.000000000000000E+00
+2.000000000000000E+00 +2.000000000000000E+00
+3.000000000000000E+00 +6.000000000000000E+00
+4.000000000000000E+00 +2.400000000000000E+01
+5.000000000000000E+00 +1.200000000000000E+02
```

**sqrt**

`sqrt(<constant or num. vec.>)`

Returns the square root of the constant if the value is a constant.

Return a numerical vector containing the square root of each element if the value is a numerical vector.

Example:

```plaintext
> x = 4;
> sqrt(x);
2
> sqrt(1+i);
(1.09868411346781+i*0.45508986056227)
> t = 0,5;
t Tableau de reels : nb reels = 6
> ts = sqrt(t);
ts Tableau de reels : nb reels = 6
> writes(t,ts);
+0.000000000000000E+00 +0.000000000000000E+00
+1.000000000000000E+00 +1.000000000000000E+00
+2.000000000000000E+00 +1.414213562373095E+00
+3.000000000000000E+00 +1.732050807568877E+00
+4.000000000000000E+00 +2.000000000000000E+00
+5.000000000000000E+00 +2.236067977499790E+00
```

**cos**

`cos(<constant or num. vec.>)`

Returns the cosine of the constant if the value is a constant.

Return a numerical vector containing the cosine of each element if the value is a numerical vector.
Example:
> cos(0.12);
0.992808635853866
> x=0;
x = 0
> cos(x);
> t=-pi,pi,pi/100;
> at=cos(t);
> cos(1+i);
(0.833730025131149-1*0.988897705762865)

\[
\text{sin} \quad [\text{Function}]
\]

\text{sin}( <\text{constant or num. vec.}> )

Return the sine of the constant if the value is a constant.
Return a numerical vector containing the sine of each element if the value is a numerical vector.

Example:
> sin(0.12);
0.119712207288912
> x = 0.12;
> sin(x);
0.119712207288912
> t=-pi,pi,pi/100;
> s=sin(t);
> sin(4*i);
(0+i*27.2899171971277)

\[
\text{tan} \quad [\text{Function}]
\]

\text{tan}( <\text{constant or num. vec.}> )

Return the tangent of the constant if the value is a constant.
Return a numerical vector containing the tangent of each element if the value is a numerical vector.

Example:
> x = 1.2;
> tan(x);
2.57215162212632
> t=-pi,pi,pi/100;
> at=tan(t);
> tan(-1+i);
(-0.271752585319512+1*95227/87854)

\[
\text{acos} \quad [\text{Function}]
\]

\text{acos}( <\text{real or real vec.}> )

Return the principal value of the arc cosine of the real number if the value is a real number.
Return a real vector containing the principal value of the arc cosine of each element if the value is a real vector.

Example:
> acos(0.12);
1.4505064440011
> x = 0.12;
> acos(x);
1.45050644400111
> t=-pi,pi,pi/100;
> at=acos(t);

asin

asin( <real or real vec.> )

Return the principal value of the arc sine of the real number if the value is a real number.
Return a real vector containing the principal value of the arc sine of each element if the value is a real vector.

Example:
> asin(0.12);
0.120289882394788
> x = 0.12;
> asin(x);
0.120289882394788
> t=-pi,pi,pi/100;
> at=asin(t);

atan

atan( <real or real vec.> )

Return the principal value of the arc tangent of the real number if the value is a real number.
Return a real vector containing the principal value of the arc tangent of each element if the value is a real vector.

Example:
> atan(2);
1.10714871779409
> x = 2;
> atan(x);
1.10714871779409
> t=-pi,pi,pi/100;
> at=atan(t);

cosh

cosh( <constant or num. vec.> )

Return the hyperbolic cosine of the constant if the value is a constant.
Return a numerical vector containing the hyperbolic cosine of each element if the value is a numerical vector.

Example:
> cosh(0.12);
1.0072086414827
> x = 0.12;
> cosh(x);
1.0072086414827
> t=-pi,pi,pi/100;
> at=cosh(t);
> cosh(1+i);
(0.833730025131149+i*0.988897705762865)
**sinh**  

sinh( <constant or num. vec.> )  

Return the hyperbolic sine of the constant if the value is a constant.  
Return a numerical vector containing the hyperbolic sine of each element if the value is a numerical vector.

Example :
> sinh(1.2);
1.50946135541217
> x = 1.2;
> sinh(x);
1.50946135541217
> t=-π,π,π/100;
> s=sinh(t);
> sinh(-1+3*i);
(1.16344036370325+i*0.217759551622152)

**tanh**

tanh( <constant or num. vec.> )

Return the hyperbolic tangent of the constant if the value is a constant.  
Return a numerical vector containing the hyperbolic tangent of each element if the value is a numerical vector.

Example :
> x = 1.2;
> tanh(x);
0.833654607012155
> tanh(1+i);
(95227/87854+i*0.271752585319512)
> t=0,10;
tTableau de reels : nb reels =11
> tanh(t);

**acosh**

acosh( <constant or num. vec.> )

Return the inverse of the hyperbolic cosine of the constant if the value is a constant.  
Return a numerical vector containing the inverse of the hyperbolic cosine of each element if the value is a numerical vector.

Example :
> acosh(1.2);
0.6223625037147786
> t=1,2,0.1;
tTableau de reels double-precision : nb reels =11
> acosh(t);
Tableau de reels double-precision : nb reels =11

**asinh**

asinh( <constant or num. vec.> )

Return the inverse of the hyperbolic sine of the constant if the value is a constant.  
Return a numerical vector containing the inverse of the hyperbolic sine of each element if the value is a numerical vector.
Example :
> asinh(1.2);  
  1.015973134179692  
> t=1,2,0.1;  
t  Tableau de reels double-precision : nb reels =11  
> asinh(t);  
  Tableau de reels double-precision : nb reels =11

`asinh`  
`asinh(constant or num. vec.)`  
Return the inverse of the hyperbolic tangent of the constant if the value is a constant.  
Return a numerical vector containing the inverse of the hyperbolic tangent of each element if the value is a numerical vector.

Example :
> atanh(0.2);  
  0.2027325540540822  
> t=0,1,0.1;  
t  Tableau de reels double-precision : nb reels =11  
> atanh(t);  
  Tableau de reels double-precision : nb reels =11

`atan2`  
`atan2(real or real vec., real or real vec.)`  
Return the principal value of the arc tangent of the first argument divided by the second argument using the signs of both arguments to determine the quadrant of the return value.

Example :
> y = 11.5;  
> x = 14;  
> atan2(y,x);  
  0.68767125603875

`mod`  
`mod(real or real vec., real or real vec.)`  
`mod(real or real vec.)`  
Return the modulo of the first and second value. The result has the same sign as the first argument and magnitude less than the magnitude of the second argument.

Example :
> mod(4,2);  
  0  
> x=4$  
> y=2$  
> mod(x,y);  
  0  
> t=21,30;  
> r=1,10;  
> f=mod(t,r);  
> g=mod(t,3);

`int`  
`int(real or real vec.)`  
Return the integer part of the real number if the value is a real.  
Return a real vector containing the integer part of each element if the value is a real vector.
Example:
> x=1.23;
> int(x);
1
> t=-pi,pi,pi/100;
> at=int(t);

log [Function]
log( <constant or num. vec.> )
Return the neperian logarithm of the constant if the value is a constant.
Return a numerical vector containing the neperian logarithm of each element if the value is a numerical vector.

Example:
> log(3);
1.09861228866811
> x = 3;
> log(x);
1.09861228866811
> r=1,10;
> at=log(r);
> log(2+i);
(0.80471895621705+i*0.463647609000806)

log10 [Function]
log10( <constant or num. vec.> )
Return the decimal logarithm (base 10) of the constant if the value is a constant.
Return a numerical vector containing the decimal logarithm of each element if the value is a numerical vector.

Example:
> log(3);
1.09861228866811
> x = 3;
> log(x);
1.09861228866811
> r=1,10;
> at=log(r);
> log(2+i);
(0.80471895621705+i*0.463647609000806)

nint [Function]
nint( <real or real vec.> )
Return the nearest integer of the real number if the value is a real.
Return a real vector containing the nearest integer of each element if the value is a real vector.

Example:
> x=1.23;
> x = 1.23
> nint(x);
1
> nint(1.78);
> t=-1,1,0.1;
t  Tableau de reels : nb reels =21
> nint(t);

**sign**

sign( <real or real vec.> )

Return the sign of the real number if the value is a real.

Return a real vector containing the sign of each element if the value is a real vector.

The function is defined with the following rules:

- \( \text{sign}(x) = +1 \) if \( x > 0 \)
- \( \text{sign}(x) = 0 \) if \( x = 0 \)
- \( \text{sign}(x) = -1 \) if \( x < 0 \)

Example :

> sign(3);
1
> sign(-3);
-1
> sign(0);
0

> t=-3,3;
t  Tableau de reels : nb reels =7
> writes(t,sign(t));
-3.0000000000000000E+00 -1.0000000000000000E+00
-2.0000000000000000E+00 -1.0000000000000000E+00
-1.0000000000000000E+00 -1.0000000000000000E+00
+0.0000000000000000E+00 +0.0000000000000000E+00
+1.0000000000000000E+00 +1.0000000000000000E+00
+2.0000000000000000E+00 +1.0000000000000000E+00
+3.0000000000000000E+00 +1.0000000000000000E+00

**histogram**

histogram(<real vec.> TY, <real vec.> TX)

Return the histogram of \( TY \) with the specified ranges \( TX \):

\( TZ=\text{histogram}(TY,TX) : TZ[N] \) contains the number of elements in \( TY \) such that \( TX[N] \leq TY[j] < TX[N+1] \).

For the last range \( TX \), the relation is : \( TX[N] \leq TY[j] \leq TX[N+1] \)

Example :

> TX=-3,9,2$
> writes("%g\n",TX);
-3
-1
1
3
5
7
9
> TZ=histogram(TY,TX)$
> writes("%g\n",TZ);
1
10.11 Conditions

\(?()::\) [Operator]

\(?(<\text{condition}>\)\)

Return a numerical vector which contains only the numbers 0 or 1. For each element of the array: if the condition at index j is true then \(<\text{name}>[j]=1\) else \(<\text{name}>[j]=0\)

\(?(<\text{condition}>): <\text{constant or num. vec.}> \text{tabvrai} : <\text{constant or num. vec.}> \text{tabfaux}\)

Return a numerical vector which contains only the numbers 0 or 1. For each element of the array: if the condition at index j is true then \(<\text{name}>[j]=\text{tabvrai}[j]\) else \(<\text{name}>[j]=\text{tabfaux}[j]\)

All vectors must have the same size.

Example:
> \(t=0.5;\)
  \(t\) double precision real vector: number of elements = 6
> \(r=5-t;\)
  \(r\) double precision real vector: number of elements = 6
> \(q=?(t<=r);\)
  \(q\) double precision real vector: number of elements = 6
> \(\text{writes}(q);\)
+1.0000000000000000E+00
+1.0000000000000000E+00
+1.0000000000000000E+00
+0.0000000000000000E+00
+0.0000000000000000E+00
+0.0000000000000000E+00
> \(l=?(t<=r):i:5;\)
  \(l\) Double precision complex vector: number of elements = 6
> \(\text{writes}(l);\)
  +0.0000000000000000E+00 +1.0000000000000000E+00
  +0.0000000000000000E+00 +1.0000000000000000E+00
  +0.0000000000000000E+00 +1.0000000000000000E+00
  +5.0000000000000000E+00 +0.0000000000000000E+00
  +5.0000000000000000E+00 +0.0000000000000000E+00
  +5.0000000000000000E+00 +0.0000000000000000E+00
> \(m=?(t>2):r:t;\)
  \(m\) double precision real vector: number of elements = 6
> \(\text{writes}(m);\)
+0.0000000000000000E+00
+1.0000000000000000E+00
+2.0000000000000000E+00
+2.0000000000000000E+00
+1.0000000000000000E+00
+0.0000000000000000E+00
>
10.12 Conversion

For the conversion from or to a numerical matrix, see Section 11.10 [ConversionMat], page 109.

10.12.1 dimtovnumR

\textit{dimtovnumR} \hspace{1cm} \textbf{[Function]}

\texttt{dimtovnumR( <array> )}

Return a numerical vector of real numbers from the array of series.

The array of series must contain only real numbers. The array must have only one dimension.

Example :
\begin{verbatim}
> dim ts[1:3];
> ts[1]=1$
> ts[2]=4$
> ts[3]=6$
> tr=dimtovnumR(ts)
> writes(tr);
+1.000000000000000E+00
+4.000000000000000E+00
+6.000000000000000E+00
> ltr=log(dimtovnumR(ts));
ltr Tableau de reels : nb reels =3
\end{verbatim}

10.12.2 dimtovnumC

\textit{dimtovnumC} \hspace{1cm} \textbf{[Function]}

\texttt{dimtovnumC( <array> )}

Return a numerical vector of complex numbers from the array of series.

The array of series must contain only complex numbers. The array must have only one dimension.

Example :
\begin{verbatim}
> dim ts[1:3];
> ts[1]=1+i$
> ts[2]=3*i$
> ts[3]=-5+i$
> tc=dimtovnumC(ts)
> writes(tc);
+1.000000000000000E+00 +1.000000000000000E+00
+0.000000000000000E+00 +3.000000000000000E+00
-5.000000000000000E+00 +1.000000000000000E+00
> ltc=exp(dimtovnumC(ts));
ltc vecteur de complexes : nb complexes =3
\end{verbatim}

10.12.3 vnumtodim

\textit{vnumtodim} \hspace{1cm} \textbf{[Function]}

\texttt{vnumtodim( <(array of) num. vec.> )}

Return an array of series (constants) from the array of numerical vectors or from the numerical vector.

Example :
\begin{verbatim}
> tr=1,4;
\end{verbatim}
Tableau de reels : nb reels = 4
> tsr=vnumtodim(tr);

tsr [1:4] nb elements = 4

> afftab(tsr);
tsr[1] = 1
tsr[2] = 2
tsr[3] = 3
tsr[4] = 4
> vnumC tc[1:2];
> tc[1]=i*tr$
> tc[2]=tr+i*tr**2$
> tsc=vnumtodim(tc);

tsc [1:4, 1:2] nb elements = 8

> afftab(tsc);
tsc[1,1] = (0+i*1)
tsc[1,2] = (1+i*1)
tsc[2,1] = (0+i*2)
tsc[2,2] = (2+i*4)
tsc[3,1] = (0+i*3)
tsc[3,2] = (3+i*9)
tsc[4,1] = (0+i*4)
tsc[4,2] = (4+i*16)

10.12.4 str

str( <real vec.> );
str( <string> format, <real vec.> );

Convert an vector of real numbers in an array of character strings. If format is specified, then each element of the vector is converted to this format.

For the description of the formats, see Section 7.6 [str], page 33.

Example :
> t = 8,10;
t double precision real vector : number of elements =3
> tabs = str("%04g", t);

tabs [1:3 ] number of elements = 3

> afftab(tabs);
tabs[1] = "0008"
tabs[2] = "0009"
tabs[3] = "0010"
11 Numerical matrices

The numerical values stored in matrices are always real or complex double-precision, quadruple-precision or multiprecision numbers depending on the current numerical mode. A numerical matrix is assumed to be a two-dimensional matrix. The numerical matrix are not resizeable. The first dimension is the lines of the matrix and the second dimension is the columns of the matrix.

11.1 Declaration

Explicit declaration for numerical matrices are only required before call to **read**, **readbin** (see Section 10.8 [Entree/SortieTabNum], page 63) and assignment of a single or several elements. Explicit declaration for array of numerical matrices are always required.

11.1.1 matrixR

```
matrixR <Procedure>

matrixR <name> ( [ <two dimensions of a matrix> MxN ] ) , ... ;  
matrixR <name> [ <dimension of an array> ] ( [ <two dimensions of a matrix> MxN ] ) , ... ;
```

It declares a real matrix of dimension MxN or an array of real matrices of dimension MxN.

After the declaration, the elements of the matrix are initialized with the value 0.

Example :

```>
> matrixR C([1:3, 1:5]);
> stat(C);
Double precision real matrix C [ 1:3 , 1:5 ].
Size of the array in bytes: 120
> bounds= 1:3,1:6;
bounds = bounds [ 1:3, 1:6 ]
> matrixR T0[1:2]([bounds]);
> stat(T0);
Array of series
T0 [ 1:2 ]
list of array’s elements :
T0 [ 1 ] =
Double precision real matrix T0 [ 1:3 , 1:6 ].
Size of the array in bytes: 144
T0 [ 2 ] =
Double precision real matrix T0 [ 1:3 , 1:6 ].
Size of the array in bytes: 144
>
```

11.1.2 matrixC

```
matrixC <Procedure>

matrixC <name> ( [ <two dimensions of a matrix> MxN ] ) , ... ;  
matrixC <name> [ <dimension of an array> ] ( [ <two dimensions of a matrix> MxN ] ) , ... ;
```

It declares a complex matrix of dimension MxN or an array of complex matrices of dimension MxN.

After the declaration, the elements of the matrix are initialized with the value 0+i*0.
Example:
> matrixC C([1:3, 1:5]);
> stat(C);
  Double precision complex matrix C [ 1:3 , 1:5 ].
  Size of the array in bytes: 240
> bounds=1:3,1:6;
bounds = bounds [ 1:3, 1:6 ]
> matrixC T0[1:2]([bounds]);
> stat(T0);
  Array of series
  T0 [ 1:2 ]
  list of array’s elements :
  T0 [ 1 ] =
  Double precision complex matrix T0 [ 1:3 , 1:6 ].
  Size of the array in bytes: 288
  T0 [ 2 ] =
  Double precision complex matrix T0 [ 1:3 , 1:6 ].
  Size of the array in bytes: 288
>

11.2 Initialization

**matrixR[,]**

< Function >
< name > = matrixR [ <real> or <real vec.> or <array of real vec.> , ... : <real> ,... ] ;
It declares and initializes a real matrix with the specified reals or real vectors.
The character : separates the lines and the character , separates the columns. All columns
must have the same line counts.

Example:
> // declare a real matrix 2x3
> tab3=matrixR[1,2,3;4,5,6];
tab3 double precision real matrix [ 1:2 , 1:3 ]
> writes(tab3);
+1.0000000000000000E+00 +2.0000000000000000E+00 +3.0000000000000000E+00 +4.0000000000000000E+00
+5.0000000000000000E+00 +6.0000000000000000E+00
>

**matrixC[,]**

< Function >
< name > = matrixC [ <complex> or <complex vec.> , ... : <complex> ,... ] ;
It declares and initializes a complex matrix with the specified complex numbers or complex
vectors.
The character : separates the lines and the character , separates the columns. All columns
must have the same line counts.

Example:
> // declare a complex matrix 2x3
> tab3=matrixC[1+2*i,3+4*i,5:2,4*i, -7+2*i];
tab3 double precision complex matrix [ 1:2 , 1:3 ]
> writes(tab3);
+1.0000000000000000E+00 +2.0000000000000000E+00 +3.0000000000000000E+00 +4.0000000000000000E+00
+5.0000000000000000E+00 +6.0000000000000000E+00
+2.0000000000000000E+00 +0.0000000000000000E+00 +0.0000000000000000E+00 +4.0000000000000000E+00
-7.0000000000000000E+00 +2.0000000000000000E+00
11.3 Display

writes

writes([<integer> : <integer> : <integer>], <string>, (array of) matrix, ...);
writes(<string>, (array of) matrix, ...);
writes((array of) matrix, ...);
equivalent to
writes({[binf:{bsup}]:{step}}, {format}, (array of) matrix, ...).

It prints, to the screen, numerical matrices or the array of numerical matrices in columns form.

From the binfth line of each numerical matrix (if binf isn’t specified, it starts from the first line).

To the bsupth line of each numerical matrix (if bsup isn’t specified, it stops to the last line of the biggest matrix).

Every step line (if step isn’t specified, the step is 1).

format is optional. This format is similar to the C format (cf. printf) and must be between double quotes ("'). The format must contain the same number of format specifiers (e.g., %g) as the number of columns.

A complex matrix will be printed on two times more columns (the first for the real part and the second for the imaginary part).

Example :
> // display a real matrix 2x3
> mat3=matrixR[1,2,3;4,5,6];
mat3 double precision real matrix [ 1:2 , 1:3 ]
> writes(mat3);
+1.0000000000000000E+00 +2.0000000000000000E+00 +3.0000000000000000E+00
+4.0000000000000000E+00 +5.0000000000000000E+00 +6.0000000000000000E+00
> // display a complex matrix 2x3
> mat4=matrixC[1+2*i,3+4*i,5:2,4*i, -7+2*i];
mat4 double precision complex matrix [ 1:2 , 1:3 ]
> writes(6*"%g "+"\n", mat4);
1 2 3 4 5 0
2 0 0 4 -7 2
>
afftab

afftab(<matrix>);

It prints, to the screen, a numerical matrix. Each line of the matrix is surrounded by the characters [].

Example :
> _affc=1$
> // display a real matrix 2x3
> mat3=matrixR[1,2,3;4,5,6];
mat3 double precision real matrix [ 1:2 , 1:3 ]
> afftab(mat3);
[ 1 2 3]
[ 4 5 6]
> // display a complex matrix 2x3
> mat4=matrixC[1+2*i,3+4*i,5:2,4*i, -7+2*i];
mat4 double precision complex matrix [ 1:2 , 1:3 ]
> afftab(mat4);
[(1+i*2) (3+i*4) 5]
[ 2 (0+i*4) (-7+i*2)]
>

### 11.4 Size

**size**

\[
\text{size( <matrix> , <integer> n )}
\]

Return the number of lines of the numerical matrix if \( n=1 \).

Return the number of columns of the numerical matrix if \( n=2 \).

**Example:**

> mat3=matrixR[1,2,3:4,5,6];
mat3 double precision real matrix [ 1:2 , 1:3 ]
> size(mat3,1);
2
> size(mat3,2);
3
>

### 11.5 Data retrieval

**[::,::]**

\[
\text{<matrix> \[ <integer> binfli : <integer> bsupli : <integer> pasli , <integer> binfcol : <integer> bsupcol : <integer> pascol ]}
\]

Return a numerical matrix which contains only the elements located between the lower bound and the upper bound with the specified step.

If the lower bound isn’t specified, then the lower bound is assumed to be 1.

If the upper bound isn’t specified, then the upper bound is assumed to be the size of the numerical vector.

If the step isn’t specified, then the step is assumed to be 1.

**Remark:** All missing combinations are allowed.

**Example:**

> M=matrixR[1,2,3:4,5,6:7,8,9];
M double precision real matrix [ 1:3 , 1:3 ]
> r=M[::2,::2];
\[
\begin{pmatrix}
+1.0000000000000000E+00 & +3.0000000000000000E+00 \\
+7.0000000000000000E+00 & +9.0000000000000000E+00
\end{pmatrix}
\]
> v=M[1:2,2:3];
v double precision real matrix [ 1:2 , 1:2 ]
> writes(v);
\[
\begin{pmatrix}
+2.0000000000000000E+00 & +3.0000000000000000E+00 \\
+5.0000000000000000E+00 & +6.0000000000000000E+00
\end{pmatrix}
\]
11.6 Input/Output

The following functions work on the numerical matrices in the same behavior as on the numerical vectors:
- write (see Section 10.8.3 [write], page 68)
- writebin (see Section 10.8.5 [writebin], page 70)

11.6.1 sauve_c

sauve_c

sauve_c(<matrix> , <filename> );

It saves the matrix to the file using the C language (version C99). The file is created in the
directory specified by _path.

sauve_c(<matrix> , <file> );

It saves the matrix to the file, previously opened in written mode, using the C language
(version C99).

Example :
> A=matrixR[9,0,7
   :1,2,3
   :4,5,6];
   A double precision real matrix [ 1:3 , 1:3 ]
> sauve_c(A, "prog.c");
>

11.6.2 sauve_fortran

sauve_fortran

sauve_fortran(<matrix> , <filename> );

It saves the matrix to the file using the Fortran language (version 77 or later). The file source
format is compatible with the fixed and free format. The file is created in the directory
specified by _path.

sauve_fortran(<matrix> , <file> );

It saves the matrix to the file, previously opened in written mode, using the Fortran language
(version 77 or later). The file source format is compatible with the fixed and free format.

Example :
> A=matrixR[9,0,7
   :1,2,3
   :4,5,6];
   A double precision real matrix [ 1:3 , 1:3 ]
> sauve_fortran(A, "prog.f");
>

11.6.3 sauve_tex

sauve_tex

sauve_tex(<matrix> , <filename> );

It saves the matrix to the file using the \TeX form. The file is created in the directory specified
by _path.

sauve_tex(<serie> , <file> );

It saves the matrix to the file, previously opened in written mode, using the \TeX form.
Example :
> A=matrixR[9,0,7
     :1,2,3
     :4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> sauve_tex(A, "essai.tex");
>
11.6.4 sauve_ml

**sauve_ml**

sauve_ml(<matrix>, <filename>);

It saves the matrix to the file using the MathML 2.0 (concept) form. The file is created in
the directory specified by _path._

sauve_ml(<matrix>, <file>);

It saves the matrix to the file, previously opened in written mode, using the MathML 2.0
(concept) form.

Example :
> A=matrixR[9,0,7
     :1,2,3
     :4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> sauve_ml(A, "essai.ml");
>
11.6.5 write

This function works on the numerical matrices in the same behavior as on the numerical vectors
(see Section 10.8.3 [write], page 68).

11.6.6 writebin

This function works on the numerical matrices in the same behavior as on the numerical vectors
(see Section 10.8.5 [writebin], page 70).

11.7 Input/Output low level

The following functions work on the numerical matrices in the same behavior as on the numerical
vectors :

– **file_write** (see Section 10.9.3 [file_write], page 72)

11.8 Standard math function

The following functions work on the numerical matrices in the same behavior as on the numerical
vectors. These operations apply on each element of the matrix.

• abs
• acos
• acosh
• arg
• asin
• asinh
• atan
• atanh
• atan2
• conj
• cos
• cosh
• erf
• erfc
• exp
• imag
• int
• log
• log10
• MAX
• MIN
• mod
• nint
• real
• sign
• sin
• sinh
• tan
• tanh

The following functions work on the numerical matrices in the same behavior as on the numerical vectors. These operations apply on the matrix.

• max
• min
• sum

11.8.1 Matrix product

\[ \text{Operator} \]

\[ \text{matrix} \times \text{matrix} \]

Compute the matrix product of the numerical matrix.

Compute \( r = a \times b \) such that \( r[i,j] = \sum_{k=1}^{\text{size}(a,2)} a[i,k] \times b[k,j] \)

Remark: the number of columns of the first matrix must be same as the number of lines of the second matrix.

Example:
> _affc=1$
> A = \text{matrixR}[1,3,5]
   :2,4,6];
A double precision real matrix [ 1:2 , 1:3 ]
> B = \text{matrixR}[5,8,11]
   :7,6,8
   :4,0,8];
B double precision real matrix [ 1:3 , 1:3 ]
\texttt{C = A\&*B;}
\texttt{C double precision real matrix \{1:2, 1:3\}}
\texttt{afftab(C);} 
\begin{tabular}{ccc}
  46 & 26 & 75 \\
  62 & 40 & 102 \\
\end{tabular}
\texttt{>}

\begin{table}[h]
\texttt{&*} \hspace{1cm} \texttt{[Operator]}
\texttt{<matrix> \&* <num. vec.>} \\
\texttt{<num. vec.> \&* <matrix>}
\caption{Operator \&*}
\end{table}

\texttt{Compute the matrix product of the numerical matrix and a numerical vector.}
\texttt{If b is the numerical vector, i computes r=a\&*b such that r[i] = \sum_{k=1}^{size(a,2)} a_{i,k} \times b_k}

### 11.8.2 Kronecker product

\texttt{kroneckerproduct} \\
\texttt{kroneckerproduct(<matrix> A ,<matrix> B)}

\texttt{It returns the Kronecker product of two numerical matrices.}

\begin{example}
\texttt{affc=1}$\$
\texttt{A = matrixR[1,2,3,4,5,6]$} \\
\texttt{B = matrixR[5,0:10,1:7,3]$} \\
\texttt{C = kroneckerproduct(A,B)$} \\
\texttt{afftab(C);} \\
\begin{tabular}{cccccc}
  5 & 0 & 10 & 0 & 15 & 0 \\
  10 & 1 & 20 & 2 & 30 & 3 \\
  7 & 3 & 14 & 6 & 21 & 9 \\
  20 & 0 & 25 & 0 & 30 & 0 \\
  40 & 4 & 50 & 5 & 60 & 6 \\
  28 & 12 & 35 & 15 & 42 & 18 \\
\end{tabular}
\end{example}

### 11.8.3 Matrix determinant

\texttt{det} \\
\texttt{det( <matrix> )}

\texttt{Computes the determinant of a square matrix.}

\texttt{Remarks : When computations are performed in double precision, the Lapack Library is used. The LU algorithm is used in all numerical precision.}

\begin{example}
\texttt{t=matrixR[9,0,7 \\
:1,2,3 \\
:4,5,6];}$\$
\texttt{t double precision real matrix \{1:3, 1:3\}} \\
\texttt{det(t);}
\end{example}

### 11.8.4 Matrix inversion

\texttt{invertmatrix} \\
\texttt{invertmatrix( <matrix> )}
Chapter 11: Numerical matrices

Computes the inverse of a square matrix.
Remarks: When computations are performed in double precision, the Lapack Library is used. The LU algorithm is used in all numerical precision.

Example:
> .affc=2$
> A=\text{matrixR}[9,0,7
:1,2,3
:4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> B=\text{invertmatrix}(A); B double precision real matrix [ 1:3 , 1:3 ]
> afftab(B);
[ 0.0625 - 0.72916667 0.29166667]
[ -0.125 - 0.54166667 0.41666667]
[ 0.0625 0.9375 - 0.375]
>

11.8.5 Matrix trace

\text{tracematrix}
\text{tracematrix( <matrix> )}
Computes the trace of a square matrix.

Example:
> A=\text{matrixR}[9,0,7
:1,2,3
:4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> t=\text{tracematrix}(A);
t = 17
>

11.8.6 Transpose

\text{transposematrix}
\text{transposematrix( <matrix> )}
Computes the transpose of a square matrix.

Example:
> .affc=1$
> A=\text{matrixR}[9,0,7
:1,2,3];
A double precision real matrix [ 1:2 , 1:3 ]
> B=\text{transposematrix}(A);
B double precision real matrix [ 1:3 , 1:2 ]
> afftab(B);
[ 9 1]
[ 0 2]
[ 7 3]
>
11.8.7 Identity matrix

identitymatrix

identitymatrix( <operation> n )
Computes the identity matrix of size n. This matrix is a n x n square matrix with ones on the main diagonal and zeros elsewhere.

Example :
> _affc=1$
> I3=identitymatrix(3);
I3 double precision real matrix [ 1:3 , 1:3 ]
> afftab(I3);
[ 1 0 0]
[ 0 1 0]
[ 0 0 1]
>

11.8.8 Eigenvalues

eigenvalues

eigenvalues( <matrix> )
Computes the eigenvalues of the square matrix using a QR algorithm (lapack library).

Example :
> A=matrixR[9,0,7 :1,2,3 :4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> B=eigenvalues(A);
B Double precision complex vector : number of elements =3
> writes(B);
+1.3769150418957313E+01 +0.0000000000000000E+00
+4.0843620348094420E+00 +0.0000000000000000E+00
-8.5351245376675389E-01 +0.0000000000000000E+00

11.8.9 Eigenvectors

eigenvectors

eigenvectors( <matrix> MAT, <matrix> TVECT, <matrix> TVAL)
eigenvectors( <matrix> MAT, <matrix> TVECT)
Compute the eigenvectors of the square matrix MAT using a QR algorithm (lapack library). It stores the eigenvectors in the matrix TVECT and the eigenvalues in the vector TVAL.

Example :
> _affc=1$
> A=matrixR[9,0,7 :1,2,3 :4,5,6];
A double precision real matrix [ 1:3 , 1:3 ]
> eigenvectors(A,_vectp, valp);
> afftab(_vectp);
[ - 0.808169 - 0.750577 - 0.484664]
[ - 0.209021 0.398522 - 0.547409]
[ - 0.550611 0.527081 0.682234]
Chapter 11: Numerical matrices

> writes(valp);
+1.3769150418957313E+01 +0.0000000000000000E+00
+4.0843620348094420E+00 +0.0000000000000000E+00
-8.5351245376675389E-01 +0.0000000000000000E+00
> // first vector
> p=vectp[:,1]$
> writes(p);
-8.0816903814944341E-01 +0.0000000000000000E+00
-2.0902130660832199E-01 +0.0000000000000000E+00
-5.5061138669696397E-01 +0.0000000000000000E+00

11.8.10 Arithmetic
The matrices must have the same number of elements on each dimension.

+ [Operator]
<matrix> + <matrix>
Add term by term two matrices.

+ [Operator]
<matrix> + <constant>
<constant> + <matrix>
Add a constant with each element of the matrix.

+ [Operator]
<matrix> + <num. vec.>
<num. vec.> + <matrix>
Add term by term the numerical vector and the matrix. The matrix must contain only one column.

* [Operator]
<matrix> * <matrix>
Multiply term by term two matrices.

* [Operator]
<matrix> * <constant>
<constant> * <matrix>
Multiply a constant with each element of the matrix.

* [Operator]
<num. vec.> * <matrix>
<num. vec.> * <matrix>
Multiply term by term the numerical vector and the matrix. The matrix must contain only one column.

- [Operator]
<matrix> - <matrix>
Subtract term by term two matrices.

- [Operator]
<matrix> - <constant>
<constant> - <matrix>
Subtract a numerical constant with each element of the matrix.
- [Operator]
  
  <matrix> - <num. vec.>
  <num. vec.> - <matrix>

  Subtract term by term the numerical vector and the matrix. The matrix must contain only one column.

/ [Operator]

  <matrix> / <matrix>

  Divide term by term two matrices.

/ [Operator]

  <matrix> / <constant>
  <constant> / <matrix>

  Divide a numerical constant with each element of the matrix.

/ [Operator]

  <matrix> / <num. vec.>
  <num. vec.> / <matrix>

  Divide term by term the numerical vector and matrix. The matrix must contain only one column.

11.9 Conditions

?(): [Operator]

  ?(<condition>)

  Return a real matrix which contains only the numbers 0 or 1. For each element of the array:
  if the condition at index i,j is true then <name> [i,j]=1 else <name> [i,j]=0

?(<condition>): <constant or matrix> tabvrai : <constant or matrix> tabfaux

  Return a numerical vector which contains only the numbers 0 or 1.
  For each element of the condition: if the condition at index i,j is true then <name> [i,j]=tabvrai[i,j] else <name> [i,j]=tabfaux[j]

  All matrix must have the same size.

Example:

> mat1=matrixR[1,2,3:4,5,6:7,8,9];
mat1 double precision real matrix [ 1:3 , 1:3 ]
> mat2=matrixR[3,0,6:5,2,7:1,4,11];
mat2 double precision real matrix [ 1:3 , 1:3 ]
> q=?(mat1<5);
q double precision real matrix [ 1:3 , 1:3 ]
> writes(3*"%g "+"\n",q);
1 1 1
1 1 0
0 0 0
> m = ?(mat1>2):mat2:-1;
m double precision real matrix [ 1:3 , 1:3 ]
> writes(3*"%g "+"\n",m);
-1 -1 6
5 2 7
1 4 11
>
11.10 Conversion

**matrixR**

\[ \text{matrixR}( \text{<object identifier> } ) \]

Return a real matrix with the object identifier. The object identifier may be an array, a numerical vector or an array of numerical vectors. The object must contain only real numbers.

Example:

```maple
> _affc=1$
> // convert an array of numbers to a real matrix
> tab3=[1,2,3
:4,5,6];
```

\[ \text{tab3} \left[ 1:2, 1:3 \right] \text{ number of elements } = 6 \]

```maple
> mat3=matrixR(tab3);
mat3 \text{ double precision real matrix } \left[ 1:2 , 1:3 \right]
> afftab(mat3);
\[ \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \end{bmatrix} \]
> // convert an numerical vector to a real matrix
> v2=exp(I*v2);
```

\[ \text{v2} \text{ double precision real vector : number of elements } =10 \]

```maple
> mat2=matrixR(v2);
mat2 \text{ double precision real matrix } \left[ 1:10 , 1:1 \right]
```

**matrixC**

\[ \text{matrixC}( \text{<object identifier> } ) \]

Return a complex matrix with the object identifier. The object identifier may be an array, a numerical vector or an array of numerical vectors. The object must contain only complex numbers.

Example:

```maple
> _affc=1$
> // convert an array of numbers to a complex matrix
> tab3=[1+2*i,3+4*i,5
:2,4*i,-7+2*i];
```

\[ \text{tab3} \left[ 1:2, 1:3 \right] \text{ number of elements } = 6 \]

```maple
> mat3=matrixC(tab3);
mat3 \text{ double precision complex matrix } \left[ 1:2 , 1:3 \right]
> afftab(mat3);
\[ \begin{bmatrix} (1+i*2) & (3+i*4) & 5 \\ 2 & (0+i*4) & (-7+i*2) \end{bmatrix} \]
> // convert an numerical vector to a complex matrix
> v2=1,10;
```

\[ \text{v2} \text{ double precision complex vector : number of elements } =10 \]

```maple
> v2 = exp(I*v2);
> mat2=matrixC(v2);
mat2 \text{ double precision complex matrix } \left[ 1:10 , 1:1 \right]
```
vnumR

vnumR( <matrix> )

Return a real vector from the matrix. The matrix must have a single column.

Example :
> // convert a real matrix to a real vector
> mat1=matrixR[2:3:5];
mat1 double precision real matrix [ 1:3 , 1:1 ]
> v1=vnumR(mat1);
v1 double precision real vector : number of elements =3
> writes(v1);
+2.0000000000000000E+00
+3.0000000000000000E+00
+5.0000000000000000E+00
>

vnumC

vnumC( <matrix> )

Return a real vector from the matrix. The matrix must have a single column.

Example :
> // convert a complex matrix to a complex vector
> mat1=matrixC[1+2*i:4*i:-7+2*i];
mat1 double precision complex matrix [ 1:3 , 1:1 ]
> v1=vnumC(mat1);
v1 Double precision complex vector : number of elements =3
> writes(v1);
+1.0000000000000000E+00 +2.0000000000000000E+00
+0.0000000000000000E+00 +4.0000000000000000E+00
-7.0000000000000000E+00 +2.0000000000000000E+00
>
12 Graphics

TRIP requires the following versions of gnuplot to work:
- on Windows, version 3.7.3 or later.
- on UNIX or MacOS X, version 3.7.0 or later.
- on Android, the application droidplot.

TRIP requires the following versions of grace to work:
- On UNIX, MacOS X or Windows, version 5.1.8 or later.

The commands plot, replot, plotf, plotps, plotps_end and plotreset use grace or gnuplot depending on the global variable _graph (see Chapter 3 _graph, page 9).

12.1 plot

plot

plot(<(array of) real vec.>TX, <(array of) real vec.>TY);
plot(<(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) string> options);
Executes gnuplot or grace and plots the contents of the numerical vector TY function of TX.
The string options is directly sent to gnuplot or grace as an argument of the command plot.
Remarks: Temporary files are created and will be destroyed at the end of the session.

plot

plot(<(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ);
plot(<(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ, <(array of) string> options);
Executes gnuplot or grace and plots in 3D the contents of the numerical vector TZ function of TY and TX.

plot

plot(<string> cmd, <(array of) real vec.>TX, <(array of) real vec.>TY);
plot(<string> cmd, <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) string> options);
It executes gnuplot or grace if necessary. It sends to gnuplot or grace the command cmd and plots the contents of the numerical vector TY function of TX.

plot

plot(<string> cmd, <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ);
plot(<string> cmd, <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ, <(array of) string> options);
It executes gnuplot or grace if necessary. It sends to gnuplot or grace the command cmd and plots in 3D the contents of the numerical vector TZ function of TY and TX.

plot

plot( (<(array of) real vec.>TX, <(array of) real vec.>TY), ... );
plot( (<(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) string> options), ... );
Executes gnuplot or grace and overlays all draws of each couplet with plotting the contents of the numerical vector TY function of TX.
plot
[Procedure]
plot( (<array of> real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ), ...);
plot( (<(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ, <(array of) string> options), ...);
Executes gnuplot or grace and overlays all draws of each triplet with plotting in 3D the contents of the numerical vector TZ function of TY and TX. session.

plot
[Procedure]
plot(<string> cmd, ( <(array of) real vec.>TX, <(array of) real vec.>TY), ...);
plot(<string> cmd, ( <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ, <(array of) string> options), ...);
It executes gnuplot or grace if necessary. It sends to gnuplot or grace the command cmd and plots the contents of each couplet of the numerical vector TY function of TX.

plot
[Procedure]
plot(<string> cmd, ( <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ), ...);
plot(<string> cmd, ( <(array of) real vec.>TX, <(array of) real vec.>TY, <(array of) real vec.>TZ, <(array of) string> options), ...);
It executes gnuplot or grace if necessary. It sends to gnuplot or grace the command cmd and plots in 3D the contents of each triplet of the vectors TZ function of TY and TX.

Example :
Tracer cos(x) pour x=-pi ` a pi avec un pas de pi/100.
> x=-pi,pi,pi/100;
x Tableau de reels : nb reels =200
> y=cos(x);
y Tableau de reels : nb reels =200
> plot(x,y);
> plot(x,y,cos(x));
> plot(x,y,"notitle w points pt 5");
> plot(x,y,"title ",x","cos(x)" ");
> t=1,10;
t Tableau de reels : nb reels =10
> plot("set xrange[2:5]",t,log(t));

> t=0,pi,pi/100;
> vnumR ta[1:3];
> ta[1]=cos(t);
> ta[2]=sin(t);
> ta[3]=cosh(t);
> dim nom[1:3];
> nom[1]="title 'cos' w l";
> nom[2]="title 'sin' w l";
> nom[3]="title 'cosh' w l";
> plot((t,ta,nom));

12.2 replot
replot
[Procedure]
replot((<array of> real vec.>TX, <(array of) real vec.>TY));
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) string options);
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\);
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), (array of) string options);
replot(string \(cmd\), (array of) real vec.\(TX\), (array of) real vec.\(TY\);
replot(string \(cmd\), (array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\);
replot(string \(cmd\), (array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), (array of) string options);
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), (array of) string options);
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), (array of) string options), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), (array of) string options), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...
replot((array of) real vec.\(TX\), (array of) real vec.\(TY\), (array of) real vec.\(TZ\), ...

This command is very similar to the command \texttt{plot} but it overlays the draws on the previous draws.

It uses the same arguments as the command \texttt{plot} (see Section 12.1 \texttt{[plot]}, page 111).

\texttt{replot} \hfill \texttt{[Procedure]}

\begin{verbatim}
replot;
\end{verbatim}

Executes gnuplot or grace and send a command to redraw all graphics.

Example:
Tracer \(\cos(x)\) et \(\sin(x)\) pour \(x=-\pi\) à \(\pi\) avec un pas de \(\pi/100\) dans la même fenêtre.
\begin{verbatim}
> x=-pi,pi,pi/100;
x  Tableau de reels : nb reels =200
> y=cos(x);
y  Tableau de reels : nb reels =200
> y1=sin(x);
y1 Tableau de reels : nb reels =200
> plot(x,y);
> replot(x,y1);
> plot(x,y,y1);
> replot(x,y,sin(2.*x));
> replot(x,sin(x),"notitle");
> replot(x,y,"w points pt 5");
\end{verbatim}
12.3 plotf

plotf

plotf(<filename> filename, <integer> ncolTX, <integer> ncolTY);

It executes gnuplot or grace and plots the contents of the column ncolTY function of ncolTX in the file filename.

Example:
Afficher la troisième colonne en fonction de la première colonne du fichier tab.out
> plotf(tab.out,1,3);

Afficher la troisième colonne en fonction de la première et deuxième colonne du fichier tab.out
> plotf(tab.out,1,2,3);

12.4 plotps

plotps

plotps <filename>;

It executes gnuplot or grace and set the terminal of gnuplot or grace in postscript.

All graphics will be stored in the specified postscript file. This file will be located in the folder specified by .path.

To close the postscript file, the command plotps_end must be executed.

Example:
Tracer cos(x) pour x=-pi à pi avec un pas de pi/100 et le stocker dans le fichier res.ps.

> x=-pi,pi,pi/100;
x nb elements réels =200
> y=cos(x);
y nb elements réels =200
> plotps res.ps;
> plot(x,y);
> plotps_end;

12.5 plotps_end

plotps_end

plotps_end;

Close the file created by plotps.

For gnuplot, it sets the terminal to its default value.

Remarks:
- on Unix, MacOS X or cygwin, the default terminal is x11.
- on Windows, the default terminal is windows.
Example :
Tracer $\cos(x)$ pour $x=-\pi$ à $\pi$ avec un pas de $\pi/100$
et le stocker dans le fichier res.ps.

```plaintext
> x=-pi,pi,pi/100;
x nb elements réels =200
> y=cos(x);
y nb elements réels =200
> plotps res.ps;
> plot(x,y);
> plotps_end;
```

### 12.6 plotreset

plotreset [Procedure]

`plotreset;`

Send a command to reinitialize gnuplot or grace.

For gnuplot, it sends the command `reset`.
For grace, it sends the command `new` followed with `redraw`.

Examples :
```plaintext
> _graph=grace;
> t=0,10;
t Tableau de reels : nb reels =11
> plot(t,t);
> plotreset;
```

### 12.7 gnuplot

gnuplot [Procedure]

`gnuplot;`

```
<gnuplot command>
<gnuplot command>@<string> @<gnuplot command>
%<trip command> \ 
<trip command>
end;
```

An alternative syntax to `gnuplot; ... end;` is `gnuplot; ... gnuplot_end;`.

TRIP takes the gnuplot commands and sends them to gnuplot (line after line). Gnuplot is executed if it wasn’t started. The prompt becomes ‘`gnuplot>’ when the user could enter the gnuplot commands.

When the first character is %, then the end of this line is one or more commands trip. This trip command could follow on several lines : the last character must be \ to indicates that the command follows on the next line.

Strings declared in trip could be send to gnuplot by surrounding with the character @.

Example :
```plaintext
> gnuplot;
gnuplot> plot 'tftf' using 1:3
gnuplot> set xrange[1:10]
```
gnuplot> replot
> >
> gnuplot;
> gnuplot> set terminal macintosh singlewin
> Terminal type set to 'macintosh'
> Options are 'nogx singlewin novertical'

> gnuplot> %plot(t,log(t),"notitle");

> gnuplot> %replot(t,exp(t),"notitle w points pt 5");

> gnuplot> set terminal macintosh multiwin
> Terminal type set to 'macintosh'
> Options are 'nogx multiwin novertical'

> ch="title 'sinus'";
> ch = "title 'sinus'"
> gnuplot;
> gnuplot> plot sin(x) @ch@
> gnuplot> end;

12.8 grace

grace

grace;
<commande grace>
<commande grace>@<string> @<commande grace>
%<commande trip> \ 
<commande trip>
end;

TRIP takes the grace commands and sends them to grace (line after line). Grace is executed if it wasn’t started. The prompt becomes ‘grace>’ when the user could enter the grace commands.

When the first character of %, then the end of this line is one or more commands trip. This trip command could follow on several lines: the last character must be \ to indicates that the command follows on the next line.

Strings declared in trip could be send to grace by surrounding with the character @.

Example :
> grace;
grace> grace> with g0
grace> read block "/USER/toto"
grace> block xy "1:2"
grace> read block "/USER/toto"
grace> block xy "1:3"
grace> redraw
grace> title "2courbes"
grace> %t=1,10;
Tableau de reels : nb reels = 10

grace> %msg "deux lignes\nfin deligne";
deux lignes fin deligne
grace> end;
> >
13 Communications

TRIP have different communication tools:
− Communication with other computer algebra systems.
− Execution of functions located in external dynamic libraries.

TRIP communicates with other computer algebra systems on the same computer. These computer algebra systems must be able to import and export computations using the MathML 2.0 protocol.

A global common session is available for each computer algebra system. Several different session for each computer algebra system may be defined.

13.1 Maple

TRIP communicates with Maple\(^1\) on all operating systems which could run maple.

The required version of maple must be equal or greater than 11.

13.1.1 maple\_put

\texttt{maple\_put} \hspace{1cm} \textbf{[Procedure]}

\texttt{maple\_put( \textless \text{object identifier} \textgreater \textit{id});}

Send the object identifier to the global maple session. This starts maple if it isn’t running.

Example:
\begin{verbatim}
> _affdist=1$
> s:=x+1;
s(x) = 1 + x
> maple\_put(s);
> maple;
> s;
1 + x
> end;
>
\end{verbatim}

\texttt{maple\_put} \hspace{1cm} \textbf{[Procedure]}

\texttt{maple\_put( \textless \text{Maple client} \textgreater \textit{session} , \textless \text{object identifier} \textgreater \textit{id});}

Send the object identifier to the maple session specified by \textit{session}, which must be started previously.

Example:
\begin{verbatim}
> _affdist=1$
> fm:=maple;
> z:=1;

> end;
> s:=x+1;
s(x) = 1 + x
> maple\_put(fm,s);
> maple(fm);
> z+s;
2 + x
> end;
>
\end{verbatim}

\(^1\) Maple is a registered trademark of Waterloo Maple Inc.
13.1.2 maple_get

maple_get

maple_get( <object identifier> id);  
Get the value of an object identifier from the global Maple session.

Example :
> _affdist=1$
> maple;
> s:=1+y;

> end;
> maple_get(s);
> s;

s(y) = 1 + y

maple_get

maple_get( <Maple client> session , <object identifier> id);  
Get the value of an object identifier from the specified Maple session.

Example :
> _affdist=1$
> fm=maple;
> s:=1+y;

> end;
> maple_get(fm, s);
> s;

s(y) = 1 + y

13.1.3 maple

maple

maple;  
<commande maple>

%<commande maple> \ 
<commande maple>
%<commande trip> \ 
<commande trip>

end;

TRIP accepts maple commands and send them to the global Maple session (line after line).
maple will be started if the global session is not already started. The prompt becomes ‘maple>’ when you could enter maple commands.

This maple command could continue on several lines : To continue the command on a new line, the last character of the line must be a \ .

If the first character is a %, then the end of the line will be considered as a TRIP command and not as a maple command. This TRIP command could continue on several lines : To continue the command on a new line, the last character of the line must be a \ .

The function maple_get retrieves objects (series, vectors, ...) from maple. The command maple_put sends objects to maple.
Example:
> maple;
> s:=gcd((x+1)*(x-1),(x-1));
\[ s := x - 1 \]
> %maple_get(s);
> %s;
s(x) =
- 1
+ 1*x
> end;
>
maple

This command is similar to the previous one but it starts a new session every time. It returns an object which specifies the session for the commands maple_put and maple_get.

An existing session could be continued by the command maple(<Maple client>); ... end;

The session may be closed by the command delete(<Maple client>);

13.2 Mathematica

TRIP communicates with Mathematica\(^2\) on all operating systems which could run mathematica.

The required version of mathematica must be equal or greater than 9.

13.2.1 mathematica_put

mathematica_put

This command continues the existing session.

Example:
> fm=maple;
> f:=gcd((x+1)*(x-1),(x-1));
\[ f := x - 1 \]
> g:=f+2;
\[ g := x + 1 \]
> end;
> maple_get(fm,g);
> delete(fm);

\(^2\) Mathematica is a registered trademark of Wolfram Research Inc.
Example:
> _affdist=1$
> s=1+x;
s(x) = 1 + x
> mathematica_put(s);
> mathematica;
> s
1 + x
> end;
>
mathematica_put
mathematica_put(<Mathematica client> session, <object identifier> id);
Send the object identifier to the mathematica session specified by session, which must be started previously.

Example:
> _affdist=1$
> fm=mathematica;
> z:=1;
> end;
> s=1+x;
s(x) = 1 + x
> mathematica_put(fm,s);
> mathematica(fm);
> z+s;
> end;
>
13.2.2 mathematica_get

mathematica_get
mathematica_get(<object identifier> id);
Get the value of an object identifier from the global mathematica session.

Example:
> _affdist=1$
> mathematica;
> s:=1+y;
> end;
> mathematica_get(s);
> s;
s(y) = 1 + y
>
mathematica_get
mathematica_get(<Mathematica client> session, <object identifier> id);
Get the value of an object identifier from the specified Mathematica session.

Example:
> _affdist=1$
> fm=mathematica;
> s:=1+y;
> end;
> mathematica_get(fm, s);
> s;
s(y) = 1 + y

13.2.3 mathematica

mathematica

[Procedure]

mathematica;
<commande mathematica>
%%<commande mathematica> \n
<commande mathematica>
%%<commande trip> \n
<commande trip>
end;

TRIP accepts mathematica commands and send them to the global session of Mathematica (line after line). mathematica will be started if the global session not already started. The prompt becomes ‘mathematica>’ when you could enter mathematica commands.

This mathematica command could continue on several lines: To continue the command on a new line, the last character of the line must be a \.

If the first character is a %, then the end of the line will be considered as a TRIP command and not as a mathematica command. This TRIP command could continue on several lines: To continue the command on a new line, the last character of the line must be a \.

The function mathematica_get retrieves objects (series, vectors, ...) from mathematica. The command mathematica_put sends objects to mathematica.

Example:
> mathematica;
> s=PolynomialGCD[(x+1)*(x-1),(x-1)]
-1 + x
> %mathematica_get(s);
> %s;
s(x) =
  1
  + 1*x
> end;

> mathematica

[Function]

<Mathematica client> = mathematica;
<commande mathematica>
end;

This command is similar to the previous one but it starts a new session every time. It returns an object which specifies the session for the commands mathematica_put and mathematica_get.

An existing session could be continued by the command mathematica(<Mathematica client>);

The session may be closed by the command delete(<Mathematica client>);.
mathematica
mathematica(<Mathematica client>);
<commande mathematica>
end;

This command continues the existing session.

Example:
> fm=mathematica;
> f=PolynomialGCD[(x+1)*(x-1),(x-1)]
  -1 + x
> g=f+2
  1 + x
> end;
> mathematica_get(fm,g);
> delete(fm);
>
13.3 Communications with other computer algebra systems

TRIP communicates with other computer algebra systems on the same or remote computer. These computer algebra systems must be able to compatible with the "Symbolic Computation Software Composability Protocol" (SCSCP) version 1.3 (http://www.symcomp.org/). It supports the OpenMath symbols from the scscp1 content dictionary (http://www.win.tue.nl/SCIEnce/cds/scscp1.html) and from the scscp2 content dictionary (http://www.win.tue.nl/SCIEnce/cds/scscp2.html). The list of OpenMath content dictionary is given in the appendice (see Appendix A [Supported OpenMath Content Dictionaries], page 211).

Multiple connections could be opened at the same time. TRIP could run in client or server mode.

Before using any functions from this module, you must execute the following command in the TRIP session. This command defines many subroutines which contain the definition of the Openmath symbols.

include libscscpserver.t;

13.3.1 SCSCP server

In order to start the SCSCP server of TRIP, you should execute the following command in a TRIP session:

include libscscpserver.t;
port = 26133;
%scscp_runserver[port];

The value of the variable of port may be changed. The default is 26133 for SCSCP servers.

The global variables of TRIP, such as _modenum, could be modified before starting the SCSCP server.

The file libscscpserver.t exports some symbols from the standard OpenMath content dictionaries. New symbols could be exported to add new functionalities to the SCSCP server. The function scscp_map_macroassymbolcd, defined in the file libscscpserver.t, must be used.
scscp_map_macroassymbolcd (string macroname, string cdname, string symbolname);

Associate the symbol symbolname of the content dictionary cdname with the macro macroname. When this symbol is found by the SCSCP server or the SCSCP client in a exchanged message, then this macro is executed.

Example:
include libscscp cooperating.t;
_modenum=NUMRATMP;
macro myscscp_evalmul(P1, P2, value)
{
    t=value,value;
    q=evalnum(P1*P2,REAL, (x,t));
    return q[1];
};
scscp_map_macroassymbolcd("myscscp_evalmul",
    "SCSCP_transient_1","scscp_muleval");
%scscp_runserver[26133];

13.3.1.1 scscp_disable_cd
see Section 13.3.2.7 [scscp_disable_cd], page 129.

13.3.1.2 Content dictionary scscp_transient_1
The SCSCP server of TRIP allow the following operations through the content dictionary scscp_transient_1:

- serversupportbinary. It returns 1 if the SCSCP server supports the binary encoding for the OpenMath objects.
- serverdisablecd(CDname). It specifies that the SCSCP server must not use the content dictionary CDname for the OpenMath objects inside the communications with this client.

Example:
> with(SCSCP);:
> with(Client);
> cookie := StorePersistent("localhost:26133", x*y+1);
    cookie := "Temp0ID1@localhost:26133"

> Retrieve("localhost:26133", cookie);
Error, (in SCSCP:-Client:-ExtractAnswer) unsupported_CD, <OMS cd = 'polyr' name = 'term'/>
> CallService("localhost", "scscp_transient_1",
    "serverdisablecd", ["polyr"]);
> CallService("localhost", "scscp_transient_1",
    "serverdisablecd", ["polyu"]);
> Retrieve("localhost:26133", cookie);
    1.0000000000000000 + 1.0000000000000000 x y

13.3.2 SCSCP client
TRIP communicates with other computer algebra systems providing a SCSCP server.

13.3.2.1 scscp_connect

scscp_connect (string computename, integer port);
Connect to the computer algebra system on the remote computer using the specified port. This function returns an sscp client object which manages that connection.

Example:
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
sc = client SCSCP connecte au serveur SCSCP localhost
> scscp_close(sc);
> stat(sc);
client SCSCP sc : deconnecte

13.3.2.2 scscp_close

scscp_close(scscp_close(<sscp client> sc));
Close the connection to the computer algebra system specified by sc.

Example:
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
sc = client SCSCP connecte au serveur SCSCP localhost
> scscp_close(sc);
> stat(sc);
client SCSCP sc : deconnecte

13.3.2.3 scscp_put

<remote object> scscp_put(scscp_put(<sscp client> sc, <operation> x));
scscp_put(<sscp client> sc, <operation> x, <string> storeoption);
Send the object identifier x to the computer algebra system using the sscp client sc, previously opened with scscp_connect. This function returns a remote object.

If storeoption isn’t specified, then its value is "persistent".
The string storeoption must be one of the following values:
- "persistent": the remote object will be kept by the server after the connection is closed.
- "session": the remote object will be destroyed by the server at the end of the connection

Example:
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
sc = client SCSCP connecté au serveur SCSCP localhost
> r=1+x;
r(x) =
    1
    +
    1*x

> remoter=scscp_put(sc, r);
remoter = objet distant "TempOID1@localhost:26133"
> scscp_close(sc);

13.3.2.4 scscp_get

scscp_get
scscp_get( <remote object> remoteobjectid );
Get the value of the object identifier from the remote computer algebra system using the object remoteobjectid, previously returned by scscp_put.

Example :
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
sc = client SCSCP connecte au serveur SCSCP localhost
> s=(1+x+y)**2;
s(x,y) =
    1
    +
    2*y
    +
    1*y**2
    +
    2*x
    +
    2*x*y
    +
    1*x**2

> remoteS = scscp_put(sc, s);
remoteS = objet distant "TempOID1@localhost:26133"
> q=scscp_get(remoteS);
q(x,y) =
    1
    +
    2*y
    +
    1*y**2
    +
    2*x
    +
    2*x*y
    +
    1*x**2

> scscp_close(sc);

13.3.2.5 scscp_delete

scscp_delete
scscp_delete( <remote object> remoteobjectid );
Destroy the value of the object identifier from the remote computer algebra system using the object remoteobjectid, previously returned by scscp_put.
Example:
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
scc = client SCSCP connecte au serveur SCSCP localhost
> s=(1+x+y)**2;
s(x,y) =
  1
  + 2*y
  + 1*y**2
  + 2*x
  + 2*x*y
  + 1*x**2

> remoteS = scscp_put(sc, s);
remoteS = objet distant "TempOID1@localhost:26133"
> delete(remoteS);
> scscp_close(sc);

13.3.2.6 scscp_execute

scscp_execute

scscp_execute( <scscp client> sc, <string> returnoption , <string> CDname , <string> remotecommand , <operation> , ... );

It executes the command remotecommand of the OpenMath CD CDname on the remote computer algebra system specified by sc. The parameters are specified just after the name of the command.

The string returnoption must be one of the following values:

− "cookie" : a remote object is returned and the result of the execution will be kept on the SCSCP server.
− "object" : the result of the execution is returned as an OpenMath object and this object is evaluated.
− "nothing" : no result or remote object are returned.

Example:
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
scc = client SCSCP connecte au serveur SCSCP localhost
> q = scscp_execute(sc,"object", "SCSCP_transcient_1", "SCSCP_MUL", 1+7*x,2+3*x);
q(x) =
  2
  + 17*x
  + 21*x**2

> qr = scscp_execute(sc,"cookie", "SCSCP_transcient_1", "SCSCP_MUL", 1+7*x,2+3*x);
13.3.2.7 scscp_disable_cd

scscp_disable_cd( <string> CDname , );

It disable the content dictionary CDname. This dictionary will not be used for the OpenMath output and for the communication with the remote computer algebra system.

Example :
> include libscscpserver.t;
Loading the SCSCP client/server library...
Registering the OpenMath CD...
> port=26133;
port = 26133
> sc=scscp_connect("localhost", port);
sc = client SCSCP connecte au serveur SCSCP localhost
> scscp_disable_cd("polyr");
> scscp_disable_cd("polyu");
> s=(1+x+y)**2$
> q=scscp_put(sc, s);
q = objet distant "TempOID1@localhost:26133"
> scscp_close(sc);

13.4 Dynamic library

TRIP can load a dynamic library and execute functions of this library. The dynamic libraries have the extensions .so, .dll, .dylib depending on the operating system. The function prototypes is required in order to perform the conversion of the parameters.

13.4.1 extern_function

extern_function( <filename> filelib, <string> declfunc );

It loads the dynamic library filelib. It adds the specific extension (.dll, .so, .dylib) of the operating system. It checks the availability of the function in this library. The function could be called as any other standard function of TRIP.

It checks the validity of the prototype of the function specified by declfunc. The prototype must be written in C language. The implementation function could be written in another language (fortran, ...).

If the dynamic library depends on other libraries, then it requires to load the other libraries with the command extern_lib.

By default, all arguments are input only. To specify that arguments are input-output, it must be specified in the string declinout. This string must contain the same number of elements as the number of parameter of the function. Each element are separated with a comma. An element could have the following values :

− in : the parameter is an input only.
− out : the parameter is an output only.
− inout : the parameter is an input and output.
For output arguments, an object identifier must be given to the function to get the value.

Remarks: the function could not have structure or derived type.

Example:
> extern_function("libm", "double j0(double);");
> r = j0(0.5);
> r = 0.9384698072408129

Example:
/* Appel de sa propre librairie libtest1 contenant test1.c */
> !$"cat test1.c"
#include <math.h>

double mylog1p(double x)
{
    return log1p(x);
}

void mylog1parray(double tabdst[], double tabsrc[], int n)
{
    int i;
    for(i=0; i<n; i++) tabdst[i]=log1p(tabsrc[i]);
}

> extern_function(libtest1,"double mylog1p(double x);"�);
> mylog1p(10);
> log(11);
> extern_function(libtest1,
    "void mylog1parray(double tabdst[], double tabsrc[], int n);",
    "inout,in,in");
> t=1,10;
t Tableau de reels : nb reels =10
> vnumR res; resize(res,10);
> mylog1parray(res,t,10);
> writes(t,res);
+1.0000000000000000E+00 +6.9314718055994529E-01
+2.0000000000000000E+00 +1.09861228886681098E+00
+3.0000000000000000E+00 +1.3862943611989066E+00
+4.0000000000000000E+00 +1.609437912424341003E+00
+5.0000000000000000E+00 +1.7917594692280550E+00
+6.0000000000000000E+00 +1.9459101490553132E+00
+7.0000000000000000E+00 +2.0794415416798357E+00
+8.0000000000000000E+00 +2.1972245773362196E+00
+9.0000000000000000E+00 +2.3025850929940459E+00
+1.0000000000000000E+01 +2.3978952727983707E+00

13.4.2 extern_lib

extern_lib( <filename> filelib);
Load the dynamic library `filelib`. It adds the specific extension (.dll, .so, .dylib) of the operating system. This function is used to load libraries required by other libraries.

Example:
```plaintext
> extern_lib("libm");
> extern_function("libm", "double j0(double)");
> r = j0(0.5);
r = 0.9384698072408129
```

### 13.4.3 extern_type

**extern_type**

```plaintext
extern_type <string> type;
```

Declares an external type `type`. The external functions can use or return pointer to objects of this type. `type` could be a C data structure whose fields may be read or modified.

Example:
```plaintext
> extern_type("t_calcephbin");
> extern_function("libcalceph",
    "t_calcephbin* calceph_open(const char *filename);";
> eph = calceph_open("inpop06c_m100_p100_littleendian.dat");
```

### 13.4.4 extern_display

**extern_display**

```plaintext
extern_display;
```

Display the list of external types or functions previously loaded with the command `extern_function` or `extern_type`.

Example:
```plaintext
> extern_function("libm", "double j0(double)");
> extern_function("libm", "double j1(double)");
> extern_display;
Liste des fonctions externes :
double j1(double);
double j0(double);
```
14 Macros

14.1 Declaration

```
macro [Procedure]
macro <name> [ <list of parameters> ] { <body> };
macro <name> { <body> };
MACRO <name> [ <list of parameters> ] { <body> };
MACRO <name> { <body> };
private macro <name> [ <list of parameters> ] { <body> };
private macro <name> { <body> };
```

Define a macro with 0 or more parameter and a body of trip code.

The list of parameters are separated by a comma. The parameter must a name. There is no limitation with the number of parameters. The macros support recursivity but a limit exists (often 70 times).

The last parameter could be ... to specify that this macro could receive one or more optional argument when it's called. To access to each optional argument, the keyword `macro_optargs[]` is used. To acces to the optional argument of index j, you just write `macro_optargs[j]`. To know the number of provided optional argument, the function `size(macro_optargs)` must be used.

The values corresponding to the ... could be given as an argument to the execution of a macro using `macro_optargs`.

The execution of the macro could be immediately stopped using the `stop` command (see Section 15.1.5 `stop`, page 141).

The macro may define local object identifiers, See Section 2.3.1 `private`, page 7.

The visibility of a macro is restricted to the source file which contains it if the declaration is prefixed by the keyword `private`. Only other macros of the same file can call it.

The following example shows the definition of a macro `a` which assigns to `r` the sum of `x` and `y`. `x` and `y` are transmitted by arguments. The macro `b` displays the contents of the current directory.

```
Example :
> macro a [x,y]
{r = x+y;};
> macro b { ! "ls";};
> macro macvar[x,y,...]
{ 
s=x+y;
  l = size(macro_optargs);
  msg("number of optional argument %g\n", l);
  for j=1 to l { msg("The optional argument %g :", j); macro_optargs[j]; };}
> %macvar[P1,P2];
s(P1,P2) =
  1*P2
+     1*P1
l =
   0
```
number of optional argument 0
> %macvar[P1,P2,"arg1", 3, 5, 7];
s(P1, P2) =
    1*P2 + 1*P1
l = 4
number of optional argument 4
The optional argument 1 :macro_optargs[1] = "arg1"
The optional argument 2 :macro_optargs[2] = 3
The optional argument 4 :macro_optargs[4] = 7

> macro macopt[x,y,...]
{
    %macvar[x, macro_optargs];
};

14.2 Execution

Execute a macro. The list of parameters is a list of parameters separated by a comma.
The parameters can be object identifier, series (result of a computation), strings of characters, arrays or numerical vectors.
The parameter can be transmitted by value or reference.
Parameters given by value can only be a result of a computation, strings of characters, arrays or numerical vectors.
Parameters given by reference can be any object identifier or an element of array. To give a parameter by reference, it must be surrounded with additional brackets []. In this case, the object identifier can be modified or created during the execution of the macro.
The execution of the macro returns a value if the command return has been called in the body of the macro. The execution of the macro continues after the command return.
The execution of the macro could be immediately stopped using the stop command (see Section 15.1.5 [stop], page 141).

Execution of the macro a and b:
Example :
> %a[1,5+3*6];
/*ou si S = x + y*/
> %a[S, 5];
>%b;
>%c[[t], [u[5]], 1+x];
Usage des chaines de caractères:
> macro nomfichier[name,j] {msg name + str(j);};
> %nomfichier["file_", 2];
file_2
> ch="file_1.2.";
Remarks: The parameters (x, y,...) keep their values. To give an expression, it mustn’t be surrounded with brackets [].

Example:
- If x is defined, its value will be kept when the macro is executed. If x is not defined, its value will be the one at the end of the macro execution.

Example:
If we want to add 'x' to 'S' variable:
> %a[S,[x]];  
est will result in r(x,y) = 1*y + 2*x

- Declaration of a matrix. Identificateur R1 does not exist.
/* creation of a matrix */
>macro matrix_create[_nom, _nbligne, _nbcol]  
{  
dim _nom[0:_nbligne, 0: _nbcol];  
for iligne=1 to _nbligne {  
  for icol=1 to _nbcol {_nom[iligne,icol]=0$};};  
/* set lines and columns to 0 */  
_nom[0, 0]=_nbligne$  
_nom[0, 1]=_nbcol$  
};  
> %matrix_create[[R1],n,3];  

Le tableau sera alors créé et initialisé.
En sortie de la macro, _nom, _nbligne, _nbcol n’existeront plus.

return  
[Procedure]
return (<operation> );
return <operation> ;
Return the result of an operation when the macro is executed.
The execution of the macro continues after the command return. Only the stop command stops the execution of the macro.
return (<operation> , <operation> , ...);
Return using a one-dimensional array the list of operations as the result of the execution of the macro. The first index of the array is 1.

Example:
> _affdist=1$  
> macro func1 { s=1+y$ return s; };  
> b=%func1;

b(y) = 1 + y
> macro func2 { v=1,10$ ch="file1"$ return (v, ch); };  
> (a, s)=%func2;

a

a double precision real vector : number of elements =10
> s;

s = "file1"
> t=%func2;
t [1:2 ] number of elements = 2

> afftab(t);
t[1] = t double precision real vector : number of elements =10

t[2] = "file1"

14.3 List of macros

@ [Procedure]

@;
Display the list of the defined macros.

Example :
> @;
Voici le nom des macros que je connais:
a [x ,y ]
b

14.4 Display the body

affmac [Procedure]

affmac <macro>;
Display the body (trip code) of the macro.

Example :
> affmac b;
mon nom est : b
!"ls";

14.5 Deletion

14.5.1 effmac

effmac [Procedure]
effmac <macro>;
Delete a macro.

Example :
> macro a[x] { return (x*2);};
> @;
Voici le nom des macros que je connais :
a [ x]
> effmac a;
> @;
je ne connais aucune macro

14.5.2 effmacros

effmacros [Procedure]
effmacros;
Remove all defined macros.
Example:
> macro a[x] { return (x*2);};
> macro b {!"ls";};
> @;
Voici le nom des macros que je connais:
a [ x]
b []
> effmacros;
> @;
je ne connais aucune macro

14.6 How to redefine a macro?
We write a new macro using the command macro.
Example:
> macro a [n] {n;};
> @;
Voici le nom des macros que je connais:
a [ n]
> affmac a;
mon nom est : a
n;

14.7 How to save a macro?
To save the macros on disk, the best solution is to use a text editor such as vi, emacs or nedit.
The name of the file which contains trip code should be ended with .t.
15 Loops and conditions

15.1 Loops

15.1.1 while

while (<condition>) do { <body> };

Execute the body loop while the condition is true.

The while loop is more yielding than for loop because condition could be sophisticated. A while loop could be immediately stopped using the stop command. For the condition description, See Section 15.2 [condition], page 141.

The loop may define local object identifiers, See Section 2.3.1 [private], page 7. They will destroyed at the end of each iteration.

Example :
> // It displays the numbers from 1 to n:
> p=1$
> while(p<=5) do { p; p=p+1$ };
> p = 1
> p = 2
> p = 3
> p = 4
> p = 5
>

15.1.2 for

for <nom> = <real> to <real> { <body> };
for <nom> = <real> to <real> step <real> { <body> };

Execute the loop "for ... until ..." (similar to the for loop in pascal, C or fortran languages).

The argument after step is the loop step. A for loop could be immediately stopped using the stop command.

The loop may define local object identifiers, See Section 2.3.1 [private], page 7. They will destroyed at the end of each iteration.

Remarks :
− The loop counter could not be i because i is the mathematical notation for complex numbers.
− The value of the loop counter could not be modified in the body loop.

Example :
> for p = 1 to 5 step 2 {p; };
> p = 1
> p = 3
> p = 5
> for p = 5 to -1 step -2 {p; };
> p = 5
> p = 3
> p = 1
p = -1

The loop for may be parallelized using the terminology OpenMP. TRIP does not parallelize the loop if global objects are written or if functions which modified the global state of TRIP are called. In these cases, a warning is printed.

If the keyword `distribute` is used, the parallelization is performed on all computing nodes if the application tripcluster runs this code.

```plaintext
for /*!trip omp parallel for */
   for <nom> = <real> to <real> { <body> }

for /*!trip omp parallel for distribute */
   for <nom> = <real> to <real> { <body> }
```

Example:

```plaintext
s=(1+x+y+z+t)**(20)
```

```plaintext
# Example
s=(1+x+y+z+t)**(20)
dim f[1:8];
n=size(f);
// parallel execution
time_s;
/*!trip omp parallel for */
for p = 1 to n { f[p]=s/p; }
time_t;
utilisateur 00.119s - reel 00.023s - systeme 00.019s - (594.73% CPU)
// sequential execution
```
if (mod(j,2)==0) then { return a; } else { return -a; };
}
s = -75

15.1.4 prod

sum
prod <nom> = <real> to <real> { <body> };
prod <nom> = <real> to <real> step <real> { <body> };

Execute the loop "prod ... until ...". It replaces the for loop statement "s=1$ for j=1 to n { s=s*...$}". The value returned by the statement return is used for the product.

The argument after step is the loop step. A prod loop could be immediately stopped using the stop command.

The loop may define local object identifiers, See Section 2.3.1 [private], page 7. They will destroyed at the end of each iteration.

Remarks:
− The loop counter could not be i because i is the mathematical notation for complex numbers.
− The value of the loop counter could not be modified in the body loop.

Example :
> p = prod j = 1 to 10 { return j$ };
p = 3628800

15.1.5 stop

stop

Immediately stops the execution of a for or while loop or of a macro.

Remarks: if stop is used outside a for or while loop or macro statement, a warning message is displayed but the execution continues.

Example :
for p = 1 to 5 { if (p>3) then {stop;} fi; p;};
1
2

15.2 condition

15.2.1 if

if

Execute the first body (after the then statement) if the condition is true. The second body is executed if the condition is false.

Example :
> if (n == 2) then {msg "VRAI";} else {msg"FAUX";};
FAUX
> n=2;
if (( n >= 0) && (n < 3)) then {msg "VRAI";} else
{msg "FAUX";};

Remarks: All object identifiers should be initialized before the execution of the condition. Because otherwise, uninitialized object identifiers will be created as variables and the condition will return false.

15.2.2 Comparison operators

!=  
<operation> != <operation>
This test returns true if the two operations are not equal.
Remarks: the test could be performed on polynomials.
<real vec.> != <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

Example :
> if (n != 2) then {} else {};
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t!=r);
> q double precision real vector : number of elements =11
> q = ?(t!=5);
> q double precision real vector : number of elements =11

==  
<operation> == <operation>
This test returns true if the two operations are equal.
If the two operands are the values NaN (not a number) , then the test returns false.
Remarks: the test could be performed on polynomials.
<real vec.> == <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

Example :
> if (n == 2) then {} else {};
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t==r);
> q double precision real vector : number of elements =11
> q = ?(t==5);
> q double precision real vector : number of elements =11
>
Chapter 15: Loops and conditions

<
<real> < <real>
This test returns true if the first real number is less than the second number.
<real vec.> < <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

Example :
> n=3$
> if (n < 2) then {} else {};
>
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t<r);
q double precision real vector : number of elements =11
> q = ?(t<5);
q double precision real vector : number of elements =11
>
>
>
>
>
>

>
<real> > <real>
This test returns true if the first real number is greater than the second number.
<real vec.> > <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

Example :
> n=3$
> if (n > 2) then {} else {};
>
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t>r);
q double precision real vector : number of elements =11
> q = ?(t>5);
q double precision real vector : number of elements =11
>

>
<=
<real> <= <real>
This test returns true if the first real number is less or equal than the second number.
<real vec.> <= <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.
Example :
> n=3$
> if (n <= 2) then {} else {};
>
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t<=r);
  q double precision real vector : number of elements =11
> q = ?(t<=5);
  q double precision real vector : number of elements =11
>
$[Operator]$

$>=$
<real> >= <real>
This test returns true if the first real number is greater or equal than the second number.
<real vec.> >= <real vec.>
This test could be used with the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62). It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

Example :
> n=3$
> if (n >= 2) then {} else {};
>
> // Condition between two vectors
> t = 0,10$
> r = 10,0,-1$
> q = ?(t>=r);
  q double precision real vector : number of elements =11
> q = ?(t>=5);
  q double precision real vector : number of elements =11
>
$[Operator]$

$&&$
<condition> && <condition>
This test returns true if the two conditions are true.
This test could be apply on numerical vectors when the operator ?:: (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62) are used. It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.
The test in parenthesis are required if the tests include several && and || .

Example :
> if ((x==2)&&(y==3)) then {} else {};
>
> // Example using two vectors
> t=0,10;
  t double precision real vector : number of elements =11
> r=10,0,-1;
  r double precision real vector : number of elements =11
Chapter 15: Loops and conditions

> q=?((t>r) && (t!=5));
q  double precision real vector : number of elements =11
>
|| [Operator]
<condition> || <condition>

This test returns true if at least one condition is true.

This test could be apply on numerical vectors when the operator (see Section 10.11 [Conditions], page 93) or the command select (see Section 10.7 [Extraction], page 62) are used. It returns a numerical vector which contains only 0 or 1. It compares each element of the two vectors. The numerical vector of real numbers must have the same size.

The test in parenthesis are required if the tests include several && and ||.

Example :
> if ((x==2)||(y==3)) then {} else {};
>
> // Example using two vectors
> t=0,10;
> r=10,0,-1;
> q=?((t>r) || (t!=5));

15.2.3 switch

switch

switch ( <operation> expr )
{
    case <operation> : { <body> };
    case <operation>, ..., <operation> : { <body> };
    else { <body> }
};

The instruction switch control complex conditional on the same value and branching operations.

expr can be a string, a serie or a constant. The value of expr are tested successively with each value of the case statements with the operator ==. When expr is equal to one of the values of a case statement, the body statement after this case is executed. The statement else is executed if no case constant-expression is equal to the value of expr. The statement else { <body> } is optional and has no semi-column after its body.

Example :

n=0;
z=0;
j=5;

switch( j )
{
    case -1 : { n=n+1; }; 
    case 0,1 : { z=z+1; }; 
}
case "mystring" : { msg "j is a string"; }
else { msg "j is invalid"; }
};
16 Library

16.1 Lapack

16.1.1 AX=B Solution

Real case for general matrices, (see [lapack_dgesv], page 147)
- Real case for band matrices, (see [lapack_dgbsv], page 148)
- Real case for tridiagonal matrices, (see [lapack_dgtsv], page 148)
- Real case for symmetric matrices, (see [lapack_dsyev], page 149)
- Real case for symmetric definite positive matrices (see [lapack_dposv], page 149)
- Real case for symmetric definite positive band matrices (see [lapack_dpbsv], page 150)
- Real case for symmetric definite positive tridiagonal matrices (see [lapack_dptsv], page 150)

Complex case for general matrices, (see [lapack_zgesv], page 151)
- Complex case for band matrices, (see [lapack_zgbsv], page 151)
- Complex case for tridiagonal matrices, (see [lapack_zgtsv], page 152)
- Complex case for symmetric matrices, (see [lapack_zsysv], page 153)
- Complex case for hermitian matrices, (see [lapack_zhesv], page 153)
- Complex case for hermitian definite positive matrices (see [lapack_zposv], page 154)
- Complex case for hermitian definite positive band matrices (see [lapack_zpbsv], page 154)
- Complex case for hermitian definite positive tridiagonal matrices (see [lapack_zptsv], page 155)

lapack_dgesv

[Function]

`lapack_dgesv(<real matrix> A ,<real matrix> B)`

It computes the solution to a real system of linear equations AX=B, where A is an N-by-N matrix and X and B are N-by-NRHS matrices. It uses the LU decomposition.

More information available at [http://www.netlib.org/lapack/double/dgesv.f](http://www.netlib.org/lapack/double/dgesv.f)

Example:

```plaintext
> // resolves a real system of linear equations AX=B
> // where A is a square matrix
> _affc=1$
> A = matrixR[22.22, -11.11:
-35.07, 78.01]$  
> B = matrixR[-88.88:
382.18]$  
> S = lapack_dgesv(A, B)$
> afftab(S);
[ -2]
[  4]
> 
```
lapack_dgbsv

lapack_dgbsv

lapack_dgbsv(<real matrix> A ,<real matrix> B)

It computes the solution to a real system of linear equations AX=B, where A is a band matrix of order N, and X and B are N-by-NRHS matrices. It uses the LU decomposition.

More information available at http://www.netlib.org/lapack/double/dgbsv.f

Example :
> // resolves a real system of linear equations AX=B
> // where A is a square matrix
> _affc=1$
> A = matrixR[
3.0, 2.1, 0, 0, 0:
3.4, 2.3, -1.0, 0, 0:
0, 3.6, -5.0, 1.9, 0:
0, 0, 7.0, -0.9, 8.0:
0, 0, 0, -6.0, 7.1$]
> B = matrixR[
2.7:
-0.5:
2.6:
0.6:
2.7]$]
> S = lapack_dgbsv(A, B);
> afftab(S);

lapack_dgtsv

lapack_dgtsv

lapack_dgtsv(<real matrix> A ,<real matrix> B)

It computes the solution to a real system of linear equations AX=B, where A is an N-by-N tridiagonal matrix and X and B are N-by-NRHS matrices.

More information available at http://www.netlib.org/lapack/double/dgtsv.f

Example :
> // example of dgtsv routine
> _affc=1$
> A = matrixR[
3.0, 2.1, 0, 0, 0:
3.4, 2.3, -1.0, 0, 0:
0, 3.6, -5.0, 1.9, 0:
0, 0, 7.0, -0.9, 8.0:
0, 0, 0, -6.0, 7.1$]
> B = matrixR[
2.7:
-0.5:
2.6:
0.6:
2.7]$]
> S = lapack_dgtsv(A, B);
> afftab(S);
lapack_dsysv

lapack_dsysv

lapack_dsysv(<real matrix> A ,<real matrix> B)

It computes the solution to a real system of linear equations \( AX=B \), where \( A \) is an N-by-N symmetric matrix and \( X \) and \( B \) are N-by-NRHS matrices.

More information available at http://www.netlib.org/lapack/double/dsysv.f

Example :

> // resolves a real system of linear equations \( AX=B \)
> // where A is a symmetric matrix
> _affc=1$
> A = matrixR[  
>   -1.81, 2 :
>   2, 1.15]$  
> B = matrixR[  
>   0.19:
>   3.15]$  
> S=lapack_dsysv(A,B)$  
> afftab(S);  
> [ 1]  
> [ 1]  
>

lapack_dposv

lapack_dposv

lapack_dposv(<real matrix> A ,<real matrix> B)

It computes the solution to a real system of linear equations \( AX=B \), where \( A \) is an N-by-N symmetric positive definite matrix and \( X \) and \( B \) are N-by-NRHS matrices.

More information available at http://www.netlib.org/lapack/double/dposv.f

Example :

> // resolves a real system of linear equations \( AX=B \)
> // where A is a symmetric positive definite matrix
> _affc=1$
> A = [  
>   4.16, -3.12:  
>   -3.12, 5.03]$  
> B = matrixR[  
>   10.40:
>   -13.18]$  
> S=lapack_dposv(A,B)$  
> afftab(S);  
> [ 1]  
> [ - 2]  
> >
lapack_dpbsv

lapack_dpbsv

lapack_dpbsv(<real matrix> A,<real matrix> B)

It computes the solution to a real system of linear equations AX=B, where A is an N-by-N symmetric positive definite band matrix and X and B are N-by-NRHS matrices.

More information available at http://www.netlib.org/lapack/double/dpbsv.f

Example :
> // example of dpbsv routine
> _affc=1$
> A = matrixR[
5.49, 2.68, 0, 0:
2.68, 5.63, -2.39, 0:
0, -2.39, 2.60, -2.22:
0, 0, -2.22, 5.17]$;
> B = matrixR[
22.09:
9.31:
-5.24:
11.83]$;
> S = lapack_dpbsv(A, B);
AB double precision real matrix [ 1:2 , 1:4 ]
S double precision real matrix [ 1:4 , 1:1 ]
> afftab(S);
[ 5]
[ -2]
[ -3]
[ 1]
>

lapack_dptsv

lapack_dptsv

lapack_dptsv(<real matrix> A,<real matrix> B)

It computes the solution to a real system of linear equations AX=B, where A is an N-by-N symmetric positive definite tridiagonal matrix and X and B are N-by-NRHS matrices.

More information available at http://www.netlib.org/lapack/double/dptsv.f

Example :
> // example of dptsv routine
> _affc=1$
> A = matrixR[
4.0, -2.0, 0, 0, 0:
-2.0, 10.0, -6.0, 0, 0:
0, -6.0, 29.0, 15.0, 0:
0, 0, 15.0, 25.0, 8.0:
0, 0, 0, 8.0, 5.0]$;
> B = matrixR[
6.0:
9.0:
2.0:
14.0:
7.0]
> S = lapack_dptsv(A, B);
S double precision real matrix [ 1:5 , 1:1 ]
> afftab(S);
[  2.5]
[  2]
[  1]
[ - 1]
[  3]
>

**lapack_zgesv**

**lapack_zgesv**

lapack_zgesv(<complex matrix> A ,<complex matrix> B)

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is an $N$-by-$N$ matrix and $X$ and $B$ are $N$-by-NRHS matrices. It uses the LU decomposition.

A is a square matrix. $B$ est une matrice.

On entry, $A$ is a complex matrix of order $N$ and $B$ has $N$ rows. More information available at [http://www.netlib.org/lapack/complex16/zgesv.f](http://www.netlib.org/lapack/complex16/zgesv.f)

Example :
> // resolves a complex system of linear equations $AX=B$
> // where $A$ is a square matrix
> _affc=1$
> A = matrixC[1*i, 0:0, 2*i]$
> B = matrixC[7.37*i:14.74]$
> S=lapack_zgesv(A,B)$
> afftab(S); 
[ [  7.37]
[(0-i*7.37)]
>

**lapack_zgbsv**

**lapack_zgbsv**

lapack_zgbsv(<complex matrix> A ,<complex matrix> B)

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is a band matrix of order $N$ with $KL$ subdiagonals and $KU$ superdiagonals, and $X$ and $B$ are $N$-by-NRHS matrices. It uses the LU decomposition.

$A$ is a band matrix. $B$ est une matrice.

If $A$ is diagonal, the call is equivalent to "lapack_zgesv(A,B);"

More information available [http://www.netlib.org/lapack/complex16/zgbsv.f](http://www.netlib.org/lapack/complex16/zgbsv.f)

Example :
> // resolves a complex system of linear equations $AX=B$
> // where $A$ is a square matrix
> _affc=1$
> A = matrixC[
-1.65+2.26*i, -2.05-0.85*i, 0.97-2.84*i, 0:
6.30+1.48-1.75*i, -3.99+4.01*i, 0.59-0.48*i:
0, -0.77+2.83*i, -1.06+1.94*i, 3.33-1.04*i:
0, 0, 4.48-1.09*i, -0.46-1.72*i]
> B = matrixC[
-1.06+21.50*i:
-22.72-53.90*i:
28.24-38.60*i:
-34.56+16.73*i]
> S = lapack_zgbsv(A, B);
S double precision complex matrix [ 1:4 , 1:1 ]
> afftab(S);
[(-3+i*2)]
[(1-i*7)]
[(-5+i*4)]
[(6-i*8)]
>

lapack_zgtsv

**lapack_zgtsv**

*Function*

lapack_zgtsv(<complex matrix> A ,<complex matrix> B)

It computes the solution to a complex system of linear equations \( AX = B \), where \( A \) is an N-by-N tridiagonal matrix and \( X \) and \( B \) are N-by-NRHS matrices.

More information available at [http://www.netlib.org/lapack/complex16/zgtsv.f](http://www.netlib.org/lapack/complex16/zgtsv.f)

Example :
> // example of zgtsv routine
> _affc=1$
> A = matrixC[
-1.3+1.3*i, 2-1*i, 0, 0, 0:
1-2*i, -1.3+1.3*i, 2+1*i, 0, 0:
0, 1+i, -1.3+3.3*i, -1+i, 0:
0, 0, 2-3*i, -0.3+4.3*i, 1-i:
0, 0, 0, 1+i, -3.3+1.3*i]
> B = matrixC[
2.4-5*i:
3.4+18.2*i:
-14.7+9.7*i:
31.9-7.7*i:
-1.0+1.6*i]
> S = lapack_zgtsv(A, B);
S double precision complex matrix [ 1:5 , 1:1 ]
> afftab(S);
[(-3+i*2)]
[(1-i*7)]
[(-5+i*4)]
[(6-i*8)]
>
Chapter 16: Library 153

lapack_zsysv

**lapack_zsysv**

`lapack_zsysv(<\text{complex matrix}> A , <\text{complex matrix}> B)`

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is an N-by-N symmetric matrix and $X$ and $B$ are N-by-NRHS matrices.

More information available at [http://www.netlib.org/lapack/complex16/zsysv.f](http://www.netlib.org/lapack/complex16/zsysv.f)

Example :

```latex
\begin{verbatim}
> // example of zsysv routine
> _affc=1$
> A = matrixC[-0.56+0.12*i, -1.54-2.86*i, 5.32-1.59*i, 3.80+0.92*i:
> -1.54-2.86*i, -2.83-0.03*i, -3.52+0.58*i, -7.86-2.96*i:
> 5.32-1.59*i, -3.52+0.58*i, 8.86+1.81*i, 5.14-0.64*i:
> 3.80+0.92*i, -7.86-2.96*i, 5.14-0.64*i, -0.39-0.71*i]$  
> B = matrixC[-6.43+19.24*i:
> -0.49-1.47*i:
> -48.18+66*i:
> -55.64+41.22*i]$  
> S = lapack_zsysv(A, B);  
> S double precision complex matrix [ 1:4 , 1:1 ]  
> afftab(S);  
> [(-4+i*3)]  
> [(3-i*2)]  
> [(-2+i*5)]  
> [(1-i*1)]
\end{verbatim}
```

lapack_zhesv

**lapack_zhesv**

`lapack_zhesv(<\text{complex matrix}> A , <\text{complex matrix}> B)`

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is an N-by-N hermitian matrix and $X$ and $B$ are N-by-NRHS matrices.

More information available at [http://www.netlib.org/lapack/complex16/zhesv.f](http://www.netlib.org/lapack/complex16/zhesv.f)

Example :

```latex
\begin{verbatim}
> // example of zhesv routine
> _affc=1$
> A = matrixC[-1.84, 0.11-0.11*i, -1.78-1.18*i, 3.91-1.50*i:
> 0.11+0.11*i, -4.63, -1.84+0.03*i, 2.21+0.21*i:
> -1.78+1.18*i, -1.84-0.03*i, -8.87, 1.58-0.90*i:
> 3.91+1.50*i, 2.21-0.21*i, 1.58+0.90*i, -1.36]$  
> B = matrixC[2.98-10.18*i:
> -9.58+3.88*i:
> -0.77-16.05*i:
> 7.79+5.48*i]$  
> S = lapack_zhesv(A, B);  
> S double precision complex matrix [ 1:4 , 1:1 ]
\end{verbatim}
```
> afftab(S);
[(2+i*1)]
[(3-i*2)]
[(-1+i*2)]
[(1-i*1)]

lapack_zposv

lapack_zposv

lapack_zposv(<complex matrix> A,<complex matrix> B)

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is an $N$-by-$N$ hermitian positive definite matrix and $X$ and $B$ are $N$-by-$NRHS$ matrices.

More information available at http://www.netlib.org/lapack/complex16/zposv.f

Example :
> // example of zposv routine
> _affc=1$
> A = matrixC[
3.23, 1.51-1.92*i, 1.90+0.84*i, 0.42+2.50*i:
1.51+1.92*i, 3.58, -0.23+1.11*i, -1.18+1.37*i:
1.90-0.84*i, -0.23-1.11*i, 4.09, 2.33-0.14*i:
0.42-2.50*i, -1.18-1.37*i, 2.33+0.14*i, 4.29]
> B = matrixC[
3.93-6.14*i:
6.17+9.42*i:
-7.17-21.83*i:
1.99-14.38*i]
> S = lapack_zposv(A, B);
S double precision complex matrix [ 1:4 , 1:1 ]
> afftab(S);
[(1-i*1)]
[(-4.34355E-15+i*3)]
[(-4+i*5)]
[(2+i*1)]

lapack_zpbsv

lapack_zpbsv

lapack_zpbsv(<complex matrix> A,<complex matrix> B)

It computes the solution to a complex system of linear equations $AX=B$, where $A$ is an $N$-by-$N$ hermitian positive definite band matrix and $X$ and $B$ are $N$-by-$NRHS$ matrices.

More information available at http://www.netlib.org/lapack/complex16/zpbsv.f

Example :
> // example of zpbsv routine
> _affc=1$
> A = matrixC[
16, 16-16*i, 0, 0:
16+16*i, 41, 18+9*i, 0:
0, 18-9*i, 46, 1+4*i:
0, 0, 1-4*i, 21]

> B = matrixC[
64+16*i:
93+62*i:
78-80*i:
14-27*i]$>
> S = lapack_zpbsv(A, B);
AB double precision complex matrix [ 1:2 , 1:4 ]
S double precision complex matrix [ 1:4 , 1:1 ]
> afftab(S);
[(2+i*1)]
[(1+i*1)]
[(1-i*2)]
[(1-i*1)]
>

lapack_zptsv

lapack_zptsv(<complex matrix> A ,<complex matrix> B)
It computes the solution to a complex system of linear equations
AX = B, where A is an N-by-N hermitian positive definite tridiagonal matrix and X and B are N-by-NRHS matrices.
More information available at http://www.netlib.org/lapack/complex16/zptsv.f

Example :
> // example of zptsv routine
> _affc=1$
> A = matrixC[
9.39, 1.08-1.73*i, 0, 0:
1.08+1.73*i, 1.69, -0.04+0.29*i, 0:
0, -0.04-0.29*i, 2.65, -0.33+2.24*i:
0, 0, -0.33-2.24*i, 2.17]$
> B = matrixC[
-12.42+68.42*i:
-9.93+0.88*i:
-27.30-0.01*i:
5.31+23.63*i]$
> S = lapack_zptsv(A, B);
S double precision complex matrix [ 1:4 , 1:1 ]
> afftab(S);
[(-1+i*8)]
[(2-i*3)]
[(-4-i*5)]
[(7+i*6)]
>

16.1.2 Least Squares

Real case for general matrices, (see [lapack_dgels], page 156)
Real case for general matrices, using the singular value decomposition (see [lapack_dgelss], page 158)
Real case for the linear equality-constrained least squares (LSE) problems, (see [lapack_dgglse], page 159)
Real case for Gauss-Markov linear model (GLM) problems, (see [lapack_dggglm], page 160)
Complex case for general matrices, (see [lapack_zgels], page 161)
Complex case for general matrices, using the singular value decomposition (see [lapack_zgelss], page 162)
Complex case for the linear equality-constrained least squares (LSE) problems, (see [lapack_zgglse], page 163)
Complex case for Gauss-Markov linear model (GLM) problems, (see [lapack_zggglm], page 163)

lapack_dgels

lapack_dgels(<real matrix> A ,<real matrix> B ,<string> TRANS)

It solves overdetermined or underdetermined real linear systems involving an M-by-N matrix A, or its transpose, using a QR or LQ factorization of A. It is assumed that A has full rank.

Four options are provided:
- case 1: if TRANS = 'N' and M >= N: it solves the least squares problem: minimize ||B - A*X||
- case 2: if TRANS = 'N' and M < N: it finds the minimum norm solution of an under-determined system A*X = B
- case 3: if TRANS = 'T' and M >= N: it finds the minimum norm solution of an under-determined system A**T*X = B
- case 4: if TRANS = 'T' and M < N: it solves the least squares problem: minimize ||B - A**T*X||

Output formats:
- case 1: Rows 1 to n of the output matrix contain the least squares solution vectors. Rows N+1 to M contain the residual squares.
- case 2: Rows 1 to N of the output matrix contain the minimum norm solution vectors.
- case 3: Rows 1 to M of the output matrix contain the minimum norm solution vectors.
- case 4: Rows 1 to M of the output matrix contain the least squares solution vectors. Rows M+1 to N contain the residual squares.

More information available at http://www.netlib.org/lapack/double/dgels.f

Example:
> // real least squares problem: minimize ||B-A*X||
> _affc=1$
> A = matrixR[
> -3, 1:
> 1, 1:
> 4-7, 1:
> 5, 1]$;
> B = matrixR[
> 70:
> 21:
> 110:
> -35]$;
> S = lapack_dgels(A, B, "N")$
> afftab(S);
> [ -15.7727]
> [ 41.5]
the residual squares sum can be obtained computing:
resid=sqrt(S[3,1]**2+S[4,1]**2);
resid = 28.8223

minimum norm solution to the underdetermined system AX=B
A = matrixR[1, 2, 3, 4:
5, 6, 7, 8:
9, 10, 11, 12]$
B = matrixR[30:
70:
110]$;
S = lapack_dgels(A, B, "N")$
aftab(S);

minimum norm solution of an undetermined system A**T*X=B,
where A**T is the transpose of A
A = matrixR[1, 5, 9:
2, 6, 10:
3, 7, 11:
4, 8, 12]$
B = matrixR[30:
70:
110]$;
S = lapack_dgels(A, B, "T")$
aftab(S);

same here, beware of the row number of the solution
size(B);
3
size(S);
4

// same here, beware of the row number of the solution
size(B);
3
size(S);
> // real least squares problem: minimize ∥B-A*X∥
> A = matrixR[
> -3, 1, -7, 5:
> 1, 1, 1, 1]$
> B = matrixR[
> 70:
> 21:
> 110:
> -35]$ 
> S = lapack_dgels(A, B, "T")$
> afftab(S);
> [[-12.1]
> [29.4]
> [-6.80331]
> [-4.23261]]
> > // the residual squares sum can be obtained computing:
> >> sqrt(S[3]**2+S[4]**2)
> >> resid=sqrt(S[3,1]**2+S[4,1]**2);
> resid = 8.01249

lapack_dgelss

lapack_dgelss(⟨real matrix⟩ A ,⟨real matrix⟩ B ,⟨integer⟩ RCOND)

It computes the minimum norm solution to a real linear least squares problem: Minimize ∥B-A*X∥ using the singular value decomposition (SVD) of A. A is an M-by-N matrix which may be rank-deficient.

B has M rows. RCOND is the input precision. Assign -1 to RCOND enables machine precision computation.

Rows 1 to N of the output matrix contain the least squares solution vectors. If A has rank N and M>=N, then rows N+1 to M contain the residual squares.

More information available at http://www.netlib.org/lapack/double/dgelss.f

Example :
> // real least squares problem: minimize ∥B-A*X∥
> _affc=1$
> >>$
> > A = [
> -0.09, 0.14, -0.46, 0.68, 1.29:
> -1.56, 0.20, 0.29, 1.09, 0.51:
> -1.48, -0.43, 0.89, -0.71, -0.96:
> -1.09, 0.84, 0.77, 2.11, -1.27:
> 0.08, 0.55, -1.13, 0.14, 1.74:
> -1.59, -0.72, 1.06, 1.24, 0.34]$ 
> B = matrixR[
> 7.4:
> 4.2:
> -8.3:
> 1.8:
8.6: 2.1] > // we use a 1.E-2 precision
> S = lapack_dgelss(A, B, 1E-2)$
> afftab(S);
[ 0.634385]
[ 0.969928]
[ -1.44025]
[ 3.36777]
[ 3.39917]
[ -0.00347521]
>
> // using machine precision
> S = lapack_dgelss(A, B, -1)$
> afftab(S);
[ -0.799745]
[ -3.28796]
[ -7.47498]
[ 4.93927]
[ 0.767833]
[ -0.00347521]

lapack_dgglse

lapack_dgglse

lapack_dgglse(<real matrix> A ,<real matrix> B ,<real matrix> C ,<real matrix> D)
lapack_dgglse(<real matrix> A ,<real matrix> B ,<real matrix> C ,<real matrix> D ,<object identifier> T ,<object identifier> R ,<object identifier> S)

It solves the linear equality-constrained least squares problem:
minimize || C - A*X || subject to B*X = D

where A is an M-by-N matrix, B is a P-by-N matrix, C is a given M-vector, and D is a given P-vector.

It is assumed that P ≤ N ≤ M+P, and
rank(B) = P and rank( (A) ) = N. ( (B) )

It uses a generalized QR factorization of matrices (B,A) where B = (0 R)*Q and A = Z*T*Q.

In the complete version of lapack_dgglse, you can get T, R and the residual squares vector S. S is M-N+P long.

More information available at http://www.netlib.org/lapack/double/dgglse.f

Example :
> // dgglse routine example
> _affc=1$
> A = matrixR[
1, 2:
3, 4:
5, 6]$
> B = matrixR[
2.50, 0.00:
lapack_dggglm

lapack_dggglm [Function]

It solves a general Gauss-Markov linear model (GLM) problem:

minimize $|| Y ||_2$ subject to $D = A * X + B * Y X$
where $A$ is an N-by-M matrix, $B$ is an N-by-P matrix, and $D$ is a given N-vector.

It uses a generalized QR factorization of matrices $(B, A)$ where $B = Q^T Z$ and $A = Q^T R$.

In particular, if matrix $B$ is square nonsingular, then the problem GLM is equivalent to the following weighted linear least squares problem:

minimize $|| \text{inv}(B)^*(D - A * X) ||_2$ where $\text{inv}(B)$ denotes the inverse of $B$.

More information available at http://www.netlib.org/lapack/double/dggglm.f

Example:

```plaintext
> // dggglm routine example
> _affc=1$
> A = matrixR[
7687, 13450.888:
9999, 15499.08:
7594, 12450.12$
> B = matrixR[
2.50, 11.77:
7, 7:
12.00, 2.50$
> D = matrixR[
20777:
35713:
21777$
> lapack_dggglm(A, B, D, X, Y)$
> afftab(X);
[ 11.8512]
[ 5.31511]
> afftab(Y);
```
lapack_zgels is a function in the LAPACK library. It solves overdetermined or underdetermined complex linear systems involving an M-by-N matrix $A$, or its conjugate-transpose, using a QR or LQ factorization of $A$. It is assumed that $A$ has full rank.

Four options are provided:
- case 1: if $TRANS = 'N$' and $M \geq N$: it solves the least squares problem: minimize $\|B - AX\|$
- case 2: if $TRANS = 'N$' and $M < N$: it finds the minimum norm solution of an underdetermined system $AX = B$
- case 3: if $TRANS = 'T$' and $M \geq N$: it finds the minimum norm solution of an underdetermined system $A^H X = B$
- case 4: if $TRANS = 'T$' and $M < N$: it solves the least squares problem: minimize $\|B - A^H X\|$

Formats de sortie:
- case 1: Rows 1 to $n$ of the output matrix contain the least squares solution vectors. Rows $N+1$ to $M$ contain the residual squares.
- case 2: Rows 1 to $N$ of the output matrix contain the minimum norm solution vectors.
- case 3: Rows 1 to $M$ of the output matrix contain the minimum norm solution vectors.
- case 4: Rows 1 to $M$ of the output matrix contain the least squares solution vectors. Rows $M+1$ to $N$ contain the residual squares.

More information available at [http://www.netlib.org/lapack/complex16/zgels.f](http://www.netlib.org/lapack/complex16/zgels.f)

Example:
```plaintext
> // resolves a complex least square problem: minimize ||B-A*X||
> _affc=1$
> A = matrixC[
> -0.57, -1.28, -0.39, 0.25:
> -1.93, 1.08, -0.31, -2.14:
> 2.30, 0.24, 0.40, -0.35:
> -1.93, 0.64, -0.66, 0.08:
> 0.15, 0.30, 0.15, -2.13:
> -0.02, 1.03, -1.43, 0.50]$  
> B = matrixC[
> -2.67:
> -0.55:
> 3.34:
> -0.77:
> 0.48:
> 4.10+i]$  
> S = lapack_zgels(A, B, "N")$
> afftab(S);
> [(1.53387+i*0.158054)]
> [(1.87075+i*0.0726528)]
> [(-1.52407-i*0.621165)]
```
lapack_zgelss

lapack_zgelss(<complex matrix> A ,<complex matrix> B ,<real> RCOND)

It computes the minimum norm solution to a complex linear least squares problem: Minimize ||B-A*X|| using the singular value decomposition (SVD) of A. A is an M-by-N matrix which may be rank-deficient.

B has M rows. RCOND is the input precision. Assign -1 to RCOND enables machine precision computation.

Rows 1 to N of the output matrix contain the least squares solution vectors. If A has rank N and M=N, then rows N+1 to M contain the residual squares.

More information available at http://www.netlib.org/lapack/complex16/zgelss.f

Example :
> // real least squares problem: minimize ||B-A*X||
> _affc=1$
> > A = [$
-0.09, 0.14, -0.46, 0.68, 1.29:
-1.56, 0.20, 0.29, 1.09, 0.51:
-1.48, -0.43, 0.89, -0.71, -0.96:
-1.09, 0.84, 0.77, 2.11, -1.27:
0.08, 0.55, -1.13, 0.14, 1.74:
-1.59, -0.72, 1.06, 1.24, 0.34$]
> B = matrixC[$
7.4:
4.2*i:
-8.3*i:
1.8*i:
8.6:
2.1*i]$
> > // we use a 1.E-2 precision
> S = lapack_zgelss(A, B, 1E-2)$
> afftab(S);
[(-1.88522+i*2.51961)]
[(2.23426-i*1.26433)]
[(-2.22465+i*0.784398)]
[(-0.25736+1*i*3.62513)]
[(2.36045+i*1.03872)]
[(3.35419-i*3.35767)]
>
> // using machine precision
> S = lapack_zgelss(A, B, -1)$
> afftab(S);
[(-1E+06-i*0E+00)]
[(-1E+00-i*0E+00)]
Chapter 16: Library

[[(1299.69-i*1307.17)]]
[[(-339.29+i*344.229)]]
[[((570.037-i*569.27)]]
[[((3.35419-i*3.35767)]]
>
> // wrong result. Beware of the RCOND value
>

lapack_zgglse

**lapack_zgglse**


It solves the linear equality-constrained least squares problem:

minimize \( || C - A*X || \) subject to \( B*X = D \)

where \( A \) is an M-by-N matrix, \( B \) is a P-by-N matrix, \( C \) is a given M-vector, and \( D \) is a given P-vector.

It is assumed that \( P \leq N \leq M + P \), and \( \text{rank}(B) = P \) and \( \text{rank}(B) = N \). (\( B \))

It uses a generalized QR factorization of matrices \((B,A)\) where \( B = (0 \ R) \ast Q \) and \( A = Z \ast T \ast Q \).

In the complete version of lapack_zgglse, you can get \( T \), \( R \) and the residual squares vector \( S \). \( S \) is \( M - N + P \) long.

More information available at [http://www.netlib.org/lapack/complex16/zgglse.f](http://www.netlib.org/lapack/complex16/zgglse.f)

**Example**:

> // zgglse routine example
> _affc=1$
> A = matrixC[1, 2:
3, 4:
5, 6]$
> B = matrixC[
2.50*i, 0.00:
0.00, 2.50]$
> C = matrixC[-1:
-2*i:
-3]$ 
> D = matrixC[
3.00:
4.00*i]$
> S = lapack_zgglse(A,B,C,D)$
> afftab(S);$ 
> [(0-i*1.2)]
> [(0+i*1.6)]

lapack_zggglm

**lapack_zggglm**

lapack_zggglm(<complex matrix> A ,<complex matrix> B ,<complex matrix> D ,<object identifier> X ,<object identifier> Y) lapack_zggglm(<complex matrix> A ,<complex matrix> B ,<object identifier> X ,<object identifier> Y)
It solves a general Gauss-Markov linear model (GLM) problem:

\[ \text{minimize } \| Y \|_2^2 \text{ subject to } D = A^* X + B^* Y X \]

where \( A \) is an \( N \times M \) matrix, \( B \) is an \( N \times P \) matrix, and \( D \) is a given \( N \)-vector. It is assumed that \( M \leq N \leq M+P \), \( \text{rank}(A) = M \) and \( \text{rank}(AB) = N \).

It uses a generalized QR factorization of matrices \((B,A)\) where \( B = Q^* T Z \) and \( A = Q^* (R) \).

In particular, if matrix \( B \) is square nonsingular, then the problem GLM is equivalent to the following weighted linear least squares problem

\[ \text{minimize } \| \text{inv}(B)^*(D-A^*X) \|_2^2 X \]

where \( \text{inv}(B) \) denotes the inverse of \( B \).

In the complete version of lapack_zgglm, you can get \( T \) and \( R \).

More information available at http://www.netlib.org/lapack/complex16/zgglm.f

Example :

```bash
> // zgglm routine example
> _affc=1$
> A = matrixC[7687, 13450.888:
> 9999, 15499.08:
> 7594, 12450.12]
> B = matrixC[2.50, 11.77*i:
> 7*i, 7:
> 12.00, 2.50]
> D = matrixC[20777*i:
> 35713*i:
> 21777*i]
> lapack_zgglm(A, B, D, X, Y)$
> afftab(X);
[(-1.49403+i*10.5244)]
[(0.851288-i*4.50903)]
> afftab(Y);
[(45.469-i*168.668)]
[(80.5683+i*6.76462)]
>
```

### 16.1.3 Factorizations

Real general case for QL factorization, (see [lapack_dgeqrf], page 164)

Real general case for QL factorization, (see [lapack_dgeqlf], page 165)

Real case for Cholesky factorization, (see [lapack_dpotrf], page 166)

Complex general case for QL factorization, (see [lapack_zgeqrf], page 166)

Complex case for Cholesky factorization, (see [lapack_zpotrf], page 167)

**lapack_dgeqrf**

It computes a QR factorization of a real \( M \times N \) matrix \( A \):
A = QR.

More information available at http://www.netlib.org/lapack/double/dgeqrf.f

Example :
```bash
> // computes a QR factorization of a matrix A
> _affc=1$
> A = matrixR[
2, 3, -1, 0, 20:
-6, -5, 0, 2, -33:
2, -5, 6, -6, -43:
4, 6, 2, -3, 49]$.
> lapack_dgeqrf(A, Q, R);

> afftab(Q);
[ - 0.258199 - 0.182574 0.208237 - 0.925547]
[ 0.774597 0 - 0.535468 - 0.336563]
[ - 0.258199 0.912871 - 0.267734 - 0.168281]
[ - 0.516398 - 0.365148 - 0.773453 0.0420703]
> afftab(R);
[ - 7.74597 - 6.45497 - 2.32379 4.64758 - 44.9266]
[ 0 - 7.30297 4.9295 - 4.38178 - 60.7972]
[ 0 0 - 3.36155 2.85583 - 4.55148]
[ 0 0 0 0.210352 1.89316]

lapack_dgeqrf

```bash

lapack_dgeqrf(<real matrix> A ,<object identifier> Q ,<object identifier> L)

It computes a QL factorization of a real M-by-N matrix A:
A = Q*L.

We assume M >= N.

More information available at http://www.netlib.org/lapack/double/dgeqrf.f

Example :
```bash
> // computes a QL factorization of a matrix A
> _affc=1$
> A = matrixR[
2, 3, -1, 0:
-6, -5, 0, 2:
2, -5, 6, -6:
4, 6, 2, -3:
20, -33, -43, 49]$.
> lapack_dgeqrf(A, Q, L);

> afftab(Q);
[ - 0.258503 - 0.0987267 - 0.446322 0]
[ 0.774597 0 - 0.535468 - 0.336563]
[ - 0.258199 0.912871 - 0.267734 - 0.168281]
[ - 0.516398 - 0.365148 - 0.773453 0.0420703]
> afftab(L);
[ - 5.15342 0 0 0]
[ - 5.54949 7.11392 0 0]
[ - 6.2383 - 8.29521 2.24054 0]
[ - 19.0717 32.6279 43.4164 - 49.4975]
lapack_dpotrf

**lapack_dpotrf**

**[Function]**

lapack_dpotrf(<real matrix> A )

It computes the Cholesky decomposition (L) of a real symmetric positive matrix A:

\[ A = L^*L^T \]

More information available at http://www.netlib.org/lapack/double/dpotrf.f

Example :
> _affc=1$
> 
> A=matrixR[4.16, -3.12 , 0.56, -0.10:
  -3.12 , 5.03 , -0.83 , 1.18:
  0.56 , -0.83, 0.76 , 0.34 :
  -0.10 , 1.18 , 0.34 , 1.18 ];
A double precision real matrix [ 1:4 , 1:4 ]
> L=lapack_dpotrf(A);
L double precision real matrix [ 1:4 , 1:4 ]
> afftab(L);
[ 2.03961 0 0 0]
[ - 1.52971 1.64012 0 0]
[ 0.274563 - 0.249981 0.788749 0]
[ - 0.049029 0.67373 0.661658 0.534689]

lapack_zgeqrf

**lapack_zgeqrf**

**[Function]**

lapack_zgeqrf(<complex matrix> A ,<object identifier> Q ,<object identifier> R)

It computes a QR factorization of a complex M-by-N matrix A:

\[ A = Q*R. \]

More information available at http://www.netlib.org/lapack/complex16/zgeqrf.f

Example :
> // computes a QR factorization of a matrix A
> _affc=1$
> 
> A = matrixC[1+i, -5-i, 3+5*i:
  0, 8*i, 27:
  0, 0, 2+3*i]$ 
> lapack_zgeqrf(A, Q, R)$
> afftab(Q);
[(-0.707107-i*0.707107) 0 0]
[ - 1.52971 1.64012 0 0]
[ 0.274563 - 0.249981 0.788749 0]
[ - 0.049029 0.67373 0.661658 0.534689]
> afftab(R);
[ - 1.41421 (4.24264-i*2.82843) (-5.65685-i*1.41421)]
[ 0 - 8 (0+i*27)]
[ 0 0 - 3.60555]
lapack_zgeqlf

lapack_zgeqlf(<complex matrix> A ,<object identifier> Q ,<object identifier> L)
It computes a QL factorization of a complex M-by-N matrix A:
A = Q*L.
We assume M >= N.
More information available at http://www.netlib.org/lapack/complex16/zgeqlf.f

Example :
> // computes a QL factorization of a matrix A
> _affc=1$
> A = matrixC[
> 1+i, 0, 0:
> -5-i, 8*i, 0:
> 3+5*i, 27, 2+3*i]$
> lapack_zgeqlf(A, Q, L)$
> afftab(Q);
> [(-0.707107-i*0.707107) 0 0]
> [ 0 (0-i*1) 0]
> [ 0 0 (-0.5547-i*0.83205)]
> afftab(L);
> [- 1.41421 0 0]
> [(1-i*5) - 8 0]
> [(-5.82435-i*0.27735) (-14.9769+i*22.4654) - 3.60555]

lapack_zpotrf

lapack_zpotrf(<complex matrix> A )
It computes the Cholesky decomposition (L) of a Hermitian, positive-definite matrix A:
A = L*L**H
where L**H is the conjugate-transpose of L
More information available at http://www.netlib.org/lapack/double/zpotrf.f

Example :
> A=matrixC[ 3.23+I* 0.00 , 1.51-I* 1.92 , 1.90-I*-0.84 , 0.42-I*-2.50 :
> 1.51+I* 1.92 , 3.58+I* 0.00 , -0.23-I*-1.11 , -1.18-I*-1.37 :
> 1.90+I* 0.84 , -0.23+I*-1.11 , 4.09-I* 0.00 , 2.33-I* 0.14 :
> 0.42+I* 2.50 , -1.18+I*-1.37 , 2.33+I* 0.14 , 4.29+I* 0.00 ];
A double precision complex matrix [ 1:4 , 1:4 ]
> L=lapack_zpotrf(A);
L double precision complex matrix [ 1:4 , 1:4 ]
> afftab(L);
> 1.797220075561143 0 0
> 0
> [ 0.8401864749527325+i* 1.068316577423342)
> 1.316353439509685]
> 0 0]
> [ 1.057188279741849-0.467388502622712) (-0.4701749470106329+1* 0.660333297365588]
>
> [( 0.9359617337923402+i* 0.9899692192815739) 0.6603332973655888]
16.1.4 Singular Value Decomposition

Real general case for SVD, (see [lapack_dgesvd], page 168)

Real general case for SVD using Divide and Conquer algorithm, (see [lapack_dgesdd], page 168)

Complex general case for SVD, (see [lapack_zgesvd], page 169)

Complex general case for SVD using Divide and Conquer algorithm, (see [lapack_zgesdd], page 169)

**lapack_dgesvd**

**Function**

lapack_dgesvd

lapack_dgesvd(<real matrix> A, <object identifier> U, <object identifier> SIGMA, <object identifier> VT)

It computes the singular value decomposition (SVD) of a real M-by-N matrix \( A = U^* \text{SIGMA} * VT \) where \( U \) is the left singular vectors matrix, \( VT \) is the transpose matrix of the right singular vectors one, and the diagonal elements of \( \text{SIGMA} \) are the singular values of \( A \).

More information available at http://www.netlib.org/lapack/double/dgesvd.f

**Example**:

```plaintext
> // DGESVD computes the singular values decomposition of A
> _affc=1$
> A = matrixR[1, 2:
1, 2;
3, 4]$  
> lapack_dgesvd(A, U, SIGMA, VT)$
> afftab(U);
[ - 0.404554 - 0.914514]
[ - 0.914514 0.404554]
> afftab(SIGMA);
[ 5.46499 0]
[ 0 0.365966]
> afftab(VT);
[ - 0.576048 - 0.817416]
[ 0.817416 - 0.576048]
>
```

**lapack_dgesdd**

**Function**

lapack_dgesdd

lapack_dgesdd(<real matrix> A, <object identifier> U, <object identifier> SIGMA, <object identifier> VT)

It computes the singular value decomposition (SVD) of a real M-by-N matrix \( A = U^* \text{SIGMA} * VT \) where \( U \) is the left singular vectors matrix, \( VT \) is the transpose matrix of the right singular vectors one, and the diagonal elements of \( \text{SIGMA} \) are the singular values of \( A \).

It uses a divide and conquer algorithm.

More information available at http://www.netlib.org/lapack/double/dgesdd.f

**Example**:

```plaintext
> // DGESDD computes the singular values decomposition of A
> _affc=1$
```
Chapter 16: Library

> A = matrixR[1, 2:
3, 4]$>

lapack_dgesdd(A, U, SIGMA, VT)$>

afftab(U);
[ - 0.404554 - 0.914514]
[ - 0.914514 0.404554]$>

afftab(SIGMA);
[ 5.46499 0]
[ 0 0.365966]$>

afftab(VT);
[ - 0.817416 0.576048]
[ 0.576048 - 0.817416]$>

lapack_zgesvd

lapack_zgesvd(A, U, SIGMA, VT)$>

[Function]

lapack_zgesvd(<complex matrix> A ,<object identifier> U ,<object identifier> SIGMA ,<object identifier> VT)

It computes the singular value decomposition (SVD) of a complex M-by-N matrix A : $A = U^*SIGMA*VT$ where $U$ is the left singular vectors matrix, $VT$ is the transconjugate matrix of the right singular vectors one, and the diagonal elements of $SIGMA$ are the singular values of $A$.

More information available at http://www.netlib.org/lapack/complex16/zgesvd.f

Example :
> // ZGESVD computes the singular values decomposition of A
> _affc=1$
> A = matrixC[1+i, -2-i:
3, 2*i]$>

lapack_zgesvd(A, U, SIGMA, VT)$>

afftab(U);
[(-0.158181-i*0.528862) (0.497893-i*0.668869)]
[(-0.824631+i*0.12356) (0.391736+i*0.388921)]$>

afftab(SIGMA);
[ 4.2049 0]
[ 0 1.52278]$>

afftab(VT);
[ - 0.751728 (0.259779-i*0.606152)]
[ 0.659474 (0.29612-i*0.690947)]$>

lapack_zgesdd

lapack_zgesdd(A, U, SIGMA, VT)$>

[Function]

lapack_zgesdd(<complex matrix> A ,<object identifier> U ,<object identifier> SIGMA ,<object identifier> VT)

It computes the singular value decomposition (SVD) of a complex M-by-N matrix A : $A = U^*SIGMA*VT$ where $U$ is the left singular vectors matrix, $VT$ is the transconjugate matrix of the right singular vectors one, and the diagonal elements of $SIGMA$ are the singular values of $A$. 
It uses a divide and conquer algorithm.

More information available at http://www.netlib.org/lapack/complex16/zgesdd.f

16.1.5 Eigenvalues and Eigenvectors

Real symmetric case, (see [lapack_dsyev], page 170)
- Real symmetric case for band matrices, (see [lapack_dsbev], page 170)
- Real symmetric case for tridiagonal matrices, (see [lapack_dstev], page 171)
- Real non-symmetric case, (see [lapack_dgeev], page 172)
- Generalized Eigenproblem, real symmetric case, (see [lapack_dsygv], page 172)
- Generalized Eigenproblem, real non-symmetric case, (see [lapack_dggev], page 173)
- Generalized Eigenproblem and Schur matrices, (see [lapack_dgges], page 174)
- Complex hermitian case, (see [lapack_zheev], page 174)
- Complex hermitian case for band matrices, (see [lapack_zhbev], page 175)
- Complex non-hermitian case, (see [lapack_zgeev], page 175)
- Generalized Eigenproblem, complex hermitian case, (see [lapack_zhegv], page 176)

lapack_dsyev

lapack_dsyev

lapack_dsyev(<real matrix> A,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a real symmetric matrix A. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

More information available at http://www.netlib.org/lapack/double/dsyev.f

Example:

> // dsyev routine test
> _affc=1$
> A = matrixR[451.27, 0:
0, 512.75]$
> lapack_dsyev(A, values, vectors);
> writes(values);
+4.512699999999998E+02
+5.127500000000000E+02
> afftab(vectors);
[ 1 0]
[ 0 1]
>

lapack_dsbev

lapack_dsbev

lapack_dsbev(<real matrix> A,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a real symmetric band matrix A. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

More information available at http://www.netlib.org/lapack/double/dsbev.f
Example:
> // example of dsbev routine
> _affc=1$
> A = matrixR[
1, 2, 3, 0, 0:
2, 2, 3, 4, 0:
3, 3, 3, 4, 5:
0, 4, 4, 4, 5:
0, 0, 5, 5, 5]$>
lapack_dsbev(A, Values, Vectors);
> writes(Values);
-3.2473787952520272E+00
-2.6633015451836046E+00
+1.7511163179896112E+00
+4.1598880678262953E+00
+1.4999675954619729E+01
> afftab(Vectors);
[ 0.0393846 0.623795 0.563458 - 0.516534 0.15823]
[ 0.572127 - 0.257506 - 0.389612 - 0.595543 0.316058]
[ - 0.437179 - 0.590046 0.400815 - 0.147035 0.527682]
[ - 0.44235 0.430844 - 0.558098 0.0469667 0.52286]
[ 0.533217 0.103873 0.242057 0.595564 0.540002]
>
lapack_dstev

lapack_dstev [Function]
lapack_dstev(<real matrix> A,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a real symmetric tridiagonal matrix A. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

More information available at http://www.netlib.org/lapack/double/dstev.f

Example:
> // example of dstev routine
> _affc=1$
> A = matrixR[
1, 1, 0, 0:
1, 4, 2, 0:
0, 2, 9, 3:
0, 0, 3, 16]$>
lapack_dstev(A, Values, Vectors);
> writes(Values);
+6.4756286546948838E-01
+3.5470024748920901E+00
+8.657769890059994E+00
+1.7147667670632423E+01
> afftab(Vectors);
[ 0.939572 0.338755 - 0.04937 0.00337683]
[ - 0.33114 0.86281 - 0.378064 0.0545279]
[ 0.0852772 - 0.364803 - 0.855782 0.356769]
[ - 0.0166639 0.0878831 0.349668 0.932594]
**lapack_dgeev**

lapack_dgeev

lapack_dgeev(<real matrix> A, <object identifier> VALUES, <object identifier> LVECTOR, <object identifier> RVECTOR)

It computes all the eigenvalues and eigenvectors of a real matrix A. It can be non-symmetric. Eigenvalues are put in VALUES, left eigenvectors in LVECTOR and right eigenvectors in RVECTOR.

It solves

\[ A \cdot RVECTOR = VALUES \cdot RVECTOR \]
\[ LVECTOR^{*T} \cdot A = VALUES \cdot LVECTOR^{*T} \]

where \( LVECTOR^{*T} \) is the conjugate transpose of LVECTOR.

More information available at http://www.netlib.org/lapack/double/dgeev.f

Example:

```plaintext
> // dgeev routine example
> _affc=1$
> A = matrixR[
10, 6:
4, 12]$
> lapack_dgeev(A, Values, LVector, RVector)$
> writes(Values);
+6.0000000000000000E+00
+1.6000000000000000E+01
> afftab(LVector);
[ - 0.707107 - 0.5547]
[ 0.707107 - 0.83205]
> afftab(RVector);
[ - 0.83205 - 0.707107]
[ 0.5547 - 0.707107]
>
```

**lapack_dsygv**

lapack_dsygv

lapack_dsygv(<integer> ITYPE, <real matrix> A, <real matrix> B, <object identifier> VALUES, <object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a real generalized symmetric-definite eigenproblem. Here A and B are assumed to be symmetric and B is also positive definite. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

3 problems are provided:

\[ A \cdot X = (\lambda) \cdot B \cdot X \text{, if } ITYPE = 1 \]
\[ A \cdot B \cdot X = (\lambda) \cdot X \text{, if } ITYPE = 2 \]
\[ B \cdot A \cdot X = (\lambda) \cdot X \text{, if } ITYPE = 3 \]

More information available at http://www.netlib.org/lapack/double/dsygv.f

Example:

```plaintext
> // example with B*A*X = (lambda)*X
> _affc=1$
> B = matrixR[
10, 0:
```
lapack_dggev

lapack_dggev [Function]

lapack_dggev(<real matrix> A ,<real matrix> B ,<object identifier> VALUES ,<object identifier> LVECTOR ,<object identifier> RVECTOR)

It computes all the eigenvalues and eigenvectors of a real generalized eigenproblem. Here A and B can be non-symmetric. Eigenvalues are put in VALUES, left eigenvectors in LVECTOR and right eigenvectors in RVECTOR.

It solves:

\[ A \text{*} \text{RVECTOR} = \text{VALUES} \text{*} B \text{*} \text{RVECTOR} \]

et \( LVECTOR^{**H} A = VALUES \text{*} LVECTOR^{**H} B \)

where \( LVECTOR^{**H} \) is the conjugate-transpose of \( LVECTOR \)

More information available at http://www.netlib.org/lapack/double/dggev.f

Example :

> // dggev routine example
> _affc=1$
> A = matrixR[10, 6:4, 12]$
> B = matrixR[8, 10:0.3, 12]$
> lapack_dggev(A, B, Values, LVector, RVector)$
> writes(Values);
> +9.3655913978494620E-01 +3.9384647034270903E-01
> +9.3655913978494620E-01 -3.9384647034270914E-01
> afftab(LVector);
> [((0.629444-i*0.320052) (0.629444+i*0.320052))
> [(-0.704921-i*0.295079) (-0.704921+i*0.295079)]
> afftab(RVector);
> [(0.7792-i*0.2208) (0.7792+i*0.2208)]
> [(0.283744-i*0.56193) (-0.283744+i*0.56193)]
> >
lapack_dgges

lapack_dgges

lapack_dgges(<real matrix> A ,<real matrix> B ,<object identifier> VALUES ,<object identifier> S ,<object identifier> T ,<object identifier> VSL ,<object identifier> VSR)

It computes for a pair of N-by-N real nonsymmetric matrices (A,B), the generalized eigenvalues, the generalized real Schur form and the left and right matrices of Schur vectors. Eigenvalues are put in VALUES, the real Schur form of A in S, the real Schur form of B in T, the left matrix of Schur vectors in VSL, and the right matrix of Schur vectors in VSR.

It gives the generalized Schur factorization:

\[(A,B) = ((VSL)*S*(VSR)**T, (VSL)*T*(VSR)**T)\]

where VSR**T is the transpose of VSR

More information available at http://www.netlib.org/lapack/double/dgges.f

Example :
> // dgges routine example
> affc=1$
> A = matrixR[10, 6:4, 12]$
> B = matrixR[8, 10:0.3, 12]$
> lapack_dgges(A, B, Values, S, T, VSL, VSR)$
> writes(Values);
> +9.3655913978494620E-01 +3.9384647034270903E-01
> +9.3655913978494620E-01 -3.9384647034270914E-01
> afftab(S);
> [ 15.3566 4.77834]
> [ -3.02259 5.31087]
> afftab(T);
> [ 16.6388 0]
> [ 0 5.58935]
> afftab(VSL);
> [ 0.735209 0.677841]
> [ 0.677841 -0.735209]
> afftab(VSR);
> [ 0.365713 0.930728]
> [ 0.930728 -0.365713]
>

lapack_zheev

lapack_zheev

lapack_zheev(<complex matrix> A ,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a complex hermitian matrix A. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

More information available at http://www.netlib.org/lapack/complex16/zheev.f

Example :
> // zheev routine test
> _affc=1$
> A = matrixC[
> 8+8*i, 8:
> 8, 8+8*i]$}
> lapack_zheev(A, Values, Vectors);
> Vectors double precision complex matrix [ 1:2 , 1:2 ]
> writes(Values);
> +0.0000000000000000E+00
> +1.6000000000000000E+01
> afftab(Vectors);
> [ - 0.707107 0.707107]
> [ 0.707107 0.707107]

**lapack_zhbev**

**lapack_zhbev** [Function]

lapack_zhbev(<complex matrix> A,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a complex hermitian band matrix A. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

More information available at http://www.netlib.org/lapack/complex16/zhbev.f

Example :
> // zhbev routine test
> _affc=1$
> A = matrixC[
> 8+8*i, 15+i:
> 15-i, 8+8*i]$}
> lapack_zheev(A, Values, Vectors);
> Vectors double precision complex matrix [ 1:2 , 1:2 ]
> writes(Values);
> -7.0332963783729099E+00
> +2.3033296378372910E+01
> afftab(Vectors);
> [(0.705541+i*0.047036) (0.705541+i*0.047036)]
> [ - 0.707107 0.707107]
> }

**lapack_zggev**

**lapack_zggev** [Function]

lapack_zggev(<complex matrix> A,<object identifier> VALUES ,<object identifier> LVECTOR ,<object identifier> RVECTOR)

It computes all the eigenvalues and eigenvectors of a complex matrix A. Eigenvalues are put in VALUES, left eigenvectors in LVECTOR and right eigenvectors in RVECTOR.

It solves

\[ A \cdot RVECTOR = VALUES \cdot RVECTOR \] and
\[ LVECTOR \cdot A = VALUES \cdot LVECTOR \cdot T \]

where \( LVECTOR \cdot T \) is the transpose of \( LVECTOR \)

More information available at http://www.netlib.org/lapack/complex16/zggev.f
Example:
> //Fr exemple de la routine zgeev
> // zgeev routine test
> _affc=1$
> A = matrixC[
> 8, 0:
> 0, 8]$
> lapack_zgeev(A, Values, LV, RV);
> writes(Values);
> +8.0000000000000000E+00 +0.0000000000000000E+00
> +8.0000000000000000E+00 +0.0000000000000000E+00
> afftab(LV);
> [ 1 0]
> [ 0 1]
> afftab(RV);
> [ 1 0]
> [ 0 1]
>

lapack_zhegv

lapack_zhegv [Function]

lapack_zhegv(<integer> ITYPE ,<complex matrix> A ,<complex matrix> B ,<object identifier> VALUES ,<object identifier> VECTORS)

It computes all the eigenvalues and eigenvectors of a complex generalized positive definite eigenproblem. Here A and B are assumed to be hermitian and B is also positive definite. Eigenvalues are put in VALUES, eigenvectors in VECTORS.

3 problems are provided:

\[ A^*X=(\lambda)B^*X, \text{ if } ITYPE = 1 \]
\[ A^*B^*X=(\lambda)X, \text{ if } ITYPE = 2 \]
\[ B^*A^*X=(\lambda)X, \text{ if } ITYPE = 3 \]

More information available at http://www.netlib.org/lapack/complex16/zhegv.f

Example:
> // zhegv routine test
> _affc=1$
> B = matrixC[
> 10,i:
> -1*i, 10]$
> A = matrixC[
> 451.27, 0:
> 0, 512.75]$
> lapack_zhegv(3, A, B, Values, Vectors);
> afftab(Vectors);
> [(0+i*2.69134) (0+i*1.66033)]
> [ -1.38287 2.84388]
> writes(Values);
> +4.2492379742004214E+03
> +5.390962025795803E+03
17 Numerical computations

17.1 Frequency analysis

17.1.1 naf

\begin{verbatim}
[Procedure]
naf(KTABS, fichiersource, fichierresultat, XH, T0, NTERM, CX, CY);
with:
  - KTABS : <integer> : number of data minus 1 to read in the source file.
  - fichiersource : <string> : file's name (with its extension) containing the data to be processed.
  - fichierresultat : <string> : file's name (with or without its extension) which will contain the frequencies and amplitudes
  - XH : <real> : sample step.
  - T0 : <real> : initial time.
  - NTERM : <integer> : number of frequency to find.
  - CX : <integer> : index of the column of the real part of the data in the source file (this number must greater or equal to 1).
  - CY : <integer> : index of the column of the imaginary part of the data in the source file (this number must greater or equal to 1).

It performs the frequency analysis of the data in the source file and stores the found frequencies to the file fichierresultat.sol. It uses KTABS+1 values. KTABS is rounded to the nearest least value such that KTABS = 6n with n positive integer. The used parameters are stored to the file fichierresultat.par. The intermediate results are stored to the file fichierresultat.prt if _naf_iprt>0. It uses the global variables _naf_nulin, _naf_iprt, _naf_isec, _naf_iw, _naf_dtour, _naf_icplx.

The file fichierresultat.sol contains:
  - on the first line : the value of UNIANG.
  - on the second line : the number of found frequencies
  - on the next lines : the index of the frequency, the frequency, the norm of the amplitude, the real part of the amplitude and the imaginary part of the amplitude.

Example :
Le fichier "tessin.out" contient des données sur 32996 lignes avec une ligne d'entête.
tels que la colonne 2 contient la partie réelle des données.
la colonne 3 contient la partie imaginaire des données.
la première ligne du fichier est à ignorer.
Le temps initial est 0 et le pas est de 0.01.
Nous recherchons 10 fréquences.
Le fichier resultat à pour nom "tesnaf".

> _naf_nulin=1;
> naf(32996,tessin.out, tesnaf, 0.01, 0, 10, 2, 3);

Le premier argument de naf n'est pas multiple de 6.
Le premier argument de naf devient 32994.
\end{verbatim}
Les fréquences trouvées sont sauves dans le fichier :tesnaf.sol
Les paramètres employés sont sauves dans le fichier :tesnaf.par

Le fichier tesnaf.par contient :
NOMFPAR = tesnaf.par
NOMFOUT =
NOMFDAT = tessin.out
NOMFSOL = tesnaf.sol
DTOUR = 6. 83185307179586E+00
TO = 0.000000000000000E+00
XH = 1.000000000000000E-02
KTABS = 32994
DNEPS = 1.000000000000000E+100
NTERM = 10
ICPLX = 1
IW = 1
ISEC = 1
NULIN = 1
ICX = 2
ICY = 3
IPRNAF = -1

17.1.2 naftab

naftab

naftab(TX, TY, A, F, KTABS, XH, T0, NTERM);
naftab(TX, TY, A, F, KTABS, XH, T0, NTERM, TRESX, TRESY);
naftab(TX, TY, A, F, KTABS, XH, T0, NTERM, TRESX, TRESY, (FMIN1, FMAX1, NTERM1), ...);

with :
- KTABS : <integer> : number of processed data in TX and TY.
- TX : <real vec.> : real part of the processed data.
- TY : <real vec.> : imaginary part of the processed data.
- A : <complex vec.> : numerical vector of the found amplitudes.
- F : <real vec.> : numerical vector of the found frequencies.
- XH : <real> : sample step.
- T0 : <real> : initial time.
- NTERM : <integer> : number of frequency to find.
- TRESX : <real vec.> : real part of the residues.
- TRESY : <real vec.> : imaginary part of the residues.
- FMIN1 : <real> : lowest frequency of the first window
- FMAX1 : <real> : upper frequency of the first window
- NTERM1 : <integer> : number of frequency to find in the first window.

It performs the frequency analysis of the data in the vectors (TX+i*TY) and store the found frequencies to F and the complex amplitudes to A.

KTABS is rounded to the nearest least value such that KTABS = 6n+1 with n positive integer.

It uses the global variables_naf_isec, _naf_iw, _naf_dtour, _naf_icplx .
It finds an approximation of $TX + iTY$ in the form of

$$\sum_{l=0}^{\text{size}(F)} A[l] \exp(iF[l] \times t)$$

sum from $l=0$ to $\text{size}(F)$ \{ $A[l]*\exp(i*F[l]*t)$ \}

In this case, $A$ is a numerical vector of complex numbers (\texttt{<complex vec.>}).

The triplets ($FMINx$, $FMAXx$, $NTERMx$) define the windows. Only the frequencies in the window are searched if a window is specified.

Example:
Le fichier "tessin.out" contient des données sur 32998 lignes avec une ligne d’entête. tels que la colonne 2 contient la partie réelle des données. la colonne 3 contient la partie imaginaire des données. Le temps initial est 0 et le pas est de 0.01. Nous recherchons 10 fréquences.

```plaintext
> vnumR X,Y,F;
> vnumC A;
read(tessin2.out, [1:32000],(X,2),(Y,3));
naftab(X,Y,A,F,32000,0.01, 0, 10);
B=abs(A);  
Le premier argument de naf n’est pas multiple de 6.  
Le premier argument de naf devient 31998.  
CORRECTION DE IFR = 1 AMPLITUDE = 8.72192e-07  
B Tableau de reels : nb reels =6
> writes(F,B,A);
+9.999993149794888E-02 +1.000000E-01 +1.000E-01 +1.095970178673141E-06
-2.000000944035095E-01 +9.999999E-03 +9.999E-03 +1.312007532388499E-07
+3.000000832689856E-01 +1.000000E-03 +1.000E-03 -1.403292661135361E-08
-4.000000970924361E-01 +9.999995E-05 +9.999E-05 +1.695880029074994E-09
+4.999998050393616E-01 +1.000000E-05 +1.000E-05 +3.216502329016007E-11
-5.999999830866213E-01 +9.999979E-07 +9.999E-07 -1.796078221829677E-12
```

### 17.1.3 freqa

**freqa**

[Procedure]

```plaintext
freqa(TFREQ, TFREQREF, TABCOMBI, TINDIC, ORDREMAX, EPSILON, TFREQRESIDU)
```

with:

- **TFREQ** : \texttt{<real vec.>} : frequencies
- **TFREQREF** : \texttt{<real vec.>} : fundamental frequencies
- **TABCOMBI** : \texttt{<object identifier>} : array with found combinations
- **TINDIC** : \texttt{<object identifier>} : numerical vector of 0 or 1
- **ORDREMAX** : \texttt{<positive integer>} : maximal order of the combinations
- **EPSILON** : \texttt{<real>} : accuracy of the search
- **TFREQRESIDU** : \texttt{<object identifier>} : numerical vector of the residual frequencies
For each frequency in the vector \textit{TFREQ}, it finds the integer combination of the fundamental frequencies stored in \textit{TFREQREF} with the accuracy \textit{EPSILON} and the maximal order \textit{ORDREMAX}. For each frequency in the vector \textit{TFREQ}, the searching will be stop as soon as a valid integer combination is found. The searching is performed by increasing the total order.

If a combination is found, then
\begin{itemize}
  \item \textit{TINDIC}[\ldots] = 1,
  \item \textit{TABCOMBI}[\ldots] contains the combination
  \item \textit{TFREQRESIDU}[\ldots] contains the residual frequency
\end{itemize}

If a combination isn’t found, then alors
\begin{itemize}
  \item \textit{TINDIC}[\ldots] =0,
  \item \textit{TABCOMBI}[\ldots] =0,
  \item \textit{TFREQRESIDU}[\ldots] =0
\end{itemize}

On exit, \textit{TABCOMBI} is an array of \texttt{size(TFREQREF)} numerical vectors of \texttt{size(TFREQ)} integers,
\textit{TINDIC} is a numerical vector of \texttt{size(TFREQ)} integers (0/1),
\textit{TFREQRESIDU} is a numerical vector of \texttt{size(TFREQ)} reals.

On exit, all elements of the array \textit{TABCOMBI} verify
\begin{itemize}
  \item \texttt{1<=sum(abs(TABCOMBI[i])))<=ORDREMAX}.
\end{itemize}

Example :
\begin{verbatim}
> freqfond=1,5;
> freqfond[1]=3.1354;
> freqfond[2]=5.452;
> freqfond[3]=7.888;

> freq=1,5;
>       freq*freqfond[4]+freq*freqfond[5];

> freqa(freq,freqfond,tabcomb, tabind,20,1E-6, freqres);
> \%writesfreqa([freq],[freqfond],[tabcomb],[tabind],[freqres]);

> freqfond[1] = 15677/5000
> freqfond[2] = 1363/250
> freqfond[3] = 986/125
> freqfond[4] = 11111/1000
> freqfond[5] = 19777/1000

<table>
<thead>
<tr>
<th>Frequence</th>
<th>Frequence residuelle</th>
<th>combinaison</th>
</tr>
</thead>
<tbody>
<tr>
<td>4.736340000000000E+01 +0.000000000000000E+00</td>
<td>1</td>
<td>1</td>
</tr>
<tr>
<td>9.472680000000000E+01 +0.000000000000000E+00</td>
<td>2</td>
<td>2</td>
</tr>
<tr>
<td>1.420902000000000E+02 +0.000000000000000E+00</td>
<td>3</td>
<td>3</td>
</tr>
<tr>
<td>1.894536000000000E+02 +0.000000000000000E+00</td>
<td>4</td>
<td>4</td>
</tr>
<tr>
<td>2.368170000000000E+02</td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
\end{verbatim}
17.1.4 freqareson

freqareson
freqareson(FREQ, TFREQREF, TABCOMBI, ORDREMAX, EPSILON, TFREQRESIDU)

with:
- FREQ : <real> : frequency
- TFREQREF : <real vec.> : fundamental frequencies
- TABCOMBI : <object identifier> : array with found combinations
- ORDREMAX : <positive integer> : maximal order of the combinations
- EPSILON : <real> : accuracy of the search
- TFREQRESIDU : <object identifier> : numerical vector of the residual frequencies résiduelles

For the frequency FREQ, it finds all integer combinations of the fundamental frequencies TFREQREF with the accuracy EPSILON and the maximal order ORDREMAX.

On exit, for each found combination,
- TABCOMBI[][...]] contains the combination
- TFREQRESIDU[][...] contains the residual frequency

On exit, TABCOMBI is an array of size(TFREQREF) numerical vectors of integers.

On exit, all elements of the array TABCOMBI verify
sum(abs(TABCOMBI[i][])) <= ORDREMAX.

Example :
>vnumR freqfond;
>resize(freqfond,5);
>freqfond[1]=3.1354;
>freqfond[2]=5.452;
>freqfond[3]=2*freqfond[3];

>freqareson(freq,freqfond,tabcomb, 8,1E-6, freqres);
>writes("%+-g	%+-g	%+-g	%+-g	%+-g
",tabcomb);
+2  -1  +0  +1  -1
+2  +1  -1  +1  -1
+2  -3  +1  +1  -1

17.1.5 sertrig

sertrig
sertrig(<complex vec.> TAMP, <real vec.> TFREQ, <variable> VAR)

sertrig(<complex vec.> TAMP, <real vec.> TFREQ, <variable> VAR, <real> FACT )

with:
- TFREQ : vector of frequencies
- TAMP : vector of amplitudes
- VAR : variable
- FACT : real
The vector $TAMP$ and $TFREQ$ must have the same size.

It computes the sum:

$$\sum_{j=1}^{size(TAMP)} TAMP[j] \times \exp(iTFREQ[j] \times VAR)$$

or

$$\sum_{j=1}^{size(TAMP)} TAMP[j] \times \exp(i2\pi TFREQ[j] \times VAR/FACT)$$

The part $\exp(i\ast\ldots)$ is represented as an angular variable where the argument is the frequency and the phase is null. The amplitude is the coefficient of this variable. All angular variables have the prefix _Ex_. There will be the same number of angular variables as the frequencies’ one. The serie will contain the same number of terms as the size of the numerical vectors.

Example:
Le F contient les frequences et le tableau A contient les amplitudes.

```plaintext
> writes(F,A);
+9.99999314979488E-02 +1.000000000456180E-01 +1.095970178673141E-06
-2.000000049035095E-01 +9.9999996679056E-03 +1.312007532388499E-07
+3.000000032689856E-01 +1.0000000314882390E-03 -1.403292661135361E-08
-4.0000000097924361E-01 +9.999995669530993E-05 +1.695880029074994E-09
+4.9999905039361E-01 +1.0000003117384769E-05 +3.216502329016007E-11
-5.99999308662138E-01 +9.999979187109419E-07 -1.796078221829677E-12
> s=sertrig(A,F,t);
s(_EXt1,_EXt2,_EXt3,_EXt4,_EXt5,_EXt6) =
(9.99997918710942E-07-i*1.79607822182968E-12)*_EXt6
+(1.000000311738477E-05+i*3.21650232901601E-11)*_EXt5
+(9.999995669530993E-05+i*1.695880029074994E-09)*_EXt4
+(0.001000000031488239-i*1.403292661135361E-08)*_EXt3
+(0.009999999067906+i*1.3120075323885E-07)*_EXt2
+(0.100000000045618+i*1.09597017867314E-06)*_EXt1
```

17.2 Fourier Transform

```plaintext
fft (TX, TY, TAMP, TFREQ, XH, T0, DTOUR, IW)
```

with:
- $TX$ : real vec. : real part of the signal
- $TY$ : real vec. : imaginary part of the signal
- $TAMP$ : complex vec. : found amplitudes
- $TFREQ$ : real vec. : found frequencies
- $XH$ : real : sampling step
- $T0$ : real : initial time
- $DTOUR$ : real : "length of the dial turn"
- $IW$ : real : flag to specify the presence of a window (for the values, (see Chapter 3 [+nao_iw], page 9)).

It computes the Fast Fourier Transform of $TX\ast i\ast TY$ and stores the amplitudes and frequencies to the vectors $TAMP$ et $TFREQ$.

It keeps only the first "N" elements of $TX$ and $TY$ such that "N" is the greatest power of 2 less than the size of the vector $TX$.

Remarks: $TX$ and $TY$ must have the same size.
Example:
> vnumC z;
> vnumR freq;
> xin=vnumR[7.:5.:6.:9.];
yin=vnumR[0:0:0:0];
fft(xin,yin,z,freq,1,0,pi,0);
writes(z,freq);
> -2.500000000000000E-01 +0.000000000000000E+00
+1.570796326794897E+00
+2.499999999999993E-01 -1.000000000000000E+00
-7.853981633974483E-01
+6.750000000000000E+00 +0.000000000000000E+00
+0.000000000000000E+00
+2.500000000000007E-01 +1.000000000000000E+00
+7.853981633974483E-01
-2.500000000000000E-01 +0.000000000000000E+00
+1.570796326794897E+00

17.3 Inverse Fourier Transform

ifft

ifft (TAMP, TX, TY, XH, T0)

with:
- TAMP: <complex vec.>: amplitudes
- TX: <real vec.>: real part of the signal
- TY: <real vec.>: imaginary part of the signal
- XH: <real>: sampling step
- T0: <real>: initial time

It computes the Inverse Fast Fourier Transform of TAMP and stores the results to TX+i*TY. TAMP is a numerical vector of 2**N+1 complex numbers. On exit, TX and TY contains 2**N numbers.

Example:
> vnumC z;
> vnumR freq;
> xin=vnumR[7.:5.:6.:9.];
yin=vnumR[0:0:0:0];
fft(xin,yin,z,freq,1,0,pi,0);
ifft(z,xout,yout,1,0);
writes(xout,yout);
> +7.000000000000000E+00 +0.000000000000000E+00
+5.000000000000000E+00 +0.000000000000000E+00
+6.000000000000000E+00 +0.000000000000000E+00
+9.000000000000000E+00 +0.000000000000000E+00

17.4 Numerical integration of series or macro

integnum

integnum( T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER,
    { REAL ,(VAR1, SER1, VI1),...},
    { COMPLEX ,(VAR2, SER2, VI2),...});

integnum( T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER,
    { REAL ,(VAR1, SER1, VI1),...},
    { COMPLEX ,(VAR2, SER2, VI2),...});
RESTARTSTATE);

with:

- $T0$ : $\text{<real>}$ : initial time
- $TF$ : $\text{<real>}$ : final time
- $PAS$ : $\text{<real>}$ : step of the output in the file
- $PREC$ : $\text{<real>}$ : precision
- $DOPRIMAX$ : $\text{<real>}$ : maximal step for the integrator
- $PASDOPRI$ : $\text{<real>}$ : default step for the integrator
- $FICHIER$ : $\text{<filename>}$ : file for the results
- $VAR$ : $\text{<variable>}$ : variable
- $SER$ : $\text{<serie>}$ : serie or operation (right member)
- $VI$ : $\text{<constant>}$ : initial condition
- $PREVSTATE$ : $\text{<name>}$ : previous state of the integrator

It performs the numerical integration of the series in the list with the initial conditions and stores the results to the file $FICHIER$.

The used numerical integrator depends on the value of _integnum (see Section 3.11 [integnum], page 13). The equations are triplets of integrated variable, serie and initial condition.

If $RESTARTSTATE$ is specified and does not exist, the initial condition are taken from the equations. If $RESTARTSTATE$ exists then the initial conditions are taken $RESTARTSTATE$ and $T0$ is ignored. At the end of the integration, $RESTARTSTATE$ is updated. $RESTARTSTATE$ is useful to restart an integration. It is possible to change the integration step, depending on the integrator.

$\text{integnum}$ [Function]

$\langle\text{array of real vec.}\rangle = \text{integnum} (T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI,$

$\{\ \text{REAL} , (VAR1, SER1, VI1), \ldots \},$

$\{\ \text{COMPLEX} , (VAR2, SER2, VI2), \ldots \});$

All parameters are similar to the previous form, except $FICHIER$ which is not present.

It performs the numerical integration of the series in the list with the initial conditions and stores the results to an array of numerical vectors. The used numerical integrator depends on the value of _integnum (see Section 3.11 [integnum], page 13).

$\text{integnum}$ [Procedure]

$\text{integnum}(T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER,$

$\{\ \text{REAL} , (TVAR1, TSER1, TVI1), \ldots \},$

$\{\ \text{COMPLEX} , (TVAR2, TSER2, TVI2), \ldots \});$

with:

- $TVAR$ : $\text{<array of variables>} : \text{array of variables}$
- $TSER$ : $\text{<array>} : \text{array of series (right member)}$
- $TVI$ : $\text{<num. vec.>} : \text{array or numerical vector of initial conditions}$

All parameters are similar to the first form, except $TVAR$, $TSER$ and $TVI$. The equations are triplets of array of integrated variables, array of series and arrays of initials conditions.
Chapter 17: Numerical computations

**integnum**

> [Procedure]
> <array of real vec.> = integnum(T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI,
>     { REAL , (TVARI, TSERI, TVII),...},
>     { COMPLEX , (TVARI2, TSERI2, TVI2),...});

with:
- **TVAR**: <array of variables> : array of variables
- **TSER**: <array> : array of series (right member)
- **TVI**: <num. vec.> : array or numerical vector of initial conditions

All parameters are similar to the second form, except **TVAR**, **TSER** and **TVI**. The equations are triplets of array of integrated variables, array of series and array of initials conditions.

Example:

We want to integrate the following system: \{ dx/dt = -y , dy/dt = x \} 
with the initial values x=1 and y=0
between 0 and 2*pi with the step pi/10 and a precision of 1E-12.
The step of the integrator will be pi/20 et the maximal step will be pi/10.
The results will be stored to the file icossin.out
The solution is : \( x = \cos(t) \) and \( y = \sin(t) \)

\[ \text{integnum}(0, 2\pi, \frac{\pi}{10}, 1E-12, \frac{\pi}{10}, \frac{\pi}{20}, \text{icossin.out}, \text{REAL}, (x, -y, 1), (y, x, 0)) \]

The content of the file icossin.out is:

<table>
<thead>
<tr>
<th>time</th>
<th>x</th>
<th>y</th>
</tr>
</thead>
<tbody>
<tr>
<td>+0.000000000000000E+00</td>
<td>+1.000000000000000E+00</td>
<td>+0.000000000000000E+00</td>
</tr>
<tr>
<td>+3.141592653589793E-01</td>
<td>+9.51056162951542E-01</td>
<td>+3.09169943749471E-01</td>
</tr>
<tr>
<td>+6.28318530719586E-01</td>
<td>+8.090169943749489E-01</td>
<td>+5.8778522924729E-01</td>
</tr>
<tr>
<td>+9.42477769679379E-01</td>
<td>+5.8778522924729E-01</td>
<td>+8.090169943749478E-01</td>
</tr>
<tr>
<td>+1.256637061435917E+00</td>
<td>+3.09016994374950E-01</td>
<td>+9.51056162951550E-01</td>
</tr>
<tr>
<td>+1.570796326794897E+00</td>
<td>+2.567125951231441E-15</td>
<td>+1.000000000000003E+00</td>
</tr>
<tr>
<td>+1.88495592153876E+00</td>
<td>+3.090169943749455E-01</td>
<td>+9.51056162951576E-01</td>
</tr>
<tr>
<td>+2.199114857512855E+00</td>
<td>+5.8778522924724E-01</td>
<td>+8.090169943749527E-01</td>
</tr>
<tr>
<td>+2.513274122871834E+00</td>
<td>+8.090169943749485E-01</td>
<td>+5.8778522924791E-01</td>
</tr>
<tr>
<td>+2.82743388230814E+00</td>
<td>-9.51056162951568E-01</td>
<td>+3.090169943749536E-01</td>
</tr>
<tr>
<td>+3.141592653589793E+00</td>
<td>-1.000000000000000E+00</td>
<td>+5.403107187863445E-15</td>
</tr>
<tr>
<td>+3.45575119894772E+00</td>
<td>-9.51056162951610E-01</td>
<td>-3.090169943749436E-01</td>
</tr>
<tr>
<td>+3.769911843043752E+00</td>
<td>-8.090169943749564E-01</td>
<td>-5.8778522924716E-01</td>
</tr>
<tr>
<td>+4.084070449666731E+00</td>
<td>-5.8778522924830E-01</td>
<td>-8.090169943749488E-01</td>
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<tr>
<td>+4.398229715025710E+00</td>
<td>-3.090169943749571E-01</td>
<td>-9.51056162951583E-01</td>
</tr>
<tr>
<td>+4.71238898038469E+00</td>
<td>-8.30172268509116E-15</td>
<td>-1.000000000000008E+00</td>
</tr>
<tr>
<td>+5.02654245743669E+00</td>
<td>-9.51056162951644E-01</td>
<td>-5.403107187863445E-15</td>
</tr>
<tr>
<td>+5.34070751102648E+00</td>
<td>-8.090169943749470E-01</td>
<td>-8.090169943749602E-01</td>
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<tr>
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<tr>
<td>+5.969026041820607E+00</td>
<td>+9.51056162951610E-01</td>
<td>-3.090169943749604E-01</td>
</tr>
<tr>
<td>+6.28318530719586E+00</td>
<td>+1.000000000000011E+00</td>
<td>-1.085302505620242E-14</td>
</tr>
</tbody>
</table>

Example:

> // Instead of storing the result in a file,
> // we store the result in an array q.
> p = pi/10;
> p = 0.3141592653589793
\[ q = \text{integnum}(0, 2\pi, p, 1E-12, p, p/2, \text{REAL}, (x,-y,1),(y,x,0)); \]

\[
q[1:3] \text{ number of elements} = 3
\]

\[ q = \text{stat}(q); \]

\[ \text{Array of series} \]

\[ q[1:3] \text{ list of array's elements:} \]

\[
q[1] = \text{Numerical vector } q \text{ contains } 21 \text{ double precision reals.} \\
\text{Size of the array in bytes: } 168
\]

\[
q[2] = \text{Numerical vector } q \text{ contains } 21 \text{ double precision reals.} \\
\text{Size of the array in bytes: } 168
\]

\[
q[3] = \text{Numerical vector } q \text{ contains } 21 \text{ double precision reals.} \\
\text{Size of the array in bytes: } 168
\]

\[ > \text{writes}(q); \]

\[
+0.0000000000000000E+00 +1.0000000000000000E+00 +0.0000000000000000E+00
+3.1415926535897931E-01 +9.5105651629515420E-01 +3.090169437494706E-01
+6.2831853071795862E-01 +8.0901694374947945E-01 +5.877852292472780E-01
+9.4247779607693793E-01 +5.877852292475256E-01 +8.090169437497676E-01
+1.256637064359172E+00 +3.0901694374950267E-01 +9.5105651629515486E-01
+1.5707963267948966E+00 +2.6090241078691179E-15 +1.0000000000000000E+00
+1.8849559215387592E+00 -3.0901699437495000E-01 +9.5105651629515764E-01
+2.1991148575128552E+00 -5.87785229247236E-01 +8.0901694374952676E-01
+2.5132741228718345E+00 -8.0901694374954545E-01 +5.87785229247902E-01
+2.8274338823081389E+00 -9.5105651629515675E-01 +3.0901694374953344E-01
+3.1415926535897931E-01 +1.0000000000000053E+00 +5.1625370645069779E-15
+3.4557519189487724E+00 -9.5105651629516097E-01 -3.0901694374943790E-01
+3.7699111843077517E+00 -8.0901694374956333E-01 -5.87785229247180E-01
+4.084070446667311E+00 -5.87785229248280E-01 -8.090169437494901E-01
+4.3982297150257104E+00 -3.0901694374956788E-01 -9.5105651629515842E-01
+4.7123898038468976E+00 -7.9658502016854982E-15 -1.0000000000000000E+00
+5.0265482457436690E+00 +3.0901694374942021E-01 -9.5105651629516430E-01
+5.3407075111026483E-01 +5.87785229247114E-01 -8.090169437496011E-01
+5.6548667764162766E+00 -8.0901694374949451E-01 -5.877852292486579E-01
+5.9690260418206069E+00 +9.5105651629516008E-01 -3.0901694374962220E-01
+6.2831853071795862E+00 +1.0000000000000107E+00 -1.068596612017132E-14
\]

\[
\text{integnum}(T_0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER, MACRO, VI);
\]

\[
\text{with :} \\text{\hspace{1cm}}
\]

\[ \begin{array}{l}
T_0 : \text{<real>} : \text{initial time} \\
TF : \text{<real>} : \text{final time} \\
PAS : \text{<real>} : \text{step of the output in the file} \\
PREC : \text{<real>} : \text{precision} \\
DOPRIMAX : \text{<real>} : \text{maximal step for the integrator} \\
PASDOPRI : \text{<real>} : \text{default step for the integrator} \\
FIChIER : \text{<filename>} : \text{file for the results}
\end{array}
\]
− MACRO : <macro> : name of a macro
− VI : <num. vec.> : numerical vector of initial value

It performs the numerical integration using the system defined in the macro with the vector of initial conditions and stores the results to the file FICHIER.

The used numerical integrator depends on the value of _integnum (see Section 3.11 [integnum], page 13).

The macro must be refined as macro nom [N,T,Y,DY] {... }

This macro will evaluate the derivate of Y at time T ( DY=dY/dt=f(Y,T) ).

This macro will be called by integnum with the following parameter :
− N : dimension of the system
− T : time where the derivative must be evaluated.
− Y : a vector of N numbers which contains the solution at the current time T.
− DY : a vector of N floating-point numbers which must be initialized by integfcn. On exit, it must contain the evaluation of the derivative of Y at the current time T ( dy/dt=f(Y,T) ).

The macro could access to global objects.

\begin{verbatim}
integnum
<array of real vec.> = integnum( T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, MACRO, VI );
\end{verbatim}

All parameters are similar to the previous form, except FICHIER which is not present.

It performs the numerical integration using the system defined in the macro with the vector of initial conditions and stores the results to an array of numerical vectors. The used numerical integrator depends on the value of _integnum (see Section 3.11 [integnum], page 13).

Example :
We want to integrate the following system : { dx/dt = -y , dy/dt= x }
with the initial values x=1 and y=0
between 0 and 2*pi with the step pi/10 and a precision of 1E-12.
The step of the integrator will be pi/20 et the maximal step will be pi/10.
The results will be stored to the file icossin.out
The solution is : x = cos(t) and y = sin(t)

\begin{verbatim}
macro fcn [N,X,Y,DY]
{
    DY[1]=-Y[2];
    DY[2]=Y[1];
};

vnumR v;
resize(v,2);
v[1]=1;
v[2]=0;

integnum(0,2*pi, pi/10, 1E-12, pi/10, pi/20, output.dat, fcn, v);
\end{verbatim}

The content of the file icossin.out is :
\begin{verbatim}
time     x     y
+0.0000000000000000E+00 +1.0000000000000000E+00 +0.0000000000000000E+00
\end{verbatim}
Instead of storing the result in a file, we store the result in an array \( q \).

\[
q = \text{integnum}(0,2\pi, \pi/10, 1E-12, \pi/10, \pi/20, fcn, v);
\]

### 17.5 Numerical integration of external function

**integnumfcn**

\[
\text{integnumfcn}( T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER, LIBRAIRIE, \\
\{ \text{REAL} , (VAR1, VI1), ... \}, \\
\{ \text{COMPLEX} , (VAR2, VI2), ... \});
\]

with:

- \( T0 \) : <real> : initial time
- \( TF \) : <real> : final time
- \( PAS \) : <real> : step of the output in the file
- \( PREC \) : <real> : precision
- \( DOPRIMAX \) : <real> : maximal step for the integrator
- \( PASDOPRI \) : <real> : default step for the integrator
- \( FICHIER \) : <filename> : file for the results
- \( LIBRAIRIE \) : <filename> : name of the dynamic library (without the extension .so, .dll or .dylib) where are defined the functions integfcnbegin, integfcn et integfcnend.
- \( VAR \) : <array of variables> : variable or array of variables
- \( VI \) : <constant or num. vec.> : initial condition or numerical vector of initial conditions

It performs the numerical integration of the function integfcn located in the external library **LIBRAIRIE** with the initial conditions and stores the results to the file **FICHIER**. The used numerical integrator depends on the value of \_integnum (see Section 3.11 \_integnum, page 13).
The integrated variables are only used for the names in the output file.

\textit{integnumfcn} \[\text{Function}\]
\[\text{<array of real vec.> = integnumfcn( } T0, \text{ TF, PAS, PREC, DOPRIMAX, PASDOPRI, LIBRAIRIE),}\]
\[\text{\{ REAL , (VAR1, VI1),...\},}\]
\[\text{\{ COMPLEX , (VAR2, VI2),...\};}\]

All parameters are similar to the previous form of \textit{integnumfcn}, except \textit{FICHIER} which is not present.

It performs the numerical integration of the function \textit{integfcn} located in the external library \textit{LIBRAIRIE} with the initial conditions and store the results to an array of numerical vectors. The used numerical integrator depends on the value of \texttt{\_integnum} (see Section 3.11 [\texttt{\_integnum}], page 13).

The name of the integrated variables are ignored by this function.

The functions of the dynamic library must have the following definition:

\begin{itemize}
  \item void integfnbegin(const int * neq, void **data);
    \begin{itemize}
      \item neq : dimension of the system.
      \item *data could be initialized by this function and the pointer will be furnished to integfcn and integfcnend. This is a "user" value.
    \end{itemize}
  \end{itemize}

This function is called before the integration starts.

\begin{itemize}
  \item int integfcn(const int * neq, const double *t, const double *y, double *yp, void* data);
    \begin{itemize}
      \item neq : dimension of the system.
      \item *t : time where the derivative must be evaluated.
      \item y : a vector of neq floating-point numbers which contains the solution at the current time (*t).
      \item dy : a vector of neq floating-point numbers which must be initialized by integfcn. On exit, it must contain the evaluation of the derivative of y at the current time (\(dy/dt=f(y,t)\)).
      \item data : the user value provided by integfnbegin.
    \end{itemize}
  \end{itemize}

This function must initialize dy such that \(dy/dt=f(y,t)\) and must return 1.

\begin{itemize}
  \item void integfnend(void *data);
    \begin{itemize}
      \item data : the user value provided by integfnbegin.
    \end{itemize}
  \end{itemize}

This function is called after the integration stops.

In order to create a dynamic library from a source file written in the language C, it should usually compiled with the following command:

\begin{itemize}
  \item under linux with gcc : gcc -fPIC -shared fcn.c -o libfcn.so
  \item under linux with intel c++ : icc -fPIC -shared fcn.c -o libfcn.so
  \item under MacOS X with gcc : gcc -shared -dynamiclib fcn.c -o libfcn.dylib
  \item under MacOS X with intel c++ : icc -fPIC -dynamiclib fcn.c -o libfcn.dylib
\end{itemize}
Example:
The dynamic library libfcn.(so,dylib,dll) contains
the results of the compilation of the following source file fcn.c:
#include <math.h>
void integfcnbgn(const int * neq, void **data)
{
}

int integfcn(const int * neq, const double *x, const double *y,
        double *yp, void* data)
{
    yp[0] = sqrt(1.E0*cos(y[1]));
    yp[1] = y[0]+y[1];
    return 1;
}

void integfcnen(void *data)
{
}

To integrate the system \( \frac{dx}{dt} = \sqrt{1+\cos(y)} \), \( \frac{dy}{dt} = x+y \)
with the initial condition \( x=1 \) and \( y=0 \)
and to store the result in the file output.dat:

> integnumfcn(0,2*pi, pi/10, 1E-12, pi/10, pi/20, output.dat,
       libfcn, REAL, (x, 1),(y,0));

To integrate the same system
but it stores the result to the array of numerical vectors q:
> q = integnumfcn(0,2*pi, pi/10, 1E-12, pi/10, pi/20,
       libfcn, REAL, (x, 1),(y,0));

integnumfcn

integnumfcn(T0, TF, PAS, PREC, DOPRIMAX, PASDOPRI, FICHIER, LIBRAIRIE,
    { REAL ,(VAR1, VII),...},
    { COMPLEX ,(VAR2, VI2),...},
    evalnum, REAL ,TB, TB_0, TB_PAS,TF, TF_0, TF_PAS )

with:
- \( T0 \) : <real> : initial time
- \( TF \) : <real> : final time
- \( PAS \) : <real> : step of the output in the file
- \( PREC \) : <real> : precision
- \( DOPRIMAX \) : <real> : maximal step for the integrator
- \( PASDOPRI \) : <real> : default step for the integrator
- \( FICHIER \) : <filename> : file for the results
- \( LIBRAIRIE \) : <filename> : name of the dynamic library (without the extension .so, .dll
   or .dylib) where are defined the functions integfcnbegin, integfcn et integfcnen.
- \( VAR \) : <(array of) variables> : variable or array of variables
Chapter 17: Numerical computations

- **$VI$**: \(<\text{constant or num. vec.}>\) initial condition or numerical vector of initial conditions
- **$TB$**: \(<\text{array of real vec.}>\) array of time series for the backward time
- **$TB\_0$**: \(<\text{real}>\) initial time of $TB$
- **$TB\_PAS$**: \(<\text{real}>\) time-step of $TB$
- **$TF$**: \(<\text{array of real vec.}>\) array of time series for the forward time (between $T0$ and $TF$
- **$TF\_0$**: \(<\text{real}>\) initial time of $TF$
- **$TF\_PAS$**: \(<\text{real}>\) time-step of $TF$

It performs the numerical integration of the function integfcn located in the external library **LIBRAIRIE** with the initial conditions and time series. It stores the results to the file **FICHIER**. The numerical integrator specified by **_integnum** must be ADAMS (see Section 3.11 [._integnum], page 13).

This function supplies to the function integfcn the extracted data from the time series $TB$ or $TF$ for the time where the second member must be evaluated. For example, if the second member must be evaluated at the time $t$, it extracts from $TB$ or $TF$ the data corresponding to the time $t$. For the beginning of the ADAMS method, it interpolates the data of $TB$ and supplies these interpolated values to the function integfcn. For the time between $T0$ and $TF$, it only extracts the values without any interpolation. For example, if $TB$ and $TF$ are arrays of 4 vectors of 1000 numbers, the function integfcn will be called with a vector of 4 numbers.

The function requires that $TB\_0$ and $TF\_0$ must have the same value as $T0$. Moreover, **PAS** must be a multiple of $TB\_PAS$ and $TF\_PAS$.

The integrated variables are only used for the names in the output file.

The functions of the dynamic library must have the following definition:

- **void integfcnbegin(const int * neq, void **data)**;
  - neq : dimension of the system.
  - *data could be initialized by this function and the pointer will be furnished to integfcnend. This is a "user" value.

  This function is called before the integration starts.

- **int integfcn(const int * neq, const double *t, const double *y, double *yp, const double* tabdata)**;
  - neq : dimension of the system.
  - *t : time where the derivative must be evaluated.
  - y : a vector of neq floating-point numbers which contains the solution at the current time (*t).
  - dy : a vector of neq floating-point numbers which must be initialized by integfcn. On exit, it must contain the evaluation of the derivative of y at the current time (dy/dt=f(y,tabdata(t),t) ).
  - tabdata : a vector with the extracted values from TB or TF at the time (*t)

  This function must initialize dy such that dy/dt=f(y,tabdata(t),t) and must return 1.

- **void integfcnend(void *data)**;
  - data : the user value provided by integfcnbegin.

  This function is called after the integration stops.
In order to create a dynamic library from a source file written in the language C, it should usually be compiled with the following command:

- under Linux with gcc: gcc -fPIC -shared fcn.c -o libfcn.so
- under Linux with Intel C++: icc -fPIC -shared fcn.c -o libfcn.so
- under MacOS X with gcc: gcc -shared -dynamiclib fcn.c -o libfcn.dylib
- under MacOS X with Intel C++: icc -fPIC -dynamiclib fcn.c -o libfcn.dylib

Example:
The dynamic library libfcn.(so,dylib,dll) contains the results of the compilation of the following source file fcn.c:

```c
#include <math.h>

void integfcnbegin(const int * neq, void **data)
{
}

int integfcn(const int * neq, const double *x, const double *y,
              double *yp, const double *tabdata)
{
    yp[0] = sqrt(1.E0+cos(y[1])*tabdata[0]);
    yp[1] = y[0]*tabdata[1]+y[1];
    return 1;
}

void integfcnend(void *data)
{
}
```

To integrate the system
{ \( \frac{dx}{dt} = \sqrt{1 + \cos(y) \sin(a \cdot t)} \), \( \frac{dy}{dt} = \cos(b \cdot t) x + y \) } with \( a=0.5 \), \( b=0.9 \) and the initial condition \( x=1 \) and \( y=0 \) and to store the result in the file output.dat:

```c
> t0 = 0;
> tf = 2*pi;
> pas=pi/10;
> vnumR tabb[1:2], tabf[1:2];
> t=-30*pas,t0,pas;
> a=0.5; b=0.9;
> tb[1]=sin(a*t); tb[2]=cos(b*t);
> t=t0, tf+2*pas,pas;
> tf[1]=sin(a*t); tf[2]=cos(b*t);
>
> integnumfcn(t0,tf, pas, 1E-12,pas, pas/2, output.dat,
            libfcn, REAL, (x, 1),(y,0),
            evalnum, REAL, tb, t0, -pas, tf, t0, pas);
```

17.6 Integral

integral

```c
integral(<num. vec.> TY, <real> XH, <integer> ORDRE, <object identifier> N)
```
integral(<num. vec.> TY, <real> XH, <integer> ORDRE, <object identifier> N, "kahan")
with:
- TY : numerical vector
- XH : step between each value in the vector
- ORDRE : integration order from 1 to 6
- N : index of the last element used in TY

It computes the integral, using Newton-Cotes formulas, of order ORDRE, of the numerical vector TY where data are equally spaced of XH.

It returns in N the index of the last element used to compute the integral. The value of N verifies: N=int((size(TY)-1)/ORDRE)*ORDRE+1

If the option "kahan" is set, the function uses the Kahan method to perform the summation (compensated summation).

Reference: For Newton-Cotes formulas, see the book "Introduction to Numerical Analysis", chapter integration, of Stoer and Burlisch.

Example:
> t=0:10000;
t  Tableau de reels : nb reels =10001
> t=t*pi/10000;
> X=-cos(t)+1;
> for j=1 to 6 { s=integral(sin(t),pi/10000,j,N); delta=s-X[N];};
s = 1.99999998355066
delta = -1.64493392240672E-08
s = 1.99999999999999
delta = -7.54951656745106E-15
s = 1.99999995065198
delta = 5.55111512312578E-15
s = 2
delta = -4.44089209850063E-16
s = 2
delta = 8.88178419700125E-16
s = 1.9999921043176
delta = 5.77315972805081E-15

17.7 Numerical iteration

iternum

iternum( nbiteration, passortie, fichierresultat ,
        { REAL , ( variable1, serie1, valeurinitiale1),...},
        { COMPLEX , ( variable2, serie2, valeurinitiale2),...});

with:
- nbiteration : <integer> : number of iterations
- passortie : <integer> : output step
- fichierresultat : <filename> : result file
- variable1 : <variable> : iterated variable
- serie1 : <operation> : serie (right member)
- \textit{valeurinitiale1} : \textit{constant} : initial value of the variable

It performs \textit{nbiteration} iterations on the equations specified in lists. It writes results to file every \textit{passsortie} iterations.

The equations are triplets of the iterated variable, serie and initial value.

The initial value could be a real or a complex number.

Example:

\begin{verbatim}
> iternum(10,1,sortie.txt,REAL,(un,un+2,3));
Le fichier "sortie.txt" contient :

\begin{verbatim}
time un
+0.000000000000000E+00 +3.000000000000000E+00
  1 +5.000000000000000E+00
  2 +7.000000000000000E+00
  3 +9.000000000000000E+00
  4 +1.100000000000000E+01
  5 +1.300000000000000E+01
  6 +1.500000000000000E+01
  7 +1.700000000000000E+01
  8 +1.900000000000000E+01
  9 +2.100000000000000E+01
 10 +2.300000000000000E+01
\end{verbatim}

> iternum(5,1,sortie.txt,REAL,(vn,un-1,3),(un,vn+un,2));
Le fichier "sortie2.txt" contient :

\begin{verbatim}
time vn un
+0.000000000000000E+00 +3.000000000000000E+00 +2.000000000000000E+00
  1 +1.000000000000000E+00 +5.000000000000000E+00
  2 +4.000000000000000E+00 +6.000000000000000E+00
  3 +5.000000000000000E+00 +1.000000000000000E+01
  4 +9.000000000000000E+00 +1.500000000000000E+01
  5 +1.400000000000000E+01 +2.400000000000000E+01
\end{verbatim}
\end{verbatim}

17.8 Interpolation

\texttt{interpol} \hspace{1cm} [Procedure]

\begin{verbatim}
interpol(LINEAR , TX, DTX, TY)
interpol(QUADRATIC , TX, DTX, TY)
interpol(SPLINE , TX, DTX, TY)
interpol(HERMITE , TX, DTX, TY, DEG)
\end{verbatim}

with:
- \texttt{TX} : \texttt{<real vec.>} : vector containing the time of the known data \texttt{DTX}.
- \texttt{DTX} : \texttt{<real vec.>} : vector of known data \texttt{DTX=f(TX)}.
- \texttt{TY} : \texttt{<real vec.>} : vector of time to find the values.
- \texttt{DEG} : \texttt{<integer>} : maximal degree of the interpolating polynomial.

It performs the linear, spline, quadratic or hermitian interpolation. It returns a numerical vector of interpolated values at the time \texttt{TY} from the known data \texttt{(TX, DTX)}.

Remarks: it assumes that \texttt{TX} et \texttt{TY} are sorted in ascending or descending order.

If \texttt{TY}[i] isn't in the interval \texttt{[TX[0], TX[size(TX)]]} , then \texttt{DTY[i]} is extrapolated.

The algorithm of the hermitian interpolation is described in the Stoer and Burlish’s book, 1993, "Introduction to Numerical Analysis, Chap 2, pages 52-57".
Example:
> x1=0,2*pi,2*pi/59;
> x2=0,3,0.5;
> sl=interpol(LINEAR,x1,cos(x1),x2);
> writes("%.2E %.4E %.4E
",x2, cos(x2), sl);
0.00E+00 1.0000E+00 1.0000E+00
5.00E-01 8.7758E-01 8.7652E-01
1.00E+00 5.4030E-01 5.3958E-01
1.50E+00 7.0737E-02 7.0719E-02
2.00E+00 -4.1615E-01 -4.1576E-01
2.50E+00 -8.0114E-01 -8.0001E-01
3.00E+00 -9.8999E-01 -9.8920E-01

17.9 Coordinate system conversions

17.9.1 ell_to_xyz

def ell_to_xyz:
    ell_to_xyz( ELL, MU, X, XP );

with:

- **ELL**: array of 6 vectors of real numbers which contain the elliptic coordinates.
  - ELL[1]: semi-major axis
  - ELL[2]: mean longitude
  - ELL[3]: eccentricity * cos(longitude of the pericenter)
  - ELL[4]: eccentricity * sin(longitude of the pericenter)
  - ELL[5]: sin(inclination/2) * cos(longitude of the ascending node)
  - ELL[6]: sin(inclination/2) * sin(longitude of the ascending node)
- **MU**: product of gravitation constant and the sum of the masses.
- **X**: array of 3 vectors of real numbers which contain the positions.
- **XP**: array of 3 vectors of real numbers which contain the velocities.

It converts the elliptic coordinates ELL to cartesian coordinates X (positions) and XP (velocities).

Example:
/* calcul d'une seule valeur */
GM= 0.2959122082855911d-03;
mu = GM*1.05;
N=2*pi/221.6;
varpi = 138*pi/180;

vnumR ell[1:6];
resize(ell,1);
ell[1][1] = (mu/n**2)**(1./3.);
ell[2][1] = varpi;
ell[3][1] = 0.54*cos(varpi);
ell[4][1] = 0.54*sin(varpi);
ell[5][1] = 0.0;
ell[6][1] = 0.0;
ell_to_xyz(ell, mu, x, xp);
writes(x);
writes(xp);

Example:
/* calcul de plusieurs valeurs contenues dans un fichier */
mu = 1;

vnumR ell[1:6];
read(file1, ell);

ell_to_xyz(ell, mu, x, xp);

writes(x);
writes(xp);

17.9.2 xyz_to_ell

xyz_to_ell(x, xp, mu, ell);[Procedure]

with:
- X : <array of real vec.> : array of 3 vectors of real numbers which contain the positions.
- XP : <array of real vec.> : array of 3 vectors of real numbers which contain the velocities.
- MU : <real> : product of gravitation constant and the sum of the masses.
- ELL : <array of real vec.> : array of 6 vectors of real numbers which contain the elliptic coordinates.
  - ELL[1] : semi-major axis
  - ELL[2] : mean longitude
  - ELL[3] : eccentricity * cos(longitude of the pericenter)
  - ELL[4] : eccentricity * sin(longitude of the pericenter)
  - ELL[5] : sin(inclination/2) * cos(longitude of the ascending node)
  - ELL[6] : sin(inclination/2) * sin(longitude of the ascending node)

It converts the cartesian coordinates X (positions) and XP (velocities) to the elliptic coordinates ELL.

Example:
mu = 1;

vnumR x[1:3], xp[1:3];
read(file2, x, xp);

xyz_to_ell(x, xp, mu, ell);

writes(ell);
18 Utilities

18.1 help

help
help;
?;
Display help.
On Unix and MacOS X operating systems, the help program uses the software info (provided by Texinfo) to display help. On Windows operating systems, the help file (chm format) is displayed.

help
help <motreserve>;
? <motreserve>;
It displays the help in the reference manual for the keyword of TRIP. On Unix and MacOS X operating systems, this function calls the program info (provided by Texinfo).

Example :
> ? vartrip;

18.2 exit

exit
quit;
exit;
Stop the execution of TRIP.

18.3 cls

cls
cls;
Clear the current screen (similar to the Unix command clear).

18.4 pause

pause
pause;
Display a message, suspend execution until the user hits a key. If the user hits the key RET, then the execution continues. If the user hits the key q+RET or e+RET, then the execution stops and the user returns to the prompt.
pause(<integer> x);
Suspend execution for x seconds.

Example :
> for j=1 to 3 { j; pause; time_s; pause(2); time_t; };
1
Appuyez sur la touche ‘return’ pour continuer
ou la touche ‘q’ ou ‘e’ pour revenir au prompt ...
18.5 msg

msg

msg <string> ;

msg(<string> textformat);

Display unformatted messages to the screen.

This message must be a string or a text between double-quotes. The messages could be on several lines.

msg

msg(<string> textformat, <real> x, ... );

Display formatted messages to the screen (with/out) real constants.

The real constants must be formatted. The format is the same as the command printf in language C (see Section 7.6 [str], page 33, for the valid conversion specifiers). This message must be a string or a text between double-quotes. The messages could be on several lines.

To display double-quotes, two double-quotes must be used.

Under the numerical mode NUMRATMP only, the integers or rational numbers are converted to double-precision floating-point numbers before they are displayed if the format is '%g', '%e' or '%f'. if the format is '%d' or '%i', the integers or rational numbers are written without any conversion.

The function msg (see Section 7.7 [msg], page 34) has the same behavior but writes the result into a string of characters.

Example :

> file="fichier1.dat"
> msg("écriture du fichier "+file;
écriture du fichier fichier1.dat
> msg "je vais faire un guillemet :
""cette chaine est encadre par des guillemets"" ;

je vais faire un guillemet :
cette chaine est encadre par des guillemets".

> msg("pi=%g pi=%E\n",pi,pi);
pi=3.14159 pi=3.14159265E+00
18.6 error

error[Procedure]

error(<string> textformat);

Raise an error and display the unformatted text to the screen.
This message must be a string or a text between double-quotes. The message could be on several lines.

Example :
> x=2;
> if (x==2) then { error("raise an error : x=2"); };
TRIP[ 3 ] : raise an error : x=2
Commande : 'if (x==2) then { error("raise an error : x=2"); }';

18.7 try

try[Procedure]

try { <body> bodytry } catch { <body> bodycatch };

Normally, it executes all statements of bodytry. If an error occurs during this execution, then it ignores the remaining statements of bodytry and it executes the statements of bodycatch.

If the global variable _info is equal to on, then an information message is displayed when the error occurs in the statements of bodytry. If the global variable _info is equal to off, then no information message is displayed when the error occurs in the statements of bodytry.

If an error occurs during the execution of the statements of bodycatch then an error is generated. This error could be caught if the current statement try-catch is inside another statement try-catch.

It is allowed to imbricate a statement try-catch inside the statements of bodytry or of bodycatch.

Example :
> // displays a warning message but it continues
> _info on;
> _info on
> b = 0;
> b = 0
> q = 2;
q = 2
> try { s = "invalid"+q; b = 2; } catch { b = 1; };
TRIP[ 1 ] : Can't add
Command : ' try { s = "invalid"+q; b = 2; } catch { b = 1; }';

Information : Continue on error (catch statement)
b = 1
>
> // does not display a warning message but it continues
> _info off;
> _info off
> b = 0;
b = 0
> q = 2;
```plaintext
q = 2
> try { s = "invalid"+q; b = 2; } catch { b = 1; }
> b = 1
```

### 18.8 delete

delete [Procedure]
delete( <object identifier> );
Delete any object (serie, variable, array, ...).

Remarks:
- Any dependencies are set to 0.
- Angular variables depending on this object will be converted to polynomial variable.
- Truncation depending on this object are deleted.

Example:
> file="fichier1.dat";
file = "fichier1.dat"
> X=expi(x,1,0);
1*X

> delete(file);
> delete(x);
La variable angulaire X est convertie en variable polynomiale.

### 18.9 reset

reset [Procedure]
reset;
Reinitialize TRIP : TRIP global variables are set to default values and delete all object identifiers in memory.

### 18.10 include

include [Procedure]
include <filename> ;
include <string> ;
It loads the file "fichier" and run it.
This file must be located in the current directory or in the directory specified by the variable _path_.
The file extension could have any value. The recomended file extension is .t .
Remarks: If the file is an object identifier of type string, then it loads the file which its name is the contents of the string.

Example:
> include ellip;
> include fct.t;
> file="fperplanumH";
file = "fperplanumH"
> include file; /*exèécute le fichier "fperplanumH" */
18.11 @@

@@ [Procedure]
@@;
Reinitialize TRIP. Then it loads and executes the last file executed by the command `include`.

18.12 vartrip

`vartrip` [Procedure]
`vartrip;`
Displays the status of all TRIP global variables.

Example :
> vartrip;

Etat des variables globales de Trip :

```plaintext
_affc = 6
_affdist = 0
_affhomog = 0
_echo = 0
_path = /exemples/
_history = /unixfiles/
_hist = ON
_naf_iprt = -1
_naf_icplx = 1
_naf_iw = 1
_naf_isec = 1
_naf_dtour = 6.283185307179586E+00
_naf_nulin = 1
_naf_tol = 1.000000000000000E-10
_time = OFF
_comment = OFF
_affvar = ( )
```

18.13 bilan

`bilan` [Procedure]
`bilan;`
Displays all object identifiers in memory with their type :

SERIE the object is a serie.
VARIABLE the object is a variable.
CONST the object is a constant.
MATRIXR the object is a numerical matrix of real numbers.
MATRIXC the object is a numerical matrix of complex numbers.
TAB the object is an array.
TABVAR the object is an array of variables.
VNUMR the object is a numerical vector of real numbers.
VNUMC the object is a numerical vector of complex numbers.
STRING the object is a string.
EXTERNAL the object is an external structure.
FILE the object is a file.
REMOTE the object is a remote object stored on a SCSCP server.
OBJECT
SCSCP CLIENT the object is a connection to a SCSCP server.

bilan mem

bilan mem;
For each serie or array of series, it displays the memory used and the number of terms.
The total number of terms and total memory used for series are displayed.

Example :
> bilan;
A VAR
ASR SERIE
ASRp SERIE
RSA SERIE
X VAR
_AsR SERIE
_CosE SERIE
_Expiv SERIE
_RsA SERIE
_RsASinv SERIE
_SinE SERIE
e VAR
ep VAR
g VAR
ga VAR
lp VAR
x VAR
> bilan mem;

<table>
<thead>
<tr>
<th>Nom</th>
<th>Memoire (octets)</th>
<th>Nb de termes</th>
</tr>
</thead>
<tbody>
<tr>
<td>ASR</td>
<td>1704</td>
<td>64</td>
</tr>
<tr>
<td>ASRp</td>
<td>1704</td>
<td>64</td>
</tr>
<tr>
<td>RSA</td>
<td>2136</td>
<td>81</td>
</tr>
<tr>
<td>_AsR</td>
<td>1704</td>
<td>64</td>
</tr>
<tr>
<td>_CosE</td>
<td>2112</td>
<td>80</td>
</tr>
<tr>
<td>_Expiv</td>
<td>1728</td>
<td>64</td>
</tr>
<tr>
<td>_RsA</td>
<td>2136</td>
<td>81</td>
</tr>
<tr>
<td>_RsACosv</td>
<td>2112</td>
<td>80</td>
</tr>
<tr>
<td>_RsAEexpiv</td>
<td>1728</td>
<td>64</td>
</tr>
<tr>
<td>_RsASinv</td>
<td>2112</td>
<td>80</td>
</tr>
<tr>
<td>_SinE</td>
<td>2112</td>
<td>80</td>
</tr>
<tr>
<td>---------</td>
<td>-----------------</td>
<td>--------------</td>
</tr>
</tbody>
</table>

21288 802

18.14 stat

stat

stat;
It displays all objects in memory. These objects are ordred by their type. For each object, it
displays a small description. For the series, it displays their contents.
Example:
> S = (1+x+y)^2$
> vnumR R$
> vnumC C$
> dim T[1:2]$
> stat;

Summary of objects:
Series:
\[
S(x,y) = 1 + 2y + 1y^2 + 2x + 2xy + 1x^2
\]

Variables:
Variable x type : 2 ordres : 2 2 2 2
dependances:
variables dependant de celle-ci :

Variable y type : 2 ordres : 3 3 3 3
dependances:
variables dependant de celle-ci :

Arrays of series:
T [1:2 ] number of elements = 2

Double precision real vectors:
Numerical vector R contains 0 double precision reals.

Double precision complex vectors:
Numerical vector C contains 0 double precision complex.

> 

[Procedure]
\[stat( <object identifier> );\]
\[stat( <object identifier> , "puismin" );\]
\[stat( <object identifier> , "puismax" );\]
\[stat( <object identifier> , "puismin" , "puismax" );\]
Depending on the object type, it displays information on number of terms and the memory used by the object.
If the option "puismin" is set, the function displays the minimal degree of each variable in the serie.
If the option "puismax" is set, the function displays the maximal degree of each variable in the serie.

Example :
> s = (1+x+y)^2$
> stat(s);
serie s ( x , y )
number of variables : 2 size of the descriptor : 96 bytes
number of terms : 6 size : 608 bytes
> stat(s,"puismin","puismax");
serie s ( x , y )
number of variables : 2 size of the descriptor : 96 bytes
number of terms : 6 size : 608 bytes
puismin : x ^ 0 , y ^ 0
puismax : x ^ 2 , y ^ 2

18.15 save_env

save_env

save_env;
Save TRIP global variables to a stack.

Example :
> _mode;
  _mode = POLP
> save_env;
> _mode=POLH;
  _mode = POLH
> rest_env;
> _mode;
  _mode = POLP

18.16 rest_env

rest_env

rest_env;
Restore the TRIP global variables saved by save_env from a stack.

Example :
> _path="/users/";
  _path = /users/
> save_env;
> _path="/users/toto/";
  _path = /users/toto/
> /*instructions utilisant le chemin specifique */;
> rest_env;
> _path;
  _path = /users/

18.17 random

random

random(<integer> x)
Return a pseudo-random number in the range from 0 and x-1.
Example:
> random(10);
   1
> random(10);
   7
> random(10);
   0
> random(10);
   9
> random(10);
   8

18.18 `nameof`

`nameof` [Function]

`nameof(<object identifier> x)`

It returns as a string the name of the object.

If the object is an expression, then the empty string "" is returned.

Example:
> c=3$
> n1 = nameof(c);
   n1 = "c"
> s="abcde"$
> n2 = nameof(s);
   n2 = "s"

18.19 `typeof`

`typeof` [Function]

`typeof(<object identifier> x)`

It returns as a string the type of the object. The returned strings are described in the command `bilan` (see Section 18.13 [bilan], page 201).

If the object doesn’t exist, then the empty string "" is returned.

Example:
> typeof(2);
   "CONST"
> typeof(1+x);
   "SERIE"
> t=1,10;
   t double precision real vector : number of elements =10
> if (typeof(t)="VNUMR") then { stat(t); }
   Numerical vector t contains 10 double precision reals.
   Size of the array in bytes: 80

18.20 `file_fullname`

`file_fullname` [Function]

`file_fullname(<string> name)`
It builds an absolute file name. All later usage of the returned string will ignore the variable \_path.

Example:
\>
\>_path;
\_path = /users/guest/data/
\> vnumR t;
\> fn=file_fullname("/tmp/mydata.txt");
\fn = nom de fichier '/tmp/mydata.txt'
\> read(fn,t);

18.21 Execution time

18.21.1 time_s

time_s

[Procedure]
time_s;

Initialize the starting time for computations in functions time_l and time_t.

18.21.2 time_l

time_l

[Procedure]
time_l;

Displays the consumed user CPU time, the elapsed real time, and the consumed system CPU time, since the last call to time_l.

If no previous call to time_l was performed, then it displays the user time used since the last call to time_s.

Example:
\>
\> time_s; for j=1 to 3 { (1+x+y+z+t+u+v)**18$ time_l; }; time_l;
utilisateur 00.313s - reel 00.183s - systeme 00.057s - (202.43% CPU)
utilisateur 00.307s - reel 00.170s - systeme 00.055s - (213.11% CPU)
utilisateur 00.311s - reel 00.173s - systeme 00.055s - (211.93% CPU)

18.21.3 time_t

time_t

[Procedure]
time_t;

Displays the consumed user CPU time, the elapsed real time, and the consumed system CPU time, since the last call to time_s.

time_t(<object identifier> usertime , <object identifier> realtime);

Displays the consumed user CPU time, the elapsed real time, and the consumed system CPU time, since the last call to time_s. usertime will contain the sum of the user and system CPU time and realtime the elapsed real time.

Example:
\>
\> time_s; for j=1 to 3 { (1+x+y+z+t+u+v)**18$ time_t; }; time_t(usertime, realtime);
utilisateur 00.310s - reel 00.178s - systeme 00.055s - (205.29% CPU)
utilisateur 00.615s - reel 00.347s - systeme 00.111s - (209.44% CPU)
utilisateur 00.928s - reel 00.519s - systeme 00.168s - (211.23% CPU)
utilisateur 00.928s - reel 00.519s - systeme 00.168s - (211.20% CPU)
18.22 Call the shell

![string> ;

Execute a shell command specified in a string.

Example :
> dir = "ls -al";
dir = "ls -al"
> ! dir;
total 8136
drw-r-sr-x 20 xxxx ttt 1536 Sep 27 17:16 .
drw-r-sr-x 22 xxxx ttt 512 Sep 06 12:16 ..
-rw-r----- 1 xxxx ttt 281 Sep 08 1998 .Guidefaults
-rw------- 1 xxxx ttt 204 Sep 27 10:46 .Xauthority

![string> ;

Execute a shell command specified in a string and return the standard output as a string. If the error output is not empty, an error is raised.

Example :
> s=str(!"uname");
s = "Darwin"
19 References

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  version 6.0.0, 2014
  Torbjorn Granlund et al.
  http://www.gmplib.org/

- **LTDL - GNU Libtool - The GNU Portable Library Tool**
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  http://www.symcomp.org/scscp

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  D. Roozemond
  http://www.win.tue.nl/SCIEnce/cds/scscp1.html

- **OpenMath content dictionary scscp2**
  D. Roozemond
  http://www.win.tue.nl/SCIEnce/cds/scscp2.html

- **SCSCP C Library - A C/C++ library for Symbolic Computation Software Composibility Protocol**
  Version 1.0.2, May 2016
  M. Gastineau
  http://www.imcce.fr/trip/scscp/
Appendix A Supported OpenMath Content Dictionaries

The following list are the OpenMath objects supported by the SCSCP server and client.

<table>
<thead>
<tr>
<th>CD</th>
<th>Symbol</th>
</tr>
</thead>
<tbody>
<tr>
<td>arith1</td>
<td>abs, divide, minus, plus, power, times, unary_minus</td>
</tr>
<tr>
<td>alg1</td>
<td>one, zero</td>
</tr>
<tr>
<td>bigfloat1</td>
<td>bigflat</td>
</tr>
<tr>
<td>arith1</td>
<td>abs, divide, minus, plus, power, times, unary_minus</td>
</tr>
<tr>
<td>complex1</td>
<td>argument, complex_cartesian, complex_polar, conjugate, imaginary, real</td>
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<tr>
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<td>C, Q, R</td>
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<tr>
<td>interval1</td>
<td>interval_cc, integer_interval, interval</td>
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<tr>
<td>linalg2</td>
<td>matrix, matrixrow, vector</td>
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<tr>
<td>list1</td>
<td>list</td>
</tr>
<tr>
<td>logic1</td>
<td>not, and, xor, or, true, false</td>
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<tr>
<td>nums1</td>
<td>e, i, infinity, NaN, pi, rational</td>
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<td>polyd1</td>
<td>poly_ring_d_named, SDMP, DMP, term</td>
</tr>
<tr>
<td>polyu</td>
<td>poly_u_rep, term</td>
</tr>
<tr>
<td>polyr</td>
<td>poly_r_rep, term</td>
</tr>
<tr>
<td>setname1</td>
<td>C, N, P, Q, R, Z</td>
</tr>
<tr>
<td>transc1</td>
<td>arccos, arccosh, arcsin, arcsinh, arctan, arctanh, cos, cosh, exp, ln, log, sin, sinh, tan, tanh</td>
</tr>
</tbody>
</table>

The following list are the OpenMath objects supported only by the SCSCP server.

<table>
<thead>
<tr>
<th>CD</th>
<th>Symbol</th>
</tr>
</thead>
<tbody>
<tr>
<td>scscp2</td>
<td>get_allowed_heads, get_transient_cd, get_signature, store, retrieve, unbind</td>
</tr>
</tbody>
</table>

The following list are the OpenMath objects supported only by the SCSCP client.

<table>
<thead>
<tr>
<th>CD</th>
<th>Symbol</th>
</tr>
</thead>
<tbody>
<tr>
<td>scscp2</td>
<td>symbol_set, signature, service_description,</td>
</tr>
</tbody>
</table>
Appendix B Index of the global variables

- _affc ............................................. 9
  _affdist ....................................... 10
  _comment ....................................... 10
  _cpu ............................................ 11
  _echo ........................................... 11
  _endian ......................................... 12
  _graph .......................................... 12
  _hist ........................................... 12
  _history ........................................ 12
  _info ........................................... 13
  _integnum ...................................... 13
  _language ....................................... 13
  _mode ........................................... 14
  _modenum ....................................... 15
  _modenum_prec .................................. 16
  _naf_dtour ..................................... 16
  _naf_icplx ..................................... 17
  _naf_iprt ...................................... 17
  _naf_iw ........................................ 18
  _naf_nulin .................................... 18
  _naf_tol ........................................ 18
  _path ........................................... 18
  _quiet .......................................... 19
  _read ........................................... 19
  _read_history .................................. 20
  _time ........................................... 21
  _userlibrary_path .............................. 21
  _naf_iv ........................................ 18
  _naf_nulin ..................................... 18
  _naf_tol ........................................ 18
  _path ........................................... 18
  _quiet .......................................... 19
  _read ........................................... 19
  _read_history .................................. 20
  _time ........................................... 21
  _userlibrary_path .............................. 21

I
  i ................................................. 22
  I ................................................. 22

P
  pi ................................................. 22
  PI ............................................... 22
Appendix C Index of the procedures

! .................................................................................. 207
!= ................................................................................. 32
% ................................................................................. 134
& ................................................................................. 44, 103
&* ................................................................................. 44, 103
( .................................................................................. 201
* ................................................................................. 25, 31, 47, 107
** .................................................................................. 25
**1 ................................................................................. 45
+ ................................................................................. 25, 31, 47, 107
, .................................................................................. 57
- .................................................................................. 25, 47, 107
/ .................................................................................. 25, 47, 107
: .................................................................................. 23, 40
< .................................................................................. 32
<= .................................................................................. 32
= .................................................................................. 39
== ................................................................................. 32
> .................................................................................. 33
>= .................................................................................. 33
?
?.................................................................................. 197
?():: ................................................................................. 93
@
@ .................................................................................. 136
@<string>@valuemyspace@ ........................................ 115, 116
[ 
[::] ................................................................. 100
[::] ................................................................. 63
[::] ................................................................. 62
^ .................................................................................. 25
\ .................................................................................. 115, 116, 120, 123
A
abs................................................................. 82
accum......................................................... 79
acos..................................................................... 87
acosh............................................................. 89
affmac.......................................................... 136
afftab............................................................ 42
arg................................................................. 82
asin................................................................. 88
asinh............................................................. 89
atan............................................................... 88
atan2............................................................. 90
atanh............................................................ 90
C
case............................................................ 145
catch............................................................. 199
close............................................................. 30
cls................................................................. 197
cof_ext ......................................................... 26
cof_num ....................................................... 27, 34
conj............................................................... 84
cos................................................................. 86
cosh............................................................. 88
crevar ........................................................... 23
D
delete ......................................... 200
deriv ......................................... 26
det ........................................... 44, 104
dim ............................................ 37
dimtovnumC .................................. 94
dimtovnumR .................................. 94
dimvar ....................................... 38
div ............................................ 26

e
ecriture ...................................... 29
effmac ....................................... 136
effmacros ................................... 136
eigenvalues .................................. 45, 106
eigenvectors ................................ 46, 106
eii_to_xyz ................................... 195
end .......................................... 115, 116, 120, 123
erf ........................................... 84
erf ........................................... 85
erf ........................................... 85
evalnum ...................................... 197
exit .......................................... 27
exp ........................................... 85
extern_display ................................ 131
extern_function ............................. 129
extern_lib .................................. 130
extern_type ................................ 131

F
fac ........................................... 29, 86
fft ........................................... 182
file_close ................................... 72
file_fullname ................................. 205
file_open .................................... 72
file_read .................................... 73
file_readappend .............................. 74
file_write ................................... 72
file_writebin ................................ 75
for ........................................... 139
freqa ........................................ 179
freqareson .................................. 181

G
gnuplot ...................................... 115
grace ........................................ 116

H
help .......................................... 197
histogram .................................... 92

I
identitymatrix ................................ 106
if .............................................. 141
ifft .......................................... 183
imag .......................................... 83
imax .......................................... 76
IMAX .......................................... 77
imin .......................................... 76
IMIN .......................................... 77
include ...................................... 200
inf ........................................... 41, 61
int ............................................ 90
integ ......................................... 26
integnum ..................................... 183
integnumfcn ................................. 188
integral ...................................... 192
interpol ...................................... 194
intersectnum ................................. 80
invertmatrix ................................ 104, 105
itemum ....................................... 193

K
kroneckerproduct ............................. 104

L
lapack_dgbsv ................................ 148
lapack_dgeev ................................. 172
lapack_dgels ................................ 156
lapack_dgelss ............................... 158
lapack_dgeqlf ............................... 165
lapack_dgeqr ............................... 164
lapack_dgesdd ............................... 168
lapack_dgesvd ............................... 147
lapack_dgesvd ................................ 168
lapack_dgges ................................ 174
lapack_dggev ................................ 173
lapack_dggglm ............................... 160
lapack_dggse ................................ 159
lapack_dgtsv ................................ 148
lapack_dpbsv ................................ 150
lapack_dpdsv ................................ 149
lapack_dpotrf ............................... 166
lapack_dptsv ................................ 150
lapack_dsbev ................................ 170
lapack_dstev ................................ 171
lapack_dsyev ................................ 170
lapack_dsysv ................................ 172
lapack_dsysv ................................ 149
lapack_zgbsv ................................ 151
lapack_zggev ................................ 175
lapack_zgels ................................ 161
lapack_zgelss ............................... 162
lapack_zgeqlf ............................... 167
lapack_zgeqr ................................ 166
lapack_zgesdd ............................... 169
lapack_zgesv ................................ 151
lapack_zgesvd ............................... 169
lapack_zggglm ............................... 163
lapack_zggse ................................ 163
lapack_zgtsv ................................ 152
lapack_zhbev ................................ 175
lapack_zheev ................................ 174
Appendix C: Index of the procedures  217

<table>
<thead>
<tr>
<th>Procedure</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>lapack_zhegv</td>
<td>176</td>
</tr>
<tr>
<td>lapack_zhesv</td>
<td>153</td>
</tr>
<tr>
<td>lapack_zpbsv</td>
<td>154</td>
</tr>
<tr>
<td>lapack_zposv</td>
<td>154</td>
</tr>
<tr>
<td>lapack_zpotrf</td>
<td>167</td>
</tr>
<tr>
<td>lapack_zptsv</td>
<td>155</td>
</tr>
<tr>
<td>lapack_zsysv</td>
<td>153</td>
</tr>
<tr>
<td>lecture</td>
<td>29</td>
</tr>
<tr>
<td>log</td>
<td>91</td>
</tr>
<tr>
<td>log10</td>
<td>91</td>
</tr>
<tr>
<td>M</td>
<td></td>
</tr>
<tr>
<td>macro</td>
<td>133</td>
</tr>
<tr>
<td>MACRO</td>
<td>133</td>
</tr>
<tr>
<td>maple</td>
<td>120</td>
</tr>
<tr>
<td>maple_get</td>
<td>120</td>
</tr>
<tr>
<td>maple_put</td>
<td>119</td>
</tr>
<tr>
<td>mathematica</td>
<td>123</td>
</tr>
<tr>
<td>mathematica_get</td>
<td>122</td>
</tr>
<tr>
<td>mathematica_put</td>
<td>121</td>
</tr>
<tr>
<td>matrixC</td>
<td>97</td>
</tr>
<tr>
<td>matrixC[,...]</td>
<td>98</td>
</tr>
<tr>
<td>matrixR</td>
<td>97</td>
</tr>
<tr>
<td>matrixR[,...]</td>
<td>98</td>
</tr>
<tr>
<td>max</td>
<td>76</td>
</tr>
<tr>
<td>MAX</td>
<td>78</td>
</tr>
<tr>
<td>min</td>
<td>76</td>
</tr>
<tr>
<td>MIN</td>
<td>77</td>
</tr>
<tr>
<td>mod</td>
<td>90</td>
</tr>
<tr>
<td>msg</td>
<td>34, 198</td>
</tr>
<tr>
<td>N</td>
<td></td>
</tr>
<tr>
<td>naf</td>
<td>177</td>
</tr>
<tr>
<td>naftab</td>
<td>178</td>
</tr>
<tr>
<td>nameof</td>
<td>205</td>
</tr>
<tr>
<td>nint</td>
<td>91</td>
</tr>
<tr>
<td>num_dim</td>
<td>41</td>
</tr>
<tr>
<td>NUMDBL</td>
<td>15</td>
</tr>
<tr>
<td>NUMDBLINT</td>
<td>15</td>
</tr>
<tr>
<td>NUMFPMP</td>
<td>15</td>
</tr>
<tr>
<td>NUMQUAD</td>
<td>15</td>
</tr>
<tr>
<td>NUMQUADINT</td>
<td>15</td>
</tr>
<tr>
<td>NUMRAT</td>
<td>15</td>
</tr>
<tr>
<td>NUMRATMP</td>
<td>15</td>
</tr>
<tr>
<td>O</td>
<td></td>
</tr>
<tr>
<td>off</td>
<td>10, 11, 12, 13, 21</td>
</tr>
<tr>
<td>on</td>
<td>10, 11, 12, 13, 21</td>
</tr>
<tr>
<td>P</td>
<td></td>
</tr>
<tr>
<td>pause</td>
<td>197</td>
</tr>
<tr>
<td>plot</td>
<td>111</td>
</tr>
<tr>
<td>plotf</td>
<td>114</td>
</tr>
<tr>
<td>plots</td>
<td>114</td>
</tr>
<tr>
<td>plotps_end</td>
<td>114</td>
</tr>
<tr>
<td>plotreset</td>
<td>115</td>
</tr>
<tr>
<td>POLP</td>
<td>14</td>
</tr>
<tr>
<td>POLPV</td>
<td>14</td>
</tr>
<tr>
<td>POLY</td>
<td>14</td>
</tr>
<tr>
<td>POLYV</td>
<td>14</td>
</tr>
<tr>
<td>print</td>
<td>29</td>
</tr>
<tr>
<td>private</td>
<td>7</td>
</tr>
<tr>
<td>prod</td>
<td>79, 141</td>
</tr>
<tr>
<td>public</td>
<td>8</td>
</tr>
<tr>
<td>Q</td>
<td></td>
</tr>
<tr>
<td>quit</td>
<td>197</td>
</tr>
<tr>
<td>R</td>
<td></td>
</tr>
<tr>
<td>random</td>
<td>204</td>
</tr>
<tr>
<td>read</td>
<td>29, 63</td>
</tr>
<tr>
<td>readappend</td>
<td>67</td>
</tr>
<tr>
<td>readbin</td>
<td>69</td>
</tr>
<tr>
<td>real</td>
<td>83</td>
</tr>
<tr>
<td>replot</td>
<td>112</td>
</tr>
<tr>
<td>reset</td>
<td>200</td>
</tr>
<tr>
<td>resize</td>
<td>61</td>
</tr>
<tr>
<td>rest_env</td>
<td>204</td>
</tr>
<tr>
<td>return</td>
<td>135</td>
</tr>
<tr>
<td>reversevm</td>
<td>81</td>
</tr>
<tr>
<td>S</td>
<td></td>
</tr>
<tr>
<td>saue_c</td>
<td>101</td>
</tr>
<tr>
<td>saue_fortran</td>
<td>101</td>
</tr>
<tr>
<td>saue_ml</td>
<td>102</td>
</tr>
<tr>
<td>saue_t</td>
<td>101</td>
</tr>
<tr>
<td>save_env</td>
<td>204</td>
</tr>
<tr>
<td>scscp_close</td>
<td>126</td>
</tr>
<tr>
<td>scscp_connect</td>
<td>125</td>
</tr>
<tr>
<td>scscp_delete</td>
<td>127</td>
</tr>
<tr>
<td>scscp_disable�</td>
<td>129</td>
</tr>
<tr>
<td>scscp_execute</td>
<td>128</td>
</tr>
<tr>
<td>scscp_get</td>
<td>127</td>
</tr>
<tr>
<td>scscp_put</td>
<td>126</td>
</tr>
<tr>
<td>select</td>
<td>62</td>
</tr>
<tr>
<td>sertrig</td>
<td>181</td>
</tr>
<tr>
<td>sign</td>
<td>92</td>
</tr>
<tr>
<td>sin</td>
<td>87</td>
</tr>
<tr>
<td>sinh</td>
<td>88</td>
</tr>
<tr>
<td>size</td>
<td>25, 35, 40</td>
</tr>
<tr>
<td>sort</td>
<td>81</td>
</tr>
<tr>
<td>sqrt</td>
<td>86</td>
</tr>
<tr>
<td>stat</td>
<td>202</td>
</tr>
<tr>
<td>stop</td>
<td>141</td>
</tr>
<tr>
<td>str</td>
<td>33, 95</td>
</tr>
<tr>
<td>sum</td>
<td>79, 140</td>
</tr>
<tr>
<td>sup</td>
<td>41, 62</td>
</tr>
<tr>
<td>switch</td>
<td>145</td>
</tr>
</tbody>
</table>
T

```
T

tabvar ........................................... 39
tan ............................................... 87
tanh ............................................... 89
time_L ........................................... 206
time_s ........................................... 206
transposematrix ................................. 105
transposevnum ................................. 82
try .............................................. 199
typeof ........................................... 205

U

unionvnum ........................................ 80

V

```
```
V

vartrip ........................................... 201
vnumC ............................................ 56
vnumC[...,...] ................................ 59
vnumQ ............................................ 57
vnumR ............................................ 55

W

```
```
V

while ............................................... 139
write ............................................ 68
writebin ........................................ 70
writes ........................................... 59, 99

X

```
```
V

xyz_to_ell ................................. 196

vnumR[...,...] ............................... 58
vnumtodim ..................................... 94
Appendix D  Full index

(Index is nonexistent)
# Table of Contents

1 Introduction ........................................................................................................... 1

2 Semantic .................................................................................................................. 3
   2.1 expression ............................................................................................................ 3
   2.2 object identifier .................................................................................................. 3
      2.2.1 serie .............................................................................................................. 5
      2.2.2 constant ......................................................................................................... 5
      2.2.3 string ............................................................................................................. 5
      2.2.4 real ............................................................................................................... 6
      2.2.5 complex ......................................................................................................... 6
      2.2.6 filename ......................................................................................................... 6
      2.2.7 file ............................................................................................................... 6
   2.3 Visibility .............................................................................................................. 6
      2.3.1 private .......................................................................................................... 7
      2.3.2 public .......................................................................................................... 8

3 Global variables ...................................................................................................... 9
   3.1 _affc .................................................................................................................... 9
   3.2 _affdist ............................................................................................................... 10
   3.3 _comment .......................................................................................................... 10
   3.4 _cpu .................................................................................................................... 11
   3.5 _echo ................................................................................................................... 11
   3.6 _ endian ............................................................................................................... 12
   3.7 _graph ............................................................................................................... 12
   3.8 _hist ..................................................................................................................... 12
   3.9 _history .............................................................................................................. 12
   3.10 _info ................................................................................................................ 13
   3.11 _integnum ......................................................................................................... 13
   3.12 _language ......................................................................................................... 13
   3.13 _mode ............................................................................................................... 14
   3.14 _modenum ....................................................................................................... 15
   3.15 _modenum_prec .............................................................................................. 16
   3.16 _naf_dtour ........................................................................................................ 16
   3.17 _naf_icplx ......................................................................................................... 17
   3.18 _naf_iprt .......................................................................................................... 17
   3.19 _naf_iisec ........................................................................................................ 17
   3.20 _naf_iw .............................................................................................................. 18
   3.21 _naf_nulin ....................................................................................................... 18
   3.22 _naf_tol .......................................................................................................... 18
   3.23 _path ................................................................................................................. 18
   3.24 _quiet ............................................................................................................... 19
   3.25 _read ................................................................................................................. 19
   3.26 _read_history ................................................................................................. 20
   3.27 _time ................................................................................................................ 21
   3.28 _userlibrary_path ......................................................................................... 21
   3.29 I ......................................................................................................................... 22
   3.30 PI ....................................................................................................................... 22
4 Variables ............................................................... 23
  4.1 crevar ............................................................ 23
  4.2 Operators .......................................................... 23

5 Series ................................................................. 25
  5.1 Operators .......................................................... 25
  5.2 Standard functions .............................................. 25
  5.3 Differentiation and integration ............................... 26
    5.3.1 deriv ........................................................ 26
    5.3.2 integ ....................................................... 26
  5.4 Euclidian division .............................................. 26
  5.5 Selection ........................................................ 26
    5.5.1 coef_ext .................................................. 26
  5.6 Evaluation ...................................................... 27
    5.6.1 coef_num ................................................ 27
    5.6.2 evalnum .................................................. 27

6 Constants .......................................................... 29
  6.1 Standard functions ............................................ 29
    6.1.1 factorial ................................................. 29
  6.2 Input/output on real numbers ................................. 29

7 Strings .............................................................. 31
  7.1 Declaration and assignment .................................... 31
  7.2 Concatenation .................................................. 31
  7.3 Repetition ...................................................... 31
  7.4 Extraction ...................................................... 31
  7.5 Comparaison .................................................... 32
  7.6 Conversion from an integer, a real or a serie to a string .. 33
  7.7 Conversion from a list of integers or reals to a string .... 34
  7.8 Conversion from a string to an integer or a real .......... 34
  7.9 Length of a string ............................................ 35

8 Arrays .............................................................. 37
  8.1 Declaration of array of series .................................. 37
  8.2 Initialization of an array of series ........................... 37
  8.3 Declaration of array of variables ............................. 38
  8.4 Initialization of an array of variables ....................... 39
  8.5 Generation of an array of variables .......................... 39
  8.6 Assignment to an array ....................................... 39
  8.7 Size of an array ............................................... 39
  8.8 bounds of an array ............................................ 40
  8.9 Display an array ............................................... 40
  8.10 Extract an element .......................................... 40
  8.11 Operations on array of series .............................. 44
    8.11.1 Matrix product ....................................... 44
    8.11.2 Matrix determinant ................................... 44
    8.11.3 Matrix inversion ..................................... 45
    8.11.4 Eigenvalues .......................................... 45
    8.11.5 Eigenvectors .......................................... 46
    8.11.6 Arithmetic ............................................ 47
  8.12 Conversion ................................................... 48
# Structures and OOP

<table>
<thead>
<tr>
<th>Section</th>
<th>Title</th>
</tr>
</thead>
<tbody>
<tr>
<td>9.1</td>
<td>Type declaration</td>
</tr>
<tr>
<td>9.2</td>
<td>Object identifier declaration</td>
</tr>
<tr>
<td>9.3</td>
<td>Access to member attributes</td>
</tr>
<tr>
<td>9.4</td>
<td>Display</td>
</tr>
<tr>
<td>9.5</td>
<td>Constructor macros</td>
</tr>
<tr>
<td>9.6</td>
<td>Member macros</td>
</tr>
<tr>
<td>9.6.1</td>
<td>Declaration</td>
</tr>
<tr>
<td>9.6.2</td>
<td>Execution</td>
</tr>
</tbody>
</table>

## Numerical vectors

<table>
<thead>
<tr>
<th>Section</th>
<th>Title</th>
</tr>
</thead>
<tbody>
<tr>
<td>10.1</td>
<td>Declaration</td>
</tr>
<tr>
<td>10.1.1</td>
<td>vnumR</td>
</tr>
<tr>
<td>10.1.2</td>
<td>vnumC</td>
</tr>
<tr>
<td>10.1.3</td>
<td>vnumQ</td>
</tr>
<tr>
<td>10.2</td>
<td>Initialization</td>
</tr>
<tr>
<td>10.3</td>
<td>Display</td>
</tr>
<tr>
<td>10.4</td>
<td>Size</td>
</tr>
<tr>
<td>10.5</td>
<td>Resize</td>
</tr>
<tr>
<td>10.6</td>
<td>Bounds</td>
</tr>
<tr>
<td>10.7</td>
<td>Data retrieval</td>
</tr>
<tr>
<td>10.7.1</td>
<td>select</td>
</tr>
<tr>
<td>10.7.2</td>
<td>Data retrieval</td>
</tr>
<tr>
<td>10.8</td>
<td>Input/Output</td>
</tr>
<tr>
<td>10.8.1</td>
<td>read</td>
</tr>
<tr>
<td>10.8.2</td>
<td>readappend</td>
</tr>
<tr>
<td>10.8.3</td>
<td>write</td>
</tr>
<tr>
<td>10.8.4</td>
<td>readbin</td>
</tr>
<tr>
<td>10.8.5</td>
<td>writebin</td>
</tr>
<tr>
<td>10.9</td>
<td>Input/Output low level</td>
</tr>
<tr>
<td>10.9.1</td>
<td>file_open</td>
</tr>
<tr>
<td>10.9.2</td>
<td>file_close</td>
</tr>
<tr>
<td>10.9.3</td>
<td>file_write</td>
</tr>
<tr>
<td>10.9.4</td>
<td>file_read</td>
</tr>
<tr>
<td>10.9.5</td>
<td>file_readappend</td>
</tr>
<tr>
<td>10.9.6</td>
<td>file_writemsg</td>
</tr>
<tr>
<td>10.9.7</td>
<td>file_writebin</td>
</tr>
<tr>
<td>10.10</td>
<td>Standard functions</td>
</tr>
<tr>
<td>10.10.1</td>
<td>minimum and maximum</td>
</tr>
<tr>
<td>10.10.2</td>
<td>addition and product</td>
</tr>
<tr>
<td>10.10.3</td>
<td>sort</td>
</tr>
<tr>
<td>10.10.4</td>
<td>transposevnum</td>
</tr>
<tr>
<td>10.10.5</td>
<td>standard math function</td>
</tr>
<tr>
<td>10.11</td>
<td>Conditions</td>
</tr>
<tr>
<td>10.12</td>
<td>Conversion</td>
</tr>
<tr>
<td>10.12.1</td>
<td>dimtovnumR</td>
</tr>
<tr>
<td>10.12.2</td>
<td>dimtovnumC</td>
</tr>
<tr>
<td>10.12.3</td>
<td>vnumtodim</td>
</tr>
<tr>
<td>10.12.4</td>
<td>str</td>
</tr>
</tbody>
</table>
11 Numerical matrices ................................................. 97
  11.1 Declaration ......................................................... 97
  11.1.1 matrixR ............................................................ 97
  11.1.2 matrixC ............................................................ 97
  11.2 Initialization ......................................................... 98
  11.3 Display ............................................................... 99
  11.4 Size .................................................................. 100
  11.5 Data retrieval ....................................................... 100
  11.6 Input/Output .......................................................... 101
    11.6.1 sauve_c ............................................................. 101
    11.6.2 sauve_fortran ..................................................... 101
    11.6.3 sauve_tex ............................................................ 101
    11.6.4 sauve_ml ............................................................. 102
    11.6.5 write ................................................................. 102
    11.6.6 writebin ............................................................. 102
  11.7 Input/Output low level ............................................. 102
  11.8 Standard math function ........................................ 102
    11.8.1 Matrix product .................................................. 103
    11.8.2 Kronecker product .............................................. 104
    11.8.3 Matrix determinant ............................................ 104
    11.8.4 Matrix inversion ............................................... 104
    11.8.5 Matrix trace ..................................................... 105
    11.8.6 Transpose ........................................................ 105
    11.8.7 Identity matrix ................................................ 106
    11.8.8 Eigenvalues ..................................................... 106
    11.8.9 Eigenvectors .................................................... 106
    11.8.10 Arithmetic .................................................... 107
  11.9 Conditions ........................................................... 108
  11.10 Conversion .......................................................... 109
12 Graphics ............................................................... 111
  12.1 plot ................................................................. 111
  12.2 replot .............................................................. 112
  12.3 plotf ................................................................. 114
  12.4 plotps ............................................................... 114
  12.5 plotps_end .......................................................... 114
  12.6 plotreset ............................................................ 115
  12.7 gnuplot ............................................................... 115
  12.8 grace ................................................................. 116
13 Communications ..................................................... 119
  13.1 Maple ................................................................. 119
    13.1.1 maple_put ....................................................... 119
    13.1.2 maple_get ...................................................... 120
    13.1.3 maple ............................................................. 120
  13.2 Mathematica ........................................................ 121
    13.2.1 mathematica_put ............................................... 121
    13.2.2 mathematica_get .............................................. 122
    13.2.3 mathematica .................................................... 123
  13.3 Communications with other computer algebra systems .... 124
    13.3.1 SCSCP server ................................................ 124
    13.3.1.1 sscp_disable_cy ........................................ 125
17 Numerical computations ........................................... 177

17.1 Frequency analysis ............................................... 177
  17.1.1 naf ............................................................... 177
  17.1.2 naftab ............................................................ 178
  17.1.3 freqa ............................................................. 179
  17.1.4 freqareson ...................................................... 181
  17.1.5 sertrig ............................................................ 181

17.2 Fourier Transform .................................................. 182

17.3 Inverse Fourier Transform ........................................ 183

17.4 Numerical integration of series or macro ..................... 183

17.5 Numerical integration of external function ................... 188

17.6 Integral ............................................................... 192

17.7 Numerical iteration .................................................. 193

17.8 Interpolation .......................................................... 194

17.9 Coordinate system conversions ................................... 195
  17.9.1 ell_to_xyz ......................................................... 195
  17.9.2 xyz_to_ell ....................................................... 196

16.1.2 Least Squares ..................................................... 155

16.1.3 Factorizations .................................................... 164

16.1.4 Singular Value Decomposition ................................. 168

16.1.5 Eigenvalues and Eigenvectors ................................ 170

lapack_zposv .......................................................... 154
lapack_zpbsv .......................................................... 154
lapack_zptsv .......................................................... 155

16.1.2 Least Squares ..................................................... 155

lapack_dgelss .......................................................... 158
lapack_dggglm .......................................................... 160
lapack_zgels .......................................................... 161
lapack_zgelss .......................................................... 162
lapack_zggglm .......................................................... 163

16.1.3 Factorizations .................................................... 164

lapack_dgeqrf .......................................................... 164
lapack_dgeqrlf .......................................................... 165
lapack_dpotrf .......................................................... 166
lapack_zgeqrf .......................................................... 166
lapack_zgeqrlf .......................................................... 167
lapack_zpotrf .......................................................... 167

16.1.4 Singular Value Decomposition ................................. 168

lapack_dgesvd .......................................................... 168
lapack_dgesdd .......................................................... 168
lapack_zgesvd .......................................................... 169
lapack_zgesdd .......................................................... 169

16.1.5 Eigenvalues and Eigenvectors ................................ 170

lapack_dsyev .......................................................... 170
lapack_dsbev .......................................................... 170
lapack_dstev .......................................................... 171
lapack_dgeev .......................................................... 172
lapack_dsyev .......................................................... 172
lapack_dgeev .......................................................... 173
lapack_dgges .......................................................... 174
lapack_zheev .......................................................... 174
lapack_zlbev .......................................................... 175
lapack_zgeev .......................................................... 175
lapack_zhegv .......................................................... 176
18 Utilities ................................................. 197
  18.1 help ..................................................... 197
  18.2 exit ...................................................... 197
  18.3 cls ....................................................... 197
  18.4 pause ..................................................... 197
  18.5 msg ....................................................... 198
  18.6 error ...................................................... 199
  18.7 try ......................................................... 199
  18.8 delete .................................................... 200
  18.9 reset ..................................................... 200
  18.10 include .................................................. 200
  18.11 @@ ........................................................ 201
  18.12 vartrip .................................................. 201
  18.13 bilan ..................................................... 201
  18.14 stat ....................................................... 202
  18.15 save_env ............................................... 204
  18.16 rest_env ............................................... 204
  18.17 random ................................................. 204
  18.18 nameof .................................................. 205
  18.19 typeof ................................................... 205
  18.20 filefullname ........................................... 205
  18.21 Execution time ........................................ 206
    18.21.1 time_s ............................................. 206
    18.21.2 time_l ............................................. 206
    18.21.3 time_t ............................................. 206
  18.22 Call the shell ......................................... 207

19 References ............................................. 209

Appendix A  Supported OpenMath Content Dictionaries ... 211
Appendix B  Index of the global variables ...................... 213
Appendix C  Index of the procedures ............................ 215
Appendix D  Full index ........................................ 219