

# A FORTRAN 90 numerical library

Alberto Ramos. Madrid, November 2006.

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# Generalities

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This is the documentation of a total of thirteen **FORTRAN 90** modules with different utilities. This code is well documented, and can be useful for several people, although the idea is *not* to produce fast, high performance code, but to have nice data structures and **INTERFACE** definitions so that complex problems can be solved fast, writing only a couple of lines of code.

The code of all these modules is *free software*, this means that you can redistribute and/or modify all the code under the terms of the GNU General Public License<sup>1</sup> as published by the Free Software Foundation; either version 2 of the License, or (at your option) any later version. Note that the code is distributed in the hope that it will be useful, but **without any warranty; without even the implied warranty of merchantability or fitness for a particular purpose**. See the GNU General Public License for more details.

The code has been written using standard **FORTRAN 90**, this means that it should run on any machine and with any compiler. In particular the code of all these modules has been compiled using GNU **gfortran**, INTEL **ifort** and DIGITAL **f90** for PC.

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The source code of all the modules as well as the last version of this document should always be available (in it's last version) at:

`http://lattice.ft.uam.es/perpag/alberto/codigo-en.php`

there is also a **sourceforge.net** project, where the last version of both the source code and the documentation should be available:

`http://sourceforge.net/projects/afnl`

Enjoy programming.

## Installation

To install this library in a Unix/Linux environment, simply edit the **Makefile** file, and set the **F90** and **F90OPT** variables to whatever your compiler and your favourite optimisation flags are. After running **make** you should obtain a file called **libf90.a**, and probably (that depends on

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<sup>1</sup><http://www.gnu.org/copyleft/gpl.html>

the particular compiler) some `.mod` files. Copy the `libf90.a` library and the `.mod` files to any place you like, and compile and link your program to that files. With GNU `gfortran` this is done using the flags `-I<path> -L<path> -lf90`, where `<path>` has to be substituted by the path you have chosen.

In other environments, you should ask the local guru/administrator about how to generate a library. In particular in a Windows environment the best option is to repartition you hard drive, eliminate Windows and install any Unix like free operating system, like Linux or FreeBSD.

# One

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## MODULE NumTypes

---

This is the documentation of the `MODULE NumTypes`, that contains the definition of Single Precision, and Double Precision data. All the other numerical modules use this data type definitions.

### 1.1 Description

The `MODULE NumTypes` provides the definition of the Single Precision and Double Precision real and complex data in a portable way. When we want to define a single precision real we *will* do it with a statement like `Real (kind=DP)`, instead of `Real (kind=4)`. What we mean with DP is defined in this module. The different data types are:

**SP:** Single precision real.

**DP:** Double precision real.

**SPC:** Single precision complex.

**DPC:** Double precision complex.

To make all the code as portable as possible, all the data definitions should make use of this module.

### 1.2 Examples

Here we will define `A` as a single precision real, `D` as a double precision real, `Ac` as a single precision complex, and `Dc` as a double precision complex.

Listing 1.1: Definition of data types.

```
1  Program Types_of_Data
2  USE NumTypes
3
4
5  Real (kind=SP) :: A
6  Real (kind=DP) :: D
```

```
7      Complex (kind=SPC)  :: Ac
      Complex (kind=DPC)  :: Dc
9
      Write(*,*) Kind(A), Kind(Aa)
11
13     End Program Types_of_Data
```

# Two

---

## MODULE Constants

---

This is the documentation of the `MODULE Constants`, that contains the definition of the most used mathematical constants. This module uses numerical types defined in the `MODULE NumTypes`.

### 2.1 Name conventions

All the real simple precision constants ends with `_SP`, the real double precision constants with `_DP`, the complex simple precision with `_SPC` and the complex double precision with `_DPC`.

If a there exist a real or complex constant of simple precison defined, then it exist other with the same name (except for the suffix) of double precision and viceversa.

### 2.2 $\pi$ -related constants

#### 2.2.1 Real

The complex  $\pi$ -related defined in this module and its values can be seen in the table (2.1)

SP Name	DP Name	Value
PI_SP	PI_DP	$\pi$
TWOPI_SP	TWOPI_DP	$2\pi$
HALFPI_SP	HALFPI_DP	$\frac{\pi}{2}$

Table 2.1:  $\pi$ -related real constants defined in the `MODULE constants`.

#### 2.2.2 Complex

The complex  $\pi$ -related defined in this module and its values can be seen in the table (2.2)

### 2.3 Square roots and log related constants

We have only real constants defined here. We can see a list of names-vlues in the table (2.3)

SPC Name	DPC Name	Value
UNITIMAG_SPC	UNITIMAG_DPC	$\iota$
PI_IMAG_SPC	PI_IMAG_DPC	$\pi\iota$
TWOPI_IMAG_SPC	TWOPI_IMAG_DPC	$2\pi\iota$
HALFPI_IMAG_SPC	HALFPI_IMAG_DPC	$\frac{\pi}{2}\iota$

Table 2.2:  $\pi$ -related complex constants defined in the MODULE constants.

SP Name	DP Name	Value
SR2_SP	SR2_DP	$\sqrt{2}$
SR3_SP	SR3_DP	$\sqrt{3}$
SRe_SP	SRe_DP	$\sqrt{e}$
SRpi_SP	SRpi_DP	$\sqrt{\pi}$
LG102_SP	LG102_DP	$\log_{10} 2$
LG103_SP	LG103_DP	$\log_{10} 3$
LG10e_SP	LG10e_DP	$\log_{10} e$
LG10pi_SP	LG10pi_DP	$\log_{10} \pi$
LGe2_SP	LGe2_DP	$\log_e 2$
LGe3_SP	LGe3_DP	$\log_e 3$
LGe10_SP	LGe10_DP	$\log_e 10$

Table 2.3: Square roots and log related constants defined in the MODULE constants.

## 2.4 Other mathematical constants

In this section we have only the Euler  $\gamma$  constant. We can see the name-value pair in the table (2.4)

SP Name	DP Name	Value
GEULER_SP	GEULER_DP	$\gamma(= 0.5772\dots)$

Table 2.4: Other mathematical constants defined in the MODULE constants.

# Three

---

## MODULE Error

---

This is the documentation of the `MODULE Error`, a set of FORTRAN 90 routines that allow to write errors.

### 3.1 Defined variables

#### 3.1.1 stderr

##### Description

This variable has the unit number of standard error.

##### Examples

Listing 3.1: Standard error unit.

```
1 Program Test
  USE Error
3
  Write(stderr,*)''This is printed in standard error.''
5
  Stop
7 End Program Test
```

### 3.2 Subroutine perror([routine], msg)

#### 3.2.1 Description

Prints the error message `msg` in standard error. If the optional argument `routine` is given, it is used as the routine where the program has crashed.

#### 3.2.2 Arguments

**routine:** Character string with arbitrary length. It should be the routine or program name where the error has occurred. It is an optional argument.

**msg:** Character string with arbitrary length. It should be the message that you want to print.

### 3.2.3 Examples

Listing 3.2: Print error message.

```

1 Program Test
  USE Error
3
  Integer :: N1, N2
5
  Write(*,*) 'Two integer numbers: '
7  Read(*,*) N1, N2
9
  If (N2 == 0) Then
    CALL Perror('Test', 'Division by zero. See the product: ')
11  Write(*,*) N1*N2
  Else
13  Write(*,*) N1/N2
  End If
15
  Stop
17 End Program Test

```

## 3.3 Subroutine abort([routine], msg)

### 3.3.1 Description

Prints the error message **msg** in standard error, and stops the program. If the optional argument **routine** is given, it is used as the routine where the program has crashed.

### 3.3.2 Arguments

**routine:** Character string with arbitrary length. It should be the routine or program name where the error has occurred. It is an optional argument.

**msg:** Character string with arbitrary length. It should be the message that you want to print.

### 3.3.3 Examples

Listing 3.3: Print error message and stop a program.

```

1 Program Test
  USE Error
3
  Integer :: N1, N2
5
  Write(*,*) 'Two integer numbers: '
7  Read(*,*) N1, N2
9
  If (N2 == 0) Then

```



```
11      CALL abort('Test', 'Division by zero')
      Else
13      Write(*,*)N1/N2
      End If

15      Stop
End Program Test
```



# Four

---

## MODULE Integration

---

This is the documentation of the `MODULE Integration`, a set of `FORTRAN 90` routines that performs numerical integration and solves the initial value problem for a specified system of first-order ordinary differential equations. This module make use of the `MODULE NumTypes`, so please read the documentation of this module *before* reading this.

### 4.1 Function Trapecio(a, b, Func, [Tol])

#### 4.1.1 Description

Calculates the integral of the function `Func` between `a` and `b` with precision `Tol` (optional) using the trapezoid rule.

#### 4.1.2 Arguments

**a, b:** Real single or double precision. The limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface
```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is `Tol = 0.01`.

### 4.1.3 Output

If the arguments are real of single (double) precision, the result will also be a real of single (double) precision. The value of the integral.

### 4.1.4 Examples

Listing 4.1: Example of integration of a function using Trapecio.

```

Program Test
2  USE NumTypes
  USE Integration
4
  Real (kind=DP) :: Tol
6
  Interface
8    Function Fint(X)
      USE NumTypes
10
      Real (kind=DP), Intent (in) :: X
12      Real (kind=DP) :: Fint
    End Function Fint
14 End Interface

16 Tol = 1.0E-6_DP
  Write(*,*) 'Integral of x**2 between 0 and 1:'
18  Write(*,*) Trapecio(0.0_DP, 1.0_DP, Fint, Tol)

20  Stop
End Program Test
22
! *****
24 ! *
Function Fint(X)
26 ! *
! *****
28
  USE NumTypes
30
  Real (kind=DP), Intent (in) :: X
32  Real (kind=DP) :: Fint

34  Fint = X**2

36  Return
End Function Fint

```

## 4.2 Function Simpson(a, b, Func, [Tol])

### 4.2.1 Description

Calculates the integral of the function `Func` between `a` and `b` with precision `Tol` (optional) using the Simpson's rule.

In general this routine is better than `Trapeccio`.

### 4.2.2 Arguments

**a, b:** Real single or double precision. The limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface
```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is `Tol = 0.01`.

### 4.2.3 Output

If the arguments are reals of single (double) precision, the result will also be a real of single (double) precision. The value of the integral.

### 4.2.4 Examples

Listing 4.2: Exmaple of integration of a function using Simpson.

```
1 Program Test
  USE NumTypes
3  USE Integration

5  Real (kind=DP) :: Tol

7  Interface
    Function Fint(X)
9    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
13 End Function Fint
```

```

15      End Interface
16
17      Tol = 1.0E-6_DP
18      Write(*,*) 'Integral of x**2 between 0 and 1:'
19      Write(*,*) Simpson(0.0_DP, 1.0_DP, Fint, Tol)
20
21      Stop
22  End Program Test
23
24  ! *****
25  ! *
26  Function Fint(X)
27  ! *****
28
29      USE NumTypes
30
31      Real (kind=DP), Intent (in) :: X
32      Real (kind=DP) :: Fint
33
34      Fint = X**2
35
36      Return
37  End Function Fint

```

### 4.3 Function TrapecioAb(a, b, Func, [Tol])

#### 4.3.1 Description

Calculates the integral of the function `Func` between `a` and `b` with precision `Tol` (optional) using the open trapezoid rule.

#### 4.3.2 Arguments

**a, b:** Single (SP) or double (DP) precision. They are the limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```

Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface

```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result.  
It is an optional parameter, and the default is  $Tol = 0.01$ .

### 4.3.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision.  
The value of the integral.

### 4.3.4 Examples

Listing 4.3: Integrating a function using the open trapezoid rule.

```

1  Program Test
   USE NumTypes
3   USE Integration

5   Real (kind=DP) :: Tol

7   Interface
   Function Fint(X)
9       USE NumTypes

11      Real (kind=DP), Intent (in) :: X
       Real (kind=DP) :: Fint
13  End Function Fint
End Interface

15
Tol = 1.0E-6_DP
17 Write(*,*) 'Integral of x**2 between 0 and 1:'
Write(*,*) TrapecioAb(0.0_DP, 1.0_DP, Fint, Tol)
19
   Stop
21 End Program Test

23 ! *****
24 ! *
25 Function Fint(X)
26 ! *
27 ! *****

29 USE NumTypes

31 Real (kind=DP), Intent (in) :: X
   Real (kind=DP) :: Fint
33
   Fint = X**2
35
   Return
37 End Function Fint

```

## 4.4 Function SimpsonAb(a, b, Func, [Tol])

### 4.4.1 Description

Calculates the integral of the function `Func` between `a` and `b` with precision `Tol` (optional) using the open Simpson's rule.

In general better than `TrapezioAb`

### 4.4.2 Arguments

**a, b:** Single (SP) or double (DP) precision. They are the limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface
```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is `Tol = 0.01`.

### 4.4.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision. The value of the integral.

### 4.4.4 Examples

Listing 4.4: Exmaple of integration using the open Simpson rule.

```
1 Program Test
  USE NumTypes
3  USE Integration

5  Real (kind=DP) :: Tol

7  Interface
    Function Fint(X)
9    USE NumTypes

11    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
13  End Function Fint
```



```

15      End Interface
17      Tol = 1.0E-6_DP
18      Write(*,*) 'Integral of x**2 between 0 and 1:'
19      Write(*,*) SimpsonAb(0.0_DP, 1.0_DP, Fint, Tol)
21      Stop
22  End Program Test
23
24  ! *****
25  ! *
26  Function Fint(X)
27  ! *
28  ! *****
29
30  USE NumTypes
31
32  Real (kind=DP), Intent (in) :: X
33  Real (kind=DP) :: Fint
34
35  Fint = X**2
36
37  Return
38 End Function Fint

```

## 4.5 Function SimpsonInfUp(a, Func, [Tol])

### 4.5.1 Description

Calculates the integral of the function **Func** between **a** and  $\infty$  with precision **Tol** (optional) using the Simpson rule and a change of variables.

### 4.5.2 Arguments

**a**: Single (SP) or double (DP) precision. They are the limits of the integral.

**Func**: The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```

Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface

```

This routine does not check if the integral exist, so the function must obviously decay fast for large  $x$  to obtain a finite value.

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is  $\text{Tol} = 0.01$ .

### 4.5.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision. The value of the integral.

### 4.5.4 examples

Listing 4.5: Integration of a function between 0 and  $\infty$ .

```

1  Program Test
   USE NumTypes
3   USE Integration

5   Real (kind=DP) :: Tol

7   Interface
   Function Fint(X)
9       USE NumTypes

11      Real (kind=DP), Intent (in) :: X
       Real (kind=DP) :: Fint
13  End Function Fint
End Interface

15  Tol = 1.0E-6_DP
17  Write(*,*) 'Integral of e**(-x**2) between 0 and infinity:'
19  Write(*,*) SimpsonInfUp(0.0_DP, Fint, Tol)

21  Stop
End Program Test

23  ! *****
25  ! *
Function Fint(X)
27  ! *****

29  USE NumTypes

31  Real (kind=DP), Intent (in) :: X
   Real (kind=DP) :: Fint

33  Fint = exp(-X**2)

35  Return
37  End Function Fint

```

## 4.6 Function SimpsonInfDw(a, Func, [Tol])

### 4.6.1 Description

Calculates the integral of the function `Func` between  $-\infty$  and `a` with precision `Tol` (optional) using the Simpson rule and a change of variables.

### 4.6.2 Arguments

**a:** Single (SP) or double (DP) precision. They are the limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface
```

This routine does not check if the integral exist, so the function must obviously decay fast for large  $-x$  to obtain a finite value.

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is `Tol = 0.01`.

### 4.6.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision. The value of the integral.

### 4.6.4 examples

Listing 4.6: Integrating a function between  $-\infty$  and 0.

```
1 Program Test
  USE NumTypes
  USE Integration
3
5 Real (kind=DP) :: Tol
7
8 Interface
  Function Fint(X)
  USE NumTypes
9
11 Real (kind=DP), Intent (in) :: X
```

```

13      Real (kind=DP) :: Fint
      End Function Fint
      End Interface

15
      Tol = 1.0E-6_DP
17      Write(*,*) 'Integral of e**(-x**2) between -infinity and 0:'
      Write(*,*) SimpsonInfDw(0.0_DP, Fint, Tol)

19
      Stop
21 End Program Test

23 ! *****
24 ! *
25 Function Fint(X)
26 ! *
27 ! *****

29 USE NumTypes

31 Real (kind=DP), Intent (in) :: X
32 Real (kind=DP) :: Fint
33
34 Fint = exp(-X**2)
35
36 Return
37 End Function Fint

```

## 4.7 Function SimpsonSingUp(a, b, Func, [Tol], gamma)

### 4.7.1 Description

Calculates the integral of the function **Func** between **a** and **b** with precision **Tol** (optional) using the Simpson's rule. The function may have an integrable singularity of the type:

$$f(x+b) \approx \frac{c}{(x-b)^\gamma} + \dots$$

with  $0 < \gamma < 1$ .

### 4.7.2 Arguments

**a, b:** Single (SP) or double (DP) precision. They are the limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```

Interface
  Function Fint(X)
    USE NumTypes

```

```

        Real (kind=DP), Intent (in) :: X
        Real (kind=DP) :: Fint
    End Function Fint
End Interface

```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is  $Tol = 0.01$ .

**gamma:** The “degree of divergence” of the function in  $x \approx b$ .

### 4.7.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision. The value of the integral.

### 4.7.4 Examples

Listing 4.7: Integrating functions with singularities in the upper limit.

```

1  Program Test
   USE NumTypes
3  USE Integration

5  Real (kind=DP) :: Tol

7  Interface
   Function Fint(X)
9      USE NumTypes

11     Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
13 End Function Fint
End Interface

15 Tol = 1.0E-6_DP
17 Write(*,*) 'Integral of 1/sqrt(-x) between -1 and 0:'
18 Write(*,*) SimpsonSingUp(-1.0_DP, 0.0_DP, Fint, Tol, 0.5_DP)
19
20 Stop
21 End Program Test

23 ! *****
24 ! *
25 Function Fint(X)
26 ! *
27 ! *****

29 USE NumTypes

31 Real (kind=DP), Intent (in) :: X
32 Real (kind=DP) :: Fint
33

```

```

35   Fint = Sqrt(-X)
37   Return
End Function Fint

```

## 4.8 Function SimpsonSingDw(a, b, Func, [Tol], gamma)

### 4.8.1 Description

Calculates the integral of the function **Func** between **a** and **b** with precision **Tol** (optional) using the Simpson's rule. The function may have an integrable singularity of the type:

$$f(x+a) \approx \frac{c}{(x-a)^\gamma} + \dots$$

with  $0 < \gamma < 1$ .

### 4.8.2 Arguments

**a, b:** Single (SP) or double (DP) precision. They are the limits of the integral.

**Func:** The function to be integrated. It must be a function of only one argument of the same type as the function itself. If it is an external function an interface block like the following should be declared:

```

Interface
  Function Fint(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X
    Real (kind=DP) :: Fint
  End Function Fint
End Interface

```

**Tol:** Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter, and the default is **Tol** = 0.01.

**gamma:** The “degree of divergence” of the function in  $x \approx a$ .

### 4.8.3 Output

If the arguments are single (double) precision, the result will also be of single (double) precision. The value of the integral.

### 4.8.4 Examples

Listing 4.8: Integrating functions with singularities in the lower limit.

```

1  Program Test
   USE NumTypes
3   USE Integration

5   Real (kind=DP) :: Tol

7   Interface
   Function Fint(X)
9       USE NumTypes

11      Real (kind=DP), Intent (in) :: X
       Real (kind=DP) :: Fint
13  End Function Fint
End Interface

15  Tol = 1.0E-6_DP
17  Write(*,*) 'Integral of 1/sqrt(x) between 0 and 1:'
19  Write(*,*) SimpsonSingDw(0.0_DP, 1.0_DP, Fint, Tol, 0.5_DP)

21  Stop
End Program Test

23  ! *****
25  ! *
Function Fint(X)
27  ! *
   ! *****

29  USE NumTypes

31  Real (kind=DP), Intent (in) :: X
   Real (kind=DP) :: Fint

33  Fint = Sqrt(X)

35  Return
37 End Function Fint

```

## 4.9 Function Euler(Init, Xo, Xfin, Feuler, [Tol])

### 4.9.1 Description

Integrate the first order set of ODE defined by the function **Feuler**, with initial conditions given by the vector **Init** in **Xo**, until **Xfin**, with a precision given by **Tol** (optional).

A set of first order ODE's is given by the first derivatives of the variables involved:

$$\frac{dy_i(x)}{dx} = f_i(y_j, x)$$

and the initial conditions:

$$y_i(x_0)$$

After the integration we get:

$$y_i(x_{\text{fin}})$$

So to define a set of first order ODE's we need the value of the derivative of the variable  $i$  in the point  $x$  (this is done by **Feuler**), a vector of initial conditions (**Init**) and the point where this initial conditions are defined (**Xo**), and finally the point where we want the solution (**Xfin**)

### 4.9.2 Arguments

**Init(:)**: Single (SP) or double (DP) precision vector of one dimension with the initial conditions.

**Xo**: Single (SP) or double (DP) precision. The point where the initial conditions are defined.

**Xfin**: Single (SP) or double (DP) precision. The point where we want the value of the functions.

**Feuler**: The function that defines the set of first order ODE's. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Feuler(X, Y) Result (Func)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X, Y(:)
    Real (kind=DP) :: Func(Size(Y))
  End Function Feuler
End Interface
```

The function must return a vector with the values of the first derivatives of the functions  $y_i(x)$  in the point  $X$ .

**Tol**: Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter.

### 4.9.3 Output

Real single or double precision (same as input) one dimensional array. The array contains the values of the functions  $y_i$  in the point **Xfin**.

### 4.9.4 Examples

This example below will integrate the set of first order ODE's defined by the equations:

$$\frac{dy_1(x)}{dx} = y_2(x); \quad \frac{dy_2(x)}{dx} = -y_1(x)$$

whose solution is:

$$y_1(x) = A \cos(x) + B \sin(x)$$



With the initial conditions  $y_1(0) = 0$ ;  $y_2(0) = 1$ , the solution is:

$$y_1(x) = \sin(x); \quad y_2(x) = \cos(x)$$

so if we plot  $y_1(1)$  and  $y_2(1)$  we will obtain the values  $\sin(1)$  and  $\cos(1)$ . In the following example, we will compare the result of integrating the differential equations with the exact values.

Listing 4.9: Integrating differential equations with Euler.

```

1 Program Test
  USE NumTypes
3  USE Integration

5  Real (kind=DP) :: Tol, In(2)

7  Interface
  Function Feuler(X, Y) Result (Func)
9    USE NumTypes

11     Real (kind=DP), Intent (in) :: X, Y(:)
    Real (kind=DP) :: Func(Size(Y))
13  End Function Feuler
End Interface

15

17  Tol = 1.0E-2_DP
  In(1) = 0.0_DP
19  In(2) = 1.0_DP
  Write(*,*) 'Values of sin(1) and cos(1): '
21  Write(*,*) Euler(In, 0.0_DP, 1.0_DP, Feuler, Tol)
  Write(*,*) Sin(1.0_DP), Cos(1.0_DP)
23

  Stop
25 End Program Test

27 ! *****
! *
29 Function FEuler(X, Y) Result (Func)
! *
31 ! *****

33     Real (kind=8), Intent (in) :: X, Y(:)
    Real (kind=8) :: Func(Size(Y))

35     Func(1) = Y(2)
37     Func(2) = -Y(1)

39     Return
End Function FEuler

```

## 4.10 Function Rgnkta(Init, Xo, Xfin, Feuler, [Tol])

### 4.10.1 Description

Integrate the first order set of ODE defined by the function **Feuler**, with initial conditions given by the vector **Init** in **Xo**, until **Xfin**, with a precision given by **Tol** (optional). This method uses a Runge-Kutta algorithm and is much more exact than the previous **Euler** function.

A set of first order ODE's is given by the first derivatives of the variables involved:

$$\frac{dy_i(x)}{dx} = f_i(y_j, x)$$

and the initial conditions:

$$y_i(x_0)$$

After the integration we get:

$$y_i(x_{\text{fin}})$$

So to define a set of first order ODE's we need the value of the derivative of the variable  $i$  in the point  $x$  (this is done by **Feuler**), a vector of initial conditions (**Init**) and the point where this initial conditions are defined (**Xo**), and finally the point where we want the solution (**Xfin**)

### 4.10.2 Arguments

**Init(:)**: Single (SP) or double (DP) precision vector of one dimension with the initial conditions.

**Xo**: Single (SP) or double (DP) precision. The point where the initial conditions are defined.

**Xfin**: Single (SP) or double (DP) precision. The point where we want the value of the functions.

**Feuler**: The function that defines the set of first order ODE's. If it is an external function an interface block like the following should be declared:

```
Interface
  Function Feuler(X, Y) Result (Func)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X, Y(:)
    Real (kind=DP) :: Func(Size(Y))
  End Function Feuler
End Interface
```

The function is the same as in the previous function.

**Tol**: Single (SP) or double (DP) precision. An estimation of the desired accuracy of the result. It is an optional parameter.

### 4.10.3 Output

Real single or double precision (same as input) one dimensional array. The array contains the values of the functions  $y_i$  in the point **Xfin**.

#### 4.10.4 Examples

This example below will integrate the set of first order ODE's defined by the equations:

$$\frac{dy_1(x)}{dx} = y_2(x); \quad \frac{dy_2(x)}{dx} = -y_1(x)$$

whose solution is:

$$y_1(x) = A \cos(x) + B \sin(x)$$

With the initial conditions  $y_1(0) = 0; y_2(0) = 1$ , we have:

$$y_1(x) = \sin(x); \quad y_2(x) = \cos(x)$$

so if we plot  $y_1(1)$  and  $y_2(1)$  we will obtain the values  $\sin(1)$  and  $y_2(1)$ . In the following example, we will compare the values obtained with **Euler**, with **Rgnkta** and the exact ones.

Listing 4.10: Integrating differential equations with the Runge-Kutta method

```

Program Test
2  USE NumTypes
  USE Integration
4
  Real (kind=DP) :: Tol, In(2)
6
  Interface
8    Function Feuler(X, Y) Result (Func)
      USE NumTypes
10
      Real (kind=DP), Intent (in) :: X, Y(:)
12    Real (kind=DP) :: Func(Size(Y))
      End Function Feuler
14 End Interface

16
  Tol = 1.0E-3_DP
18  In(1) = 0.0_DP
  In(2) = 1.0_DP
20  Write(*,*) 'Values of sin(1) and cos(1): '
  Write(*,*) ' Euler: '
22  Write(*,*) Euler(In, 0.0_DP, 1.0_DP, Feuler, Tol)
  Write(*,*) ' Runge-Kutta: '
24  Write(*,*) Rgnkta(In, 0.0_DP, 1.0_DP, Feuler, Tol)
  Write(*,*) ' Exact: '
26  Write(*,*) Sin(1.0_DP), Cos(1.0_DP)

28  Stop
End Program Test
30
! *****
32 ! *
  Function FEuler(X, Y) Result (Func)
34 ! *
! *****
36
```

```
38      Real (kind=8), Intent (in) :: X, Y(:)
      Real (kind=8) :: Func(Size(Y))

40      Func(1) = Y(2)
      Func(2) = -Y(1)

42      Return

44      End Function FEuler
```

# Five

---

## MODULE Optimization

---

This is the documentation of the MODULE `Optimization`, a set of routines to Optimise (maximise or minimise) functions of one or several variables. Lot of work is needed to improve this module (conjugate gradient, simplex, etc...).

### 5.1 Function `Step(X, FStep[, Tol])`

#### 5.1.1 Description

The function `Step(X, FStep, Tol)` returns the position of the minimum of the Function `Fstep` with an optional precision `Tol`.

#### 5.1.2 Arguments

**X:** Real single or double precision. An initial guess of the position of the minimum.

**Fstep:** The function that we want to minimise. It can be a function of one or several variables.

In the case of one variable functions an interface like the following should be declared

```
Interface
  Function Fstep(Xo)
    USE NumTypes

    Real (kind=DP), Intent (in) :: Xo
    Real (kind=DP) :: Fstep
  End Function Fstep
End Interface
```

In the case of a function of several variables, the interface block should be like the following

```
Interface
  Function Fstep(Xo)
    USE NumTypes
```

```

        Real (kind=DP), Intent (in) :: Xo(:)
        Real (kind=DP) :: Fstep
    End Function Fstep
End Interface

```

Tol: Real single or double precision. As estimation of the precision of the result. The default value is  $10^{-3}$ .

### 5.1.3 Output

Real Single or double precision (same as the output). The position of a minimum of Fstep.

### 5.1.4 Example

Listing 5.1: Minimising a function.

```

Program TestMin
2
  USE NumTypes
4  USE Optimization

6  Integer, Parameter :: Ndim = 4
  Real (kind=DP) :: XoM(Ndim), Xmin(Ndim)
8
  Interface
10    Function FstepM(Xo)
      USE NumTypes
12
      Real (kind=DP), Intent (in) :: Xo(:)
14    Real (kind=DP) :: FstepM
    End Function FstepM
16  End Interface

18
  ! Initial guess of the position of the minimum
20  XoM(1) = 1.373_DP
  XoM(2) = 1.373_DP
22  XoM(3) = 1.373_DP
  XoM(4) = 1.373_DP
24

26  Write(*,*)'Initial Position: '
  Do I = 1, Ndim
28    Write(*, '(1A,1I4,1A,1ES33.25)') 'Variable ', I, " : ", XoM(I)
  End Do
30

  Xmin = Step(XoM, FstepM, 1.0E-7_DP)
32  Write(*,*)
  Write(*,*)'Position of the minimum: '
34  Do I = 1, Ndim
    Write(*, '(1A,1I4,1A,1ES33.25)') 'Variable ', I, " : ", Xmin(I)

```

```

36      End Do
38
39      Stop
40  End Program TestMin
41
42  ! *****
43  ! *
44  Function FstepM(Xo)
45  ! *
46  ! *****
47
48  USE NumTypes
49
50  Real (kind=DP), Intent (in) :: Xo(:)
51  Real (kind=DP) :: FstepM
52
53  FstepM = (Xo(1)-1.0_DP)**2 + &
54           & (Xo(2)-2.0_DP)**2 + &
55           & (Xo(3)+3.0_DP)**4 + &
56           & (Xo(4)-4.0_DP)**8
57
58  Return
59 End Function FstepM

```

## 5.2 Function MaxPosition(FVal, IpX, IpY)

### 5.2.1 Description

Given a two dimensional array of values `FVal(:, :)`, the function `MaxPosition(FVal, IpX, IpY)` returns the number of local maxima of `FVal(:, :)` and its positions in the one dimensional arrays `IpX(:)` and `IpY(:)`.

### 5.2.2 Arguments

`FVal(:, :)`: Real single or double precision. The values of a function in a two dimensional grid of points.

`IpX(:)`: Integer. A one dimensional array that contains the value of  $X$  for the positions of the maxima.

`IpY(:)`: Integer. A one dimensional array that contains the value of  $Y$  for the positions of the maxima.

### 5.2.3 Output

Integer. The number of local maxima of the input. `FVal(:, :)`.

### 5.2.4 Example

Listing 5.2: Example of the usage of the routine MaxPosition.

```

1  Program MaxLoc
2
3      USE NumTypes
4      USE Constants
5      USE Optimization
6      USE Error
7
8      IMPLICIT NONE
9
10     Integer :: I, J, IsX, IsY, Nmax
11     Character (len=200) :: Filename
12     Real (kind=DP) :: DnullX, DnullY
13     Real (kind=DP), Allocatable :: F(:, :)
14     Integer :: IpX(10), IpY(10)
15
16     Write(stderr, *) "SizeX, SizeY, Filename"
17     Read(*, *) IsX, IsY, Filename
18
19     Allocate(F(IsX, IsY))
20     Open (Unit=666, File=Trim(Filename), Action="READ")
21     Do I = 1, IsX
22         Do J = 1, IsY
23             Read(666, *) DnullX, DnullY, F(I, J)
24             Write(stderr, *) DnullX, DnullY, F(I, J)
25         End Do
26     End Do
27     Close(666)
28
29     Nmax = MaxPosition(F, IpX, IpY)
30     Write(*, *) "# Number of maxima: ", Nmax
31     Write(*, *) "# Positions of the maxima: "
32     Do I = 1, Nmax
33         Write(*, *) IpX(I), IpY(I)
34     End Do
35
36     Stop
37 End Program MaxLoc

```



# Six

---

## MODULE Linear

---

This is the documentation of the `MODULE Linear`, a set of `FORTRAN 90` routines to solve linear systems of equations. This module make use of the `MODULE NumTypes`, and `MODULE Error` so please read the documentation of these modules *before* reading this.

### 6.1 Subroutine `Pivoting(M,Ipiv,Idet)`

#### 6.1.1 Description

Permute the rows of  $M$  so that the biggest elements (in absolute value) of  $M$  are in the diagonal.

#### 6.1.2 Arguments

`M(:, :)`: Real or complex single or double precision two dimensional array. Initially it contains the matrix to permute, after calling the routine, it contains the permuted matrix. *Note that  $M$  is overwritten when calling this routine.*

`Ipiv(:)`: Integer one dimensional array. It returns the permutation of rows made to  $M$ .

`Idet`: Integer. If the number of permutations is odd, `Idet` = 1, if it is even `Idet` = -1

#### 6.1.3 Examples

Listing 6.1: Pivoting data of a matrix label

```
1 Program TestLinear
3   USE NumTypes
3   USE Linear
5
5   Integer, Parameter :: Nord = 4
7
7   Real (kind=DP) :: M(Nord,Nord), L(Nord,Nord), U(Nord,Nord), &
9       & Mcp(Nord,Nord)
9   Integer :: Ipiv(Nord), Iperm
11
```

```

13  ! Fill M of random numbers
    CALL Random_Number(M)
15
16  Write(*,*) 'Original M: '
17  Do I = 1, Nord
    Write(*, '(100ES10.3)')(M(I,J), J = 1, Nord)
19  End Do

21  CALL Pivoting (M, Ipiv, Iperm)
    Write(*,*) 'Permuted M: '
23  Do I = 1, Nord
    Write(*, '(100ES10.3)')(M(I,J), J = 1, Nord)
25  End Do

27  Stop
End Program TestLinear

```

## 6.2 Subroutine LU(M, Ipiv, Idet)

### 6.2.1 Description

Make the LU decomposition of matrix  $M$ . That is to say, given a matrix  $M$ , this function returns two matrix  $L$  and  $U$ , such that

$$M = LU \quad (6.1)$$

where  $L$  is lower triangular, and  $U$  upper triangular.

$$L = \begin{pmatrix} 1 & 0 & 0 & \dots \\ L_{21} & 1 & 0 & \dots \\ L_{31} & L_{32} & 1 & \dots \\ \vdots & \vdots & \vdots & \ddots \end{pmatrix}; \quad U = \begin{pmatrix} U_{11} & U_{12} & U_{13} & \dots \\ 0 & U_{22} & U_{23} & \dots \\ 0 & 0 & U_{33} & \dots \\ \vdots & \vdots & \vdots & \ddots \end{pmatrix} \quad (6.2)$$

The rows of  $M$  are permuted so that the biggest possible elements are on the diagonal (this makes the problem more stable). The two matrices  $L$  and  $U$  are returned *overwriting*  $M$ .

### 6.2.2 Arguments

**M(:, :):** Real or complex single or double precision two dimensional array. Initially it contains the matrix to decompose, after calling the routine, it contains  $L$  in its lower part, and  $U$  in its upper part. *Note that  $M$  is overwritten when calling this routine.*

**Ipiv(:):** Integer one dimensional array. It returns the permutation of rows made to  $M$ .

**Idet:** Integer. If the number of permutations is odd, **Idet** = 1, if it is even **Idet** = -1

## 6.2.3 Examples

Listing 6.2: Making the LU decomposition.

```

Program TestLinear
2
3  USE NumTypes
4  USE Linear
5
6  Integer, Parameter :: Nord = 4
7
8  Real (kind=DP) :: M(Nord,Nord), L(Nord,Nord), U(Nord,Nord), &
    & Mcp(Nord,Nord)
9  Integer :: I piv(Nord), Iperm
10
11
12  ! Fill M of random numbers, and make a copy
13  CALL Random_Number(M)
14  Mcp = M
15  L = 0.0_DP
16  U = 0.0_DP
17
18  ! Make the LU decomposition and fill the matrices
19  ! L and U
20  CALL Lu(M, I piv, Iperm)
21
22  Do I = 1, Nord
23      L(I,I) = 1.0_DP
24      U(I,I) = M(I,I)
25      Do J = I+1, Nord
26          L(J,I) = M(J,I)
27          U(I,J) = M(I,J)
28      End Do
29  End Do
30
31  ! Now Make the product and see that it is the original matrix with
32  ! some rows permuted
33  Write(*,*) 'M: '
34  Do I = 1, Nord
35      Write(*, '(100ES10.3)')(Mcp(I,J), J = 1, Nord)
36  End Do
37
38  Write(*,*) 'L: '
39  Do I = 1, Nord
40      Write(*, '(100ES10.3)')(L(I,J), J = 1, Nord)
41  End Do
42  Write(*,*) 'U: '
43  Do I = 1, Nord
44      Write(*, '(100ES10.3)')(U(I,J), J = 1, Nord)
45  End Do
46
47  M = MatMul(L,U)
48  Write(*,*) 'LU (Same as M with some rows permuted): '
49  Do I = 1, Nord

```

```

50      Write(*,'(100ES10.3)')(M(I,J), J = 1, Nord)
      End Do
52
54      Stop
      End Program TestLinear

```

## 6.3 Subroutine LUsolve(M, b)

### 6.3.1 Description

Solves the linear system of equations

$$\begin{aligned}
 M_{11}x_1 + M_{12}x_2 + M_{13}x_3 + M_{14}x_4 + \dots &= b_1 \\
 M_{21}x_1 + M_{22}x_2 + M_{23}x_3 + M_{24}x_4 + \dots &= b_2 \\
 &\vdots
 \end{aligned}
 \tag{6.3}$$

### 6.3.2 Arguments

**M(:, :):** Real or complex single or double precision two dimensional array. The matrix of coefficients. *M is overwritten when solving the system.*

**b(:):** Real or complex single or double precision one dimensional array. The independent terms before calling the routine, and the solution of the linear system of equations after calling it. *Note that b is overwritten when calling this routine.*

### 6.3.3 Examples

Listing 6.3: Solving systems of linear equations.

```

1  Program TestLinear
3
4      USE NumTypes
5      USE Linear
6
7      Integer, Parameter :: Nord = 10
8
9      Real (kind=DP) :: M(Nord,Nord), L(Nord,Nord), U(Nord,Nord), &
10         & Mcp(Nord,Nord), b(Nord), bcp(Nord)
11      Integer :: I piv(Nord), Iperm
12
13      ! Fill M and b of random numbers, and make a copy of both
14      CALL Random_Number(M)
15      CALL Random_Number(b)
16      Mcp = M
17      bcp = b
18
19      ! Solve the linear system
20      CALL LUsolve(M,b)

```

```

21  ! Check that it is a solution:
    b = MatMul(Mcp,b)
23  Write(*,*)'b: '
    Write(*,'(100ES10.3)')(Abs(bcp(I)-b(I)), I = 1, Nord)
25
27  Stop
End Program TestLinear

```

## 6.4 Function Det(M)

### 6.4.1 Description

Computes the determinant of the matrix  $M$ .

### 6.4.2 Arguments

$M(:, :)$ : Real or complex, simple or double precision two dimensional array. The matrix whose determinant we want to know.

### 6.4.3 Output

The value of the determinant. Same precision as the input argument.

### 6.4.4 Examples

Listing 6.4: Computing the determinant of a matrix.

```

Program TestLinear
2
  USE NumTypes
4  USE Linear

6  Integer, Parameter :: Nord = 10

8  Real (kind=DP) :: M(Nord,Nord), L(Nord,Nord), U(Nord,Nord), &
    & Mcp(Nord,Nord), b(Nord), bcp(Nord)
10 Integer :: I piv(Nord), Iperm

12
14  ! Fill M of randoms numbers
    CALL Random_Number(M)

16  ! Now compute the determinant.
    Write(*,'(ES15.8)') Det(M)
18
20  Stop
End Program TestLinear

```



# Seven

---

## MODULE NonNum

---

This is the documentation of the `MODULE NonNum`, a set of FORTRAN 90 routines to sort and search. This module make use of the `MODULE NumTypes`, and `MODULE Error` so please read the documentation of these modules *before* reading this.

### 7.1 Subroutine Swap(X,Ind1,Ind2)

#### 7.1.1 Description

Swaps elements `Ind1` and `Ind2` of the array `X(:)`.

#### 7.1.2 Arguments

`X(:)`: Integer, real single or real double precision one dimensional array. *Note that X is overwritten when calling this routine.*

`Ind1, Ind2`: Integer. The elements that we want to permute.

#### 7.1.3 Examples

Listing 7.1: Sorting data.

```
1 Program TestNN
3   USE NumTypes
4   USE NonNumeric
5
6   Integer, Parameter :: Nmax = 10
7   Integer :: Ima(Nmax), I
8
9   ! Fill Ima(:)
11  Forall (I=1:Nmax) Ima(I) = I
12
13  ! Plot the numbers
14  Do I = 1, Nmax
15      Write(*,'(1000I10)') Ima(I)
```

```

17      End Do
18      ! Swap first and last elemetns of Ima(:) and plot them.
19      CALL Swap(Ima, 1, Nmax)
20      Write(*,*)'# With first and last elements permuted: '
21      Do I = 1, Nmax
22          Write(*, '(1000I10) ') Ima(I)
23      End Do
24
25      Stop
27 End Program TestNN

```

## 7.2 Subroutine Insrt(X[, Ipt])

### 7.2.1 Description

Sort the elements of  $X(:)$  in ascendant order<sup>1</sup>.

### 7.2.2 Arguments

$X(:)$ : Integer, real single or real double precision one dimensional array. Initially it contains unsorted numbers, and after calling the routine, it contains the sorted elements. *Note that  $X$  is overwritten when calling this routine.*

$Ipt(:)$ : Integer vector, Optional. It returns the permutation made to  $X(:)$  to sort it.

### 7.2.3 Examples

Listing 7.2: Sorting data.

```

1 Program TestNN
2
3     USE NumTypes
4     USE NonNumeric
5
6     Integer, Parameter :: Nmax = 10
7     Integer :: Ima(Nmax)
8     Real (kind=DP) :: X(Nmax), Y(Nmax)
9
10
11     ! Fill X(:) with random data, and define Y(:)
12     CALL Random_Number(X)
13     Y = Sin(12.34_DP*(X-0.5_DP))
14
15     ! Plot an unsorted data table
16     Do I = 1, Nmax
17         Write(*, '(1000ES13.5) ') X(I), Y(I)

```

<sup>1</sup>This routine uses *insertion sort*, and is much slower than `Qsort`. To sort more than 10 elements, use `Qsort` unless you know what you are doing.



```

19      End Do
20
21      ! Sort them, and plot the table again. Same points, but this time
22      ! sorted
23      CALL Insrt(X, Ima)
24      Write(*,*)'# Again, this time sorted: '
25      Do I = 1, Nmax
26          Write(*,'(1000ES13.5)')X(I), Y(Ima(I))
27      End Do
28
29      Stop
End Program TestNN

```

## 7.3 Subroutine Qsort(X[, Ipt])

### 7.3.1 Description

Sort the elements of  $X(:)$  in ascendant order.

### 7.3.2 Arguments

$X(:)$ : Integer, real single or real double precision one dimensional array. Initially it contains unsorted numbers, and after calling the routine, it contains the sorted elements. *Note that  $X$  is overwritten when calling this routine.*

$Ipt(:)$ : Integer vector, Optional. It returns the permutation made to  $X(:)$  to sort it.

### 7.3.3 Examples

Listing 7.3: Sorting data.

```

Program TestNN
2
3      USE NumTypes
4      USE NonNumeric
5
6      Integer, Parameter :: Nmax = 10
7      Integer :: Ima(Nmax)
8      Real (kind=DP) :: X(Nmax), Y(Nmax)
9
10
11     ! Fill X(:) with random data, and define Y(:)
12     CALL Random_Number(X)
13     Y = Sin(12.34_DP*(X-0.5_DP))
14
15     ! Plot an unsorted data table
16     Do I = 1, Nmax
17         Write(*,'(1000ES13.5)')X(I), Y(I)
18     End Do

```

```

20      ! Sort them, and plot the table again. Same points, but this time
21      ! sorted
22      CALL Qsort(X, Ima)
23      Write(*,*)'# Again, this time sorted: '
24      Do I = 1, Nmax
25          Write(*, '(1000ES13.5) ')X(I), Y(Ima(I))
26      End Do
27
28      Stop
30 End Program TestNN

```

## 7.4 Function Locate( $X$ , $X_0$ [, $Iin$ ])

### 7.4.1 Description

Given a *sorted* vector of elements  $X(:)$ , and a point  $X_0$ , **Locate** returns the position  $n$  such that  $X(n) < X_0 < X(n+1)$ . If  $X_0$  is less than all the elements of  $X(:)$ , **Locate** returns 0, and if it is greater than all the elements of  $X(:)$ , it returns the number of elements of  $X(:)$

### 7.4.2 Arguments

$X(:)$ : Integer, real single or real double precision one dimensional *sorted* array.

$X_0$ : Integer, real single or real double precision number, but the same type as  $X(:)$ . Point that we want to locate in the sorted vector  $X(:)$ .

$Iin$ : Integer, Optional. Initial guess of the position.

### 7.4.3 Output

Integer. The position  $n$  such that

$$X(n) < X_0 < X(n+1)$$

### 7.4.4 Examples

Listing 7.4: Searching data position in an ordered list.

```

Program TestNN
2
3     USE NumTypes
4     USE NonNumeric
5
6     Integer, Parameter :: Nmax = 100
7     Integer :: Ima(Nmax), Idx
8     Real (kind=DP) :: X(Nmax), Y(Nmax), X0
9
10
11     ! Fill X(:) with random data, and set X0 to some arbitrary value.
12     CALL Random_Number(X)

```

```
14      X0 = 0.276546754_DP
16      ! Sort X(:), find the position of X0, and plot the neighborr
16      ! elements.
18      CALL Qsort(X)
18      Idx = Locate(X, X0)
20      Write(*, '(1A,1ES33.25)') 'Searched element: ', X0
20      Write(*, '(1A,1ES33.25)') 'Previous element in the list: ', X(Idx)
20      Write(*, '(1A,1ES33.25)') 'Next element in the list: ', X(Idx+1)
22
22      Stop
24 End Program TestNN
```



# Eight

---

## MODULE SpecialFunc

---

This is the documentation of the `MODULE SpecialFunc`, a set of `FORTRAN 90` routines to compute the value of some functions. This module make use of the `MODULE NumTypes`, `MODULE Constants`, `MODULE Error` so please read the documentation of these modules *before* reading this.

### 8.1 Function GammaLn(X)

#### 8.1.1 Description

Compute  $\log(\Gamma(X))$ .

#### 8.1.2 Arguments

**X:** Double (DP) precision. The point in which we want to know the value of  $\Gamma(X)$ .

#### 8.1.3 Output

A real Double precision (DP).

#### 8.1.4 Examples

This program should write the factorial of the first 100 numbers.

Listing 8.1: Computing the Gamma Function.

```
1 Program TestSpecialFunc
2
3   USE NumTypes
4   USE SpecialFunc
5
6   Integer :: q
7
8
9   Do q = 1, 100
10    Write(*, '(1A13,1I4,1A3,1ES33.25)') 'Factorial of:', q, ' = ', &
```

```

12      & exp(GammaLn(Real(q+1,kind=DP)))
13  End Do
14
15  Stop
16 End Program TestSpecialFunc

```

## 8.2 Function Theta(i, z, tau[, Prec])

### 8.2.1 Description

Compute the value of the  $i^{\text{th}}$  Jacobi theta function ( $i = 1, 2, 3, 4$ ) with nome  $q = e^{i\pi\tau}$

$$\vartheta_i(z|\tau) \quad (8.1)$$

For a definition and properties of these functions take a look [2], here we will only say that following the conventions of the cited reference, our Theta functions have quasi-periods  $\pi$  and  $\tau\pi$ .

### 8.2.2 Arguments

- i:** Integer. Which theta function we want to compute.  $i$  must have one of the following values: 1, 2, 3, 4.
- z:** Complex Double Precision (DPC) or Complex Single Precision (SPC). The point in which we want to compute the Theta function.
- tau:** Complex, with the same precision as **z**. is the quasi period of the Theta function. must be in the upper half plane ( $\text{Im}(\tau) > 0$ ).
- Prec:** Real, Optional. If **z** is DPC (SPC), **Prec** must be double precision (single precision). An estimation of the desired precision of the result. The default value is  $1 \times 10^{-3}$

### 8.2.3 Output

If **z** is Double Precision Complex (SPC), the the result will be Double Precision Complex (SPC).

### 8.2.4 Examples

Listing 8.2: Computing the Jacobi Theta functions.

```

2  Program TestSpecialFunc
3
4  USE NumTypes
5  USE SpecialFunc
6
7  Complex (DPC) :: Z, tau
8

```

```

10  Z = Cmplx(0.546734, 2.76457643, kind=DPC)
    tau = Cmplx(0.0_DP, 3.76387540_DP)

12  ! Check the quasi-periodicity of the Third
    ! Jacobi Theta function.
14  Write(*,*) Theta(3, Z, tau)
    Write(*,*) Theta(3, Z+Cmplx(PI_DP), tau)
16  Write(*,*) Theta(3, Z+PI_DP*tau, tau) * &
        &exp(PI*IMAG_DPC*tau + 2.0_DP*UNITIMAG_DPC*Z)
18
20  Stop
End Program TestSpecialFunc

```

## 8.3 Function ThetaChar(a, b, z, tau[, Prec])

### 8.3.1 Description

Computes the value of the Theta function with Characteristics  $(a, b)$  and quasi-periods  $(\pi, \pi\tau)$  in the point  $z$ :

$$\vartheta \left[ \begin{array}{c} a \\ b \end{array} \right] (z|\tau) \quad (8.2)$$

### 8.3.2 Arguments

**a, b:** Complex or Real, Single or double precision. The two characteristics of the Theta function.

**z:** Complex (Single or Double precision). The point in the complex plane.

**tau:** Complex (Single or Double precision). The quasi-period of the theta function. Must have  $\text{Im}(\tau) > 0$ .

**Prec:** Real (Single or Double precision). Optional. An estimation of the desired precision of the value of the theta function.

### 8.3.3 Output

Complex Single or Double precision, the same as the input values.

### 8.3.4 Examples

Listing 8.3: Computing the Jacobi Theta functions with characteristics.

```

1  Program TestSpecialFunc
3
3  USE NumTypes
  USE SpecialFunc
5
  Real(kind=DP) :: Deriv, X1, X2

```

```

7  Complex (DPC) :: Wmas, Wmenos, Z, tau
   Integer :: q, s
9
11  Z = Cmplx(0.546734, 2.76457643, kind=DPC)
   tau = Cmplx(0.0_DP, 3.76387540_DP)
13
15  Write(*,*) 'Theta 1:'
   Write(*,*) Theta(1, Z, tau)
17  Write(*,*) - ThetaChar(0.5_DP, 0.5_DP, Z, tau)
   Write(*,*) 'Theta 2:'
19  Write(*,*) Theta(2, Z, tau)
   Write(*,*) ThetaChar(0.5_DP, 0.0_DP, Z, tau)
21  Write(*,*) 'Theta 3:'
   Write(*,*) Theta(3, Z, tau)
23  Write(*,*) ThetaChar(0.0_DP, 0.0_DP, Z, tau)
   Write(*,*) 'Theta 4:'
25  Write(*,*) Theta(4, Z, tau)
   Write(*,*) ThetaChar(0.0_DP, 0.5_DP, Z, tau)
27
29  Stop
   End Program TestSpecialFunc

```

## 8.4 Function Hermite(n,x[, Dval])

### 8.4.1 Description

Returns the value of the  $n^{\text{th}}$  Hermite polynomial in the point  $X$ . If **Dval** is specified, the value of the Derivative of the  $n^{\text{th}}$  Hermite polynomial in the point  $X$  is also returned.

### 8.4.2 Arguments

**n**: Integer. Which Hermite polynomial wants to compute.

**x**: Real (Single or Double precision). The point in which we want to compute the Polynomial.

**Dval**: Real (Single or Double precision). Optional. If specified, it stores the value of the Derivative of the Polynomials.

### 8.4.3 Output

Real single or double precision (same as input). The value of the  $n^{\text{th}}$  Hermite Polynomial in the point  $X$ .

### 8.4.4 Examples

Listing 8.4: Computing the first 31 Hermite numbers.

```

Program TestSpecialFunc

```



```

2      USE NumTypes
4      USE SpecialFunc

6      Integer :: q

8
10     Write(*,*) 'The first 31 Hermite Numbers'
10     Write(*,*) 'http://www.research.att.com/~njas/sequences/A067994'
12     Do q = 1, 31
12         Write(*, '(1I4,1ES33.25)') q, Hermite(q, 0.0_DP)
14     End Do

14     Stop
16 End Program TestSpecialFunc

```

## 8.5 Function HermiteFunc(n, x[, Dval])

### 8.5.1 Description

Returns the value of the  $n^{\text{th}}$  Hermite function

$$\frac{1}{\sqrt{n!2^n\sqrt{\pi}}}e^{-x^2/2}H_n(x) \quad (8.3)$$

in the point  $X$ . If **Dval** is specified, the value of the Derivative of the  $n^{\text{th}}$  Hermite function in the point  $X$  is also returned.

### 8.5.2 Arguments

**n:** Integer. Which Hermite function wants to compute.

**x:** Real (Single or Double precision). The point in which we want to compute the Polynomial.

**Dval:** Real (Single or Double precision). Optional. If specified, it stores the value of the Derivative of the function.

### 8.5.3 Output

Real single or double precision (same as input). The value of the  $n^{\text{th}}$  Hermite function in the point  $X$ .

### 8.5.4 Examples

Listing 8.5: Compute the Hermite functions.

```

Program TestSpecialFunc

2      USE NumTypes
4      USE SpecialFunc

```

```

6  Real(kind=DP) :: Deriv, X1, X2, Sum
   Complex (DPC) :: Wmas, Wmenos, Z, tau
8  Integer :: q, s

10
12  Write(*,*) 'A (really bad) proof of orthonormality:'
   X1 = -10.0_DP
   Sum = 0.0_DP
14  Do q = -1000, 1000
       Sum = Sum + HermiteFunc(6,X1)**2
16     X1 = X1 + 0.01_DP
   End Do

18
20  Write(*, '(1 ES33.25) ') Sum*0.01_DP

22  Stop
End Program TestSpecialFunc

```

## 8.6 Function Basis(X1, X2, n, s, q, itau[, Prec])

### 8.6.1 Description

Return the value of the basis elements of the Hilbert space  $\mathcal{H}_q$  of quasi-periodic functions

$$|n, s\rangle = e^{i\frac{f}{2}x_1x_2} \sum_{k \in s+q\mathbb{Z}} e^{-u^2/2} H_n(u) e^{2\pi i k \frac{x_1}{l_1}} \quad n = 0, \dots, \infty; s = 1, \dots, q \quad (8.4)$$

defined in the appendix of [1] (look there for more details and properties).

### 8.6.2 Arguments

**X1,X2:** Real (Single or Double precision). The point in the Torus.

**n,s:** Integer. Specify which element of the basis.

**q:** Integer. Specify the Hilbert space  $\mathcal{H}_q$ .

**itau:** Real (Single or Double precision). Specify the ratio of quasi-periods:  $\text{itau} = l_2/l_1$ .

**Prec:** Real (Single or Double precision). Optional. An estimation of the desired precision.

### 8.6.3 Output

Complex single or double precision, depends of the input arguments.

### 8.6.4 Examples

Listing 8.6: Computing the bassi of a special Hilbert space (details in [1]).

```

2  Program TestSpecialFunc

```

```

4  USE NumTypes
   USE SpecialFunc

6  Real(kind=DP) :: X1, X2
   Complex (DPC) :: Wmas, Wmenos,
8  Integer :: I, q, s

10
11  Write(*,*)'Looking at the quasi-periodicity properties:'
12  X1 = 0.97834D0
13  X2 = 0.873873D0
14  q = 4
15  s = 3
16  Do I = 0, 8
17     Wmas = Basis( X1, X2+1.0_DP, I, s, q, 1.0_DP, 1.0D-15) * &
18           & exp(PLIMAG_DPC*X1*q)
19     Wmenos = Basis( X1+1.0_DP, X2, I, s, q, 1.0_DP, 1.0D-15) * &
20           & exp(-PLIMAG_DPC*X2*q)
21     Write(*,'(1I3,2ES33.25)')I, Basis( X1, X2, I, s, q, 1.0_DP, 1.0D-15)
22     Write(*,'(1I3,2ES33.25)')I, Wmas
23     Write(*,'(1I3,2ES33.25)')I, Wmenos
24 End Do

26
27 Stop
28 End Program TestSpecialFunc

```

## 8.7 Function Factorial(N)

### 8.7.1 Description

Compute  $N!$ . Better (faster and more accurate for small numbers) than the use of `GammaLn` to compute the factorial of a number.

### 8.7.2 Arguments

**N:** Integer. The number to compute the factorial.

### 8.7.3 Output

A real Double precision (DP).

### 8.7.4 Examples

This program should write the factorial of the first 100 numbers.

Listing 8.7: Computing the factorial.

```

2  Program TestSpecialFunc

   USE NumTypes

```

```
4  USE SpecialFunc
6
8  Integer :: q
10 Do q = 1, 100
12   Write(*, '(1A13,1I4,1A3,1ES33.25)') 'Factorial of:', q, ' = ', &
    & Factorial(q)
14 End Do
16 Stop
End Program TestSpecialFunc
```

# Nine

---

## MODULE Statistics

---

This is the documentation of the `MODULE Statistics`, a set of FORTRAN 90 routines to perform statistical description of data. This module make use of the `MODULE NumTypes`, `MODULE Constants`, `MODULE Error` and `MODULE Linear` so please read the documentation of these modules *before* reading this.

### 9.1 Function Mean(X)

#### 9.1.1 Description

Compute the mean value of the numbers stored in `X(:)`.

#### 9.1.2 Arguments

`X(:)`: Double (DP) or simple (SP) precision one dimensional array. The values whose mean we want to compute.

#### 9.1.3 Output

A real double or simple precision (same type as the input). The mean of the values.

#### 9.1.4 Examples

Listing 9.1: Computing the Mean of a vector of numbers.

```
1 Program Tests
2
3   USE NumTypes
4   USE Error
5   USE Statistics
6
7   Integer, Parameter :: Nmax = 100
8   Real (kind=DP) :: X(Nmax)
9
10  CALL Random_Number(X)
11  Write (*, '(ES33.25) ') Mean(X)
```

```

12      Stop
14 End Program Tests

```

## 9.2 Function Median(X)

### 9.2.1 Description

Compute the median value of the numbers stored in `X(:)`.

### 9.2.2 Arguments

`X(:)`: Double (DP) or simple (SP) precision one dimensional array. The values whose median we want to compute.

### 9.2.3 Output

A real double or simple precision (same type as the input). The median of the values.

### 9.2.4 Examples

Listing 9.2: Computing the Median of a vector of numbers.

```

Program Tests
2
   USE NumTypes
4   USE Error
   USE Statistics
6
   Integer, Parameter :: Nmax = 5
8   Real (kind=SP) :: X(Nmax) = (/1.0, 1.0, 2.0, 4.0, 1.5/)
10  Write(*, '(ES33.25) ') Median(X)
12
   Stop
14 End Program Tests

```

## 9.3 Function Var(X)

### 9.3.1 Description

Compute the variance of a vector of numbers `X(:)`

### 9.3.2 Arguments

`X(:)`: Double (DP) or simple (SP) precision one dimensional array. The values whose variance we want to compute.

### 9.3.3 Output

A real double or simple precision (same type as the input). The variance of the values.

### 9.3.4 Examples

Listing 9.3: Computing the Variance of a set of numbers.

```

1 Program Tests
2
3   USE NumTypes
4   USE Error
5   USE Statistics
6
7   Integer, Parameter :: Nmax = 100, Npinta = 100, Npar = 4
8   Real (kind=DP) :: X(Nmax), Y(Nmax), Yerr(Nmax), &
9       & Coef(Npar), Cerr(Npar), Corr, Xd(Nmax,2)
10
11
12   CALL Random_Number(X)
13   Write(*, '( ES33.25 ) ') Var(X)
14
15
16   Stop
17 End Program Tests

```

## 9.4 Function Stddev(X)

### 9.4.1 Description

Computes the standard deviation of the numbers stored in the vector **X(:)**.

### 9.4.2 Arguments

**X(:)**: Double (DP) or simple (SP) precision one dimensional array. The values whose standard deviation we want to compute.

### 9.4.3 Output

Real Single or Double precision, the same as the input values. The standard deviation of the values.

### 9.4.4 Examples

Listing 9.4: Computing the standard deviation.

```

1 Program Tests
2
3   USE NumTypes
4   USE Error
5   USE Statistics

```

```

7  Integer, Parameter :: Nmax = 100, Npinta = 100, Npar = 4
Real (kind=DP) :: X(Nmax), Y(Nmax), Yer(Nmax), &
9      & Coef(Npar), Cerr(Npar), Corr, Xd(Nmax,2)

11
CALL Random_Number(X)
13 Write(*,'(ES33.25)') Stddev(X)

15
Stop
17 End Program Tests

```

## 9.5 Function Moment(X, k)

### 9.5.1 Description

Returns the  $k^{th}$  moment of the values stored in the vector  $X(:)$ .

### 9.5.2 Arguments

$X(:)$ : Real (Single or Double precision). The numbers whose  $k^{th}$  moment we want to compute.

$k$ : Integer. Which moment we want to compute.

### 9.5.3 Output

Real single or double precision. The  $k^{th}$  moment of the numbers.

### 9.5.4 Examples

Listing 9.5: Computing the  $k^{th}$  moment of a data set.

```

1  Program Tests

3  USE NumTypes
USE Error
5  USE Statistics

7  Integer, Parameter :: Nmax = 100, Npinta = 100, Npar = 4
Real (kind=DP) :: X(Nmax), Y(Nmax), Yer(Nmax), &
9      & Coef(Npar), Cerr(Npar), Corr, Xd(Nmax,2)

11
CALL Random_Number(X)
13 Write(*,*) 'We should obtain the same numbers twice: '
Write(*,'(ES33.25)') Moment(X,2), Var(X)

15
Stop
17 End Program Test

```



## 9.6 Subroutine Histogram(Val, Ndiv, Ntics, Vmin, Vmax, h)

### 9.6.1 Description

Given a set of points `Val(:)`, this routine makes `Ndiv` divisions between the minimum and the greatest value of `Val` (respectively returned in `Vmin` and `Vmax`), each of size `h` (also returned), and returns in the integer vector `Nticks(:)` the number of points that are in each interval.

### 9.6.2 Arguments

**Val(:):** Real (Single or Double precision) one dimensional array. The original values.

**Ndiv:** Integer. The number of divisions.

**Nticks:** Integer one dimensional array. `Ndiv(I)` Tells how many points of `Val(:)` are between  $Vmin + (I - 1)h$  and  $Vmin + Ih$ .

**Vmin, Vmax:** Real (Single or Double precision). The minimum and maximum values of `Val`.

**h:** Real (Single or Double precision). After calling the routine has the step of the division.

### 9.6.3 Examples

Listing 9.6: Making Histograms.

```

1  Program Tests
3
3  USE NumTypes
3  USE Error
5  USE Statistics
7
7  Integer, Parameter :: Nmax = 500000, Npinta = 100, Npar = 4, Ndiv = 100
7  Real (kind=DP) :: X(Nmax), Y(Nmax), Yer(Nmax), &
9      & Coef(Npar), Cerr(Npar), Corr, Xd(Nmax,2), &
9      & Xmin, Xmax, h, Xac
11 Integer :: Ntics(Ndiv)
13
13 CALL Normal(X, 1.23_DP, 0.345_DP)
13 CALL Histogram(X, Ndiv, Ntics, Xmin, Xmax, h)
15
15 Do I = 1, Ndiv
17     Xac = Xmin + (I-1)*h
17     Write(*, '(1ES33.25,1I)') Xac, Ntics(I)
19 End Do
21
21 Stop
End Program Tests

```

## 9.7 Subroutine LinearReg(X, Y, Yerr, [Func], Coef, Cerr, ChisqrV)

### 9.7.1 Description

Given a set of points  $X(:)$  and  $Y(:)$ , this routine performs a linear fit to a set of functions defined by `Func`.

$$Y = \sum_i a_i f_i(X)$$

This routine also performs multi-dimensional fitting, in which case the points are specified as  $X(:, :)$ , where the first argument tells which point, and the second which variable.

### 9.7.2 Arguments

**X(:, :):** Real single or double precision one dimensional array (for a one dimensional fit) or two dimensional array (for a multidimensional fit). The independent variables. For a multidimensional fit, the first argument tells which point, and the second which variable. So the size of the array should be `X(Npoints, Ndim)`.

**Y(:):** Real single or double precision one dimensional array. The dependent variable.

**Yerr(:):** Real single or double precision one dimensional array. The errors of the points. If you don't have them, you should put all of them to some non-zero value.

**Func:** Optional. This routine defines the functions to fit. An interface like this should be provided

```
Interface
  Function Func(Xx, i)

    USE NumTypes

    Real (kind=SP), Intent (in) :: Xx
    Integer, Intent (in) :: i
    Real (kind=SP) :: Func

  End Function Func
End Interface
```

if you want to perform a one dimensional fitting, and like this

```
Interface
  Function Func(Xx, i)

    USE NumTypes

    Real (kind=SP), Intent (in) :: Xx(:)
    Integer, Intent (in) :: i
```

```

      Real (kind=SP) :: Func

      End Function Func
End Interface

```

if it is a multidimensional fitting. Since you are making a fitting to a function of the type

$$Y = \sum_i a_i f_i(X)$$

the values  $f_i(X)$  are given by this function as **Func(X, I)**. If the functions are not specified (i.e. you don't put this argument), a fit to a polynomial is made (this only work for one-dimensional fittings).

**Coef(:):** Real single or double precision one dimensional array. The parameters that you want to determine.

**Cerr(:):** Real single or double precision one dimensional array. The errors in the parameters.

**ChiSqr:** Real single or double precision. The  $\chi^2$  per degree of freedom of the fit.

### 9.7.3 Examples

Listing 9.7: Doing linear regressions.

```

Program Tests
2
  USE NumTypes
4  USE Error
  USE Statistics
6
  Integer, Parameter :: Nmax = 200, Npinta = 100, Npar = 4, Ndiv = 100
8  Real (kind=DP) :: X(Nmax), Y(Nmax), Yerr(Nmax), &
    & Coef(Npar), Cerr(Npar), Corr, Xd(Nmax,2), &
10    & Xmin, Xmax, h, Xac
  Integer :: Ntics(Ndiv)
12
  Interface
14    Function Fd(Xx, i)
      USE NumTypes
16
      Real (kind=DP), Intent (in) :: Xx(:)
      Integer, Intent (in) :: i
      Real (kind=DP) :: Fd
20
22    End Function FD
  End Interface
24
26  CALL Random_Number(Xd)
  Xd(:, :) = 10.0_DP*(Xd(:, :) - 0.8_DP)

```

```

28      CALL Normal(Yer, 0.0_DP, 1.0E-3_DP)
30      Y(:) = 12.34_DP*Xd(:,1)*sin(Xd(:,2)) - 2.23_DP + &
           & 0.67_DP*Xd(:,1)**2*Xd(:,2) + 0.23_DP*Xd(:,1) + Yer(:)
32
34      CALL LinearReg(Xd, Y, Yer, Fd, Coef, Cerr, Corr)
36
37      ! This should print the adjusted parameters,
38      ! that have values: 12.34, -2.23, 0.67, 0.23
39      Do I = 1, Npar
40          Write(*, '(2ES33.25)') Coef(I), Cerr(I)
41      End Do
42
43      ! This prints the ChiSqr, that should be very
44      ! close to 1.
45      Write(*, '(1A,1ES33.25)') 'ChiSqr of the Fit: ', Corr
46
47      Stop
48  End Program Tests
49
50  ! *****
51  ! *
52  Function Fd(X, i)
53  ! *
54  ! *****
55
56      USE NumTypes
57
58      Real (kind=DP), Intent (in) :: X(:)
59      Integer, Intent (in) :: i
60      Real (kind=DP) :: Fd
61
62      If (I==1) Then
63          Fd = 1.0_DP
64      Else If (I==2) Then
65          Fd = X(1)*sin(X(2))
66      Else If (I==3) Then
67          Fd = X(1)**2*X(2)
68      Else If (I==4) Then
69          Fd = X(1)
70      End If
71
72      Return
73  End Function FD

```

## 9.8 Subroutine NonLinearReg(X, Y, Yerr, Func, Coef, Cerr, ChisqrV)

### 9.8.1 Description

Given a set of points  $X(:)$  and  $Y(:)$ , this routine performs a non-linear fit to a set of functions defined by **Func**.

This routine also performs multi-dimensional fitting, in which case the points are specified as  $X(:, :)$ , where the first argument tells which point, and the second which variable.

This routine uses the Levenberg-Marquardt algorithm to perform the optimisation<sup>1</sup>.

### 9.8.2 Arguments

**X(:, :)**: Real single or double precision one dimensional array (for a one dimensional fit) or two dimensional array (for a multidimensional fit). The independent variables. For a multidimensional fit, the first argument tells which point, and the second which variable. So the size of the array should be  $X(Npoints, Ndim)$ .

**Y(:)**: Real single or double precision one dimensional array. The dependent variable.

**Yerr(:)**: Real single or double precision one dimensional array. The errors of the points. If you don't have them, you should put all of them to some non-zero value.

**Func**: This routine define the functions to fit. An interface like this should be provided

```
Interface
```

```
Subroutine Func(X, Cf, Valf, ValD)
```

```
USE NumTypes
```

```
Real (kind=SP), Intent (in) :: X, Cf(:)
```

```
Real (kind=SP), Intent (out) :: Valf, ValD(Size(Cf))
```

```
End Subroutine Func
```

```
End Interface
```

if you want to perform a one dimensional fitting, and like this

```
Interface
```

```
Subroutine Func(X, Cf, Valf, ValD)
```

```
USE NumTypes
```

```
Real (kind=SP), Intent (in) :: X(:), Cf(:)
```

```
Real (kind=SP), Intent (out) :: Valf, ValD(Size(Cf))
```

```
End Subroutine Func
```

```
End Interface
```

---

<sup>1</sup>[http://en.wikipedia.org/wiki/Levenberg-Marquardt\\_algorithm](http://en.wikipedia.org/wiki/Levenberg-Marquardt_algorithm)

if it is a multidimensional fitting.

This routine returns the values of the function at  $X$  for some values of the parameters given in **Cf** in the variable **Valf**, and a vector with the derivatives (respect with the parameters) in **ValD**(:).

**Coef**(:): Real single or double precision one dimensional array. Output. The parameters that you want to determine.

**Cerr**(:): Real single or double precision one dimensional array. Output. The errors in the parameters.

**ChiSqr**: Real single or double precision. The  $\chi^2$  per degree of freedom of the fit.

### 9.8.3 Examples

In this example we will fit some generated data, with a Normal noise to the function

$$f(x_1, x_2; a, b) = \sin(ax_1) + bx_1x_2$$

We will generate the data with the values  $a = 2.0$  and  $b = 0.2$ , so our fitting routine *should* return this values within errors.

The derivatives of the function (repect the parameters  $a$  and  $b$ ), as well as the value of function are given by the user routine **Func**. The derivatives are:

$$\begin{aligned} \frac{\partial f(x_1, x_2; a, b)}{\partial a} &= x_1 \cos ax_1 \\ \frac{\partial f(x_1, x_2; a, b)}{\partial b} &= x_1x_2 \end{aligned}$$

Listing 9.8: Doing non linear regressions.

```

1  Program NLFit
3      USE NumTypes
4      USE Statistics
5
6      Interface
7          Subroutine Func(X, Cf, Valf, ValD)
8
9              USE NumTypes
10
11             Real (kind=DP), Intent (in) :: X(:), Cf(:)
12             Real (kind=DP), Intent (out) :: Valf, ValD(Size(Cf))
13
14         End Subroutine Func
15     End Interface
16
17     Integer, Parameter :: Np = 20, Ndim = 2
18     Real (kind=DP) :: X(Np, Ndim), Y(Np), Ye(Np), Co(2), Vd(2), Ce(2), Ch
19
20     ! First Fill the data

```

```

21  CALL Random_Number(X)
    X = 2.0_DP*(X - 0.5_DP)
23  Co(1) = 2.0_DP
    Co(2) = 0.2_DP
25  CALL Normal(Ye, 0.0_DP, 0.5_DP)
    Do I = 1, Np
27      CALL Func(X(I,:), Co, Y(I), Vd)
        Y(I) = Y(I) + Ye(I)
29  End Do

31
    ! Now Perform the non linear fit
33  Co = 1.0_DP
    CALL NonLinearReg(X, Y, Abs(Ye), Func, Co, Ce, Ch)
35  Do I = 1, Npar
        Write(*, '(1A,100ES33.25)') 'Parameter and error: ', Co(I), Ce(I)
37  End Do
    Write(*, '(1A,100ES33.25)') &
39  & 'Chi Square per degree of freedom of the Fit: ', Ch

41
    Stop
43 End Program NLFit

45 Subroutine Func(X, Cf, Valf, ValD)

47  USE NumTypes

49  Real (kind=DP), Intent (in) :: X(:), Cf(:)
    Real (kind=DP), Intent (out) :: Valf, ValD(Size(Cf))

51  Valf = Sin(Cf(1)*X(1)) + Cf(2)*X(1)*X(2)
53  ValD(1) = X(1)*Cos(Cf(1)*X(1))
    ValD(2) = X(1)*X(2)
55
57 End Subroutine Func

```

## 9.9 Subroutine SetLuxLevel(Ilevel)

### 9.9.1 Description

This function changes the behaviour of the “intrinsic” `Random_Number` subroutine. To call the intrinsic `FORTTRAN 90` routine, you should set the Luxury level to zero. Other values makes the `Random_Number` subroutine use the Marsaglia and Zaman algorithm modified by M. Lüscher [3]. This part of the code has some inspiration in the `FORTTRAN 90` implementation of Allan Miller. **Note:** This routine affects the way all the other routines that use pseudo random number generators works. That is to say, all the routines of this module that follow this one.

### 9.9.2 Arguments

**Ilevel:** Integer. Set the way the random number generator works. We have the following options:

- **Ilevel** = 0. The default intrinsic **FROTRAN 90** routine is used. This is usually considered a *bad*, non portable pseudo-random number generator, that you do not want to use when correlations are important.
- **Ilevel** = 1. This is the original Marsaglia and Zaman algorithm. Probably better than the intrinsic procedure, but still with large correlations.
- **Ilevel** = 2. The M. Lüscher modified version of the algorithm with  $p = 48$ . Still fail many tests, but a very good pseudo-number generator for many things.
- **Ilevel** = 3. The M. Lüscher modified version of the algorithm with  $p = 97$ . Theoretically still defective (fail the serial correlation test).
- **Ilevel** = 4. The M. Lüscher modified version of the algorithm with  $p = 218$ . Any theoretically possible correlation has a very small probability of being observed. A value very similar to this has been used for large scale high precision lattice gauge theory computations. This corresponds with the Lüscher recommended level 0.
- **Ilevel** = 5. The M. Lüscher modified version of the algorithm with  $p = 404$ . Good for anything you need the pseudo-random numbers (but cryptography, of course). It seems completely pointless to use values of  $p$  higher than this. This corresponds with the Lüscher recommended level 1. This is the **default value**.
- **Ilevel** = 6. The M. Lüscher modified version of the algorithm with  $p = 794$ . Good for anything you need the pseudo-random numbers (but cryptography, of course). Probably waste of computer resources. This corresponds with the Lüscher recommended level 2.
- Other values of **Ilevel** ( $> 24$ ). Sets the value of  $p$  in the M. Lüscher modified version of the algorithm. Don't use it, unless you know what you are doing.

### 9.9.3 Examples

Listing 9.9: Setting the Luxury level of the pseudo random number generator.

```

1  Program Tests
3      USE NumTypes
4      USE Statistics
5
6      Integer, Parameter :: NN = 70
7      Real (kind=SP) :: rr(NN)
8
9      CALL SetLuxLevel(5)
10     CALL Random_Number(rr)
11     Write(*,*) rr
12
13     Stop
14
15 End Program Tests

```



## 9.10 Subroutine PutLuxSeed([ISeed(25)])

### 9.10.1 Description

Restarts the position in the Luxury pseudo-random number generator.

### 9.10.2 Arguments

**ISeed(25):** Integer one dimensional array of 25 elements. Optional. If present, restarts the generator from an output of **GetLuxSeed**. If not present, initialises the random number generator.

### 9.10.3 Examples

Listing 9.10: Using a previously saved point in the generating process.

```

1 Program Tests
3   USE NumTypes
3   USE Statistics
5
5   Integer, Parameter :: NN = 70
7   Real (kind=SP) :: rr(NN)
7   Integer :: Sd(25)
9
11  Open(Unit=69, File="seed.dat")
11  Read(69,*) Sd
13  CALL PutLuxSeed(Sd)
13  CALL Random_Number(rr)
15  Write(*,*) rr
17
17  Stop
19 End Program Tests

```

## 9.11 Subroutine GetLuxSeed(ISeed(25))

### 9.11.1 Description

Saves the position in the Luxury pseudo-random number generator.

### 9.11.2 Arguments

**ISeed(25):** Integer one dimensional array of 25 elements. Saves the position in the random number generator.

### 9.11.3 Examples

Listing 9.11: Saving a point in the generating process.

```

1  Program Tests
3      USE NumTypes
   USE Statistics
5
   Integer, Parameter :: NN = 70
7   Real (kind=SP) :: rr(NN)
   Integer :: Sd(25)
9
11  CALL Random_Number(rr)
   CALL GetLuxSeed(Sd)
13  Open(Unit=69, File="seed.dat")
   Write(69,*)Sd
15
17  Stop
   End Program Tests

```

## 9.12 Subroutine Normal(X, [Rm], [Rsig])

### 9.12.1 Description

Fills `X(:)` with numbers from a normal distribution with mean `Rm`, and standard deviation `Rsig`. The parameters `Rm` and `Rsig` are optional. If they are not given the mean will be 0, and the standard deviation 1.

### 9.12.2 Arguments

**X(:):** Real (Single or Double precision) one dimensional array. A vector that will be filled with numbers according to the normal distribution.

**Rm:** Real (Single or Double precision), Optional. The mean of the normal distribution. If not present the default value is 0.

**Rsig:** Real (Single or Double precision), Optional. The standard deviation of the normal distribution. If not present the default value is 1.

### 9.12.3 Examples

Listing 9.12: Obtaining numbers with a normal distribution.

```

2  Program Tests
   USE NumTypes
4   USE Error
   USE Statistics

```

```

6      Integer, Parameter :: Nmax = 100
8      Real (kind=DP) :: X(Nmax)

10
12      CALL Normal(X, 1.23_DP, 0.345_DP)
14      ! Now compute the mean and standard deviation of the data
16      Write(*,*) 'We should obtain 1.23 and 0.345: '
18      Write(*, '(ES33.25) ') Mean(X), Stddev(X)

      Stop
End Program Tests

```

## 9.13 Subroutine FishTipp(X, Rm, Rb)

### 9.13.1 Description

Fills  $X(:)$  with numbers from a Fisher-Tippett distribution with parameters<sup>2</sup>  $Rm$ , and  $2Rb^2$ .

### 9.13.2 Arguments

$X(:)$ : Real (Single or Double precision) one dimensional array. A vector that will be filled with numbers according to the normal distribution.

$Rm$ : Real (Single or Double precision).

$Rb$ : Real (Single or Double precision).

### 9.13.3 Examples

Listing 9.13: Obtaining numbers with a Fisher-Tippett distribution.

```

Program Tests

2      USE NumTypes
4      USE Error
6      USE Statistics

8      Integer, Parameter :: Nmax = 100
10     Real (kind=DP) :: X(Nmax)

12     CALL FishTipp(X, 2.00_DP, 1.00_DP)
14     ! Now compute the mean and standard deviation of the data
16     Write(*,*) 'We should obtain 2.57721... and 1.2782...: '
18     Write(*, '(ES33.25) ') Mean(X), Stddev(X)

```

<sup>2</sup>More info about this distribution in the Wikipedia: [http://en.wikipedia.org/wiki/Fisher-Tippett\\_distribution](http://en.wikipedia.org/wiki/Fisher-Tippett_distribution)

```

18      Stop
      End Program Tests

```

## 9.14 Subroutine Laplace(X, Rm, Rb)

### 9.14.1 Description

Fills `X(:)` with numbers from a Laplace distribution with mean `Rm`, and variance  $2Rb^2$ .

### 9.14.2 Arguments

`X(:)`: Real (Single or Double precision) one dimensional array. A vector that will be filled with numbers according to the Laplace distribution.

`Rm`: Real (Single or Double precision). The mean of the Laplace distribution.

`Rb`: Real (Single or Double precision). The width of the Laplace distribution (i.e. The variance is  $2Rb^2$ ).

### 9.14.3 Examples

Listing 9.14: Obtaining numbers with a Laplace distribution.

```

Program Tests
2
  USE NumTypes
4  USE Error
  USE Statistics
6
  Integer, Parameter :: Nmax = 100
8  Real (kind=DP) :: X(Nmax)
10
  CALL Laplace(X, 1.23_DP, 1.0_DP)
12  ! Now compute the mean and standard deviation of the data
  Write(*,*) 'We should obtain 1.23 and sqrt(2): '
14  Write(*, '(ES33.25) ') Mean(X), Stddev(X)
16
  Stop
18 End Program Tests

```

## 9.15 Subroutine/Function Irand([Irnd], N, M)

### 9.15.1 Description

If present, fills `Irnd(:)` with random integer numbers between `N` and `M` with an uniform distribution. If `Irnd(:)` is not present returns a integer random number between `N` and `M`.

### 9.15.2 Arguments

**Irnd(:):** Integer, Optional. A vector that will be filled with integer numbers according to a uniform distribution.

**N:** Integer. The minimum number that we can obtain.

**M:** Integer. The maximum number that we can obtain.

### 9.15.3 Examples

Listing 9.15: Obtaining integer random numbers.

```

1 Program Tests
2
3   USE NumTypes
4   USE Error
5   USE Statistics
6
7   Integer, Parameter :: Nmax = 100
8   Integer :: Irnd(Nmax)
9
10
11  CALL Irand(Irnd, 0, 1)
12  ! Now compute the mean
13  Write(*,*) 'We should obtain 0.5: '
14  Write(*, '(ES33.25) ') Mean(Real(Irnd(:), kind=DP))
15
16
17  Stop
18 End Program Tests

```

## 9.16 Subroutine Permutation(Idx)

### 9.16.1 Description

Returns a random permutation of N elements. It uses the Knuth shuffle algorithm<sup>3</sup>.

### 9.16.2 Arguments

**Idx(:):** Integer one dimensional array. Output. The random permutation.

Listing 9.16: Obtaining a permutation.

```

1 Program Tests
2
3   USE NumTypes
4   USE Error
5   USE Statistics

```

<sup>3</sup>[http://en.wikipedia.org/wiki/Knuth\\_shuffle](http://en.wikipedia.org/wiki/Knuth_shuffle)

```

6      Integer, Parameter :: Nmax = 10
8      Integer :: Id(Nmax)

10
12     CALL Permutation(Id)
12     Write(*, '(100I3)')(Id(I), I = 1, Nmax)

14     Stop

16 End Program Tests

```

## 9.17 Subroutine BootStrap(Ibt)

### 9.17.1 Description

Generates  $N_b$  bootstrap sequence of  $N$  numbers each. These bootstraps are returned in the two dimensional integer array  $Ibt(:, :)$  of size  $N \times N_b$ .

### 9.17.2 Arguments

$Ibt(:, :)$ : Integer. A two dimensional array of size  $N \times N_b$ , where  $N_b$  is the number of bootstraps, and  $N$  is the range of each bootstrap.

### 9.17.3 Examples

Listing 9.17: Resampling some data.

```

Program Tests
2
3     USE NumTypes
4     USE Error
5     USE Statistics
6
7     Integer, Parameter :: Nmax = 8, Nbt = 5
8     Real (kind=DP) :: X(Nmax)
9     Integer :: Ib(Nmax, Nbt), I, J
10
11
12     CALL Random_Number(X)
13     ! Generate 5 bootstraps
14     CALL BootStrap(Ib)
15
16     ! Write the original sample, and the bootstraps
17     Write(*, '(1000ES33.25)')(X(J), J=1, Nmax)
18     Do I = 1, Nbt
19         Write(*, '(1000ES33.25)')(X(Ib(J)), J=1, Nmax)
20     End Do
21
22     Stop
End Program Tests

```

## 9.18 Subroutine SaveBstrp(Ibt, Filename)

### 9.18.1 Description

Saves the bootstrap stored in **Ibt** and saves it in the file **Filename**.

### 9.18.2 Arguments

**Ibt(:):** Integer. A two dimensional array of size  $N \times N_b$ , where  $N_b$  is the number of bootstraps, and  $N$  is the range of each bootstrap.

**Filename:** Character (len=\*). A file name to save the resampling data.

### 9.18.3 Examples

Listing 9.18: Reading the resampling info.

```

1  Program Tests
3      USE NumTypes
4      USE Error
5      USE Statistics
7      Integer, Parameter :: Nmax = 8, Nbt = 5
8      Integer :: Ib(Nmax, Nbt)
9
11     ! Generate 5 bootstraps
12     CALL BootStrap(Ib)
13
14     ! Save it
15     SaveBstrp(Ibt, 'example.bst')
17     Stop
End Program Tests

```

## 9.19 Subroutine ReadBstrp(Ibt, Filename)

### 9.19.1 Description

Reads the bootstrap stored in the file **Filename**, and returns it in **Ibt**.

### 9.19.2 Arguments

**Ibt(:):** Integer. A two dimensional array of size  $N \times N_b$ , where  $N_b$  is the number of bootstraps, and  $N$  is the range of each bootstrap.

**Filename:** Character (len=\*). A file name to read the resampling data.

### 9.19.3 Examples

Listing 9.19: Saving the resampling info.

```

1  Program Tests
2
3      USE NumTypes
4      USE Error
5      USE Statistics
6
7      Integer, Parameter :: Nmax = 8, Nbt = 5
8      Integer :: Ib(Nmax, Nbt)
9
10     ! Read a saved Bootstrap.
11     ReadBstrp(Ibt, 'example.bst')
12
13     Stop
14 End Program Tests

```

## 9.20 Subroutine EstBstrp(Data, Ibt, Func, Val, Err[, Rest])

### 9.20.1 Description

Estimates using the Bootstrap method the average and error of an estimator given as a user supplied function.

### 9.20.2 Arguments

**Data(:):** Double precision Real. A one dimensional array with the original sampling.

**ibt(:,):** Integer two dimensional array. The bootstrap that we want to use to make the estimation.

**Func:** A user supplied function that returns the value of the estimator. An interface block of the following type should be defined.

```

Interface
  Function Func(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X(:)
    Real (kind=DP) :: Func

  End Function Func
End Interface

```

**Val:** Double precision real. Output. The value of the estimation of the parameter.

**Err:** Double precision real. Output. An estimation of the error in the estimation of the parameter.



**Rest:** Double precision real one dimensional array (same dimension as the number of bootstraps in Ibt). Output. The value of the estimator for each resampling.

### 9.20.3 Examples

Listing 9.20: Estimating the average.

```

1  Program Tests
3      USE NumTypes
4      USE Error
5      USE Statistics
7      Integer, Parameter :: Nmax = 100, Nbt = 50
8      Integer :: Ib(Nmax, Nbt)
9      Real (kind=DP) :: Avg, Err, Data(Nmax), Rest(Nbt)
11     Interface
12         Function F(X)
13             USE NumTypes
15             Real (kind=DP), Intent (in) :: X(:)
16             Real (kind=DP) :: F
17         End Function F
18     End Interface
19
21     ! Read a saved Bootstrap, and the data from a file
22     ReadBstrp(Ibt, 'example.bst')
23     Open (Unit=22, File="data.dat")
24     Read(22,*)Data
25     Close(22)
27     ! And estimate the average
28     CALL EstBstrp(Data, Ibt, F, Avg, Err, Rest)
29
30     ! Print the Average of each resampling
31     Do I = 1, Nbt
32         Write(*,*)I, Rest(I)
33     End Do
35     Stop
36 End Program Tests
37
38 Function F(X)
39     USE NumTypes
40     USE Statistics
41
43     Real (kind=DP), Intent (in) :: X(:)
44     Real (kind=DP) :: F
45
46     F = Mean(X)

```

47

```
End Function F
```

## 9.21 Subroutine BstrpConfInt(Data, Ibt, alpha, Func, dmin, dpls))

### 9.21.1 Description

Gives an Confidence interval for the estimator given as the user supplied function `Func`, such that

$$\mathcal{P}(dmin < Func(Data) < dpls) = 1 - 2\alpha$$

### 9.21.2 Arguments

**Data(:):** Double precision Real. A one dimensional array with the original sampling.

**ibt(:,):** Integer two dimensional array. The bootstrap that we want to use to make the estimation.

**alpha:** Double precision real. The level of the confidence interval.

**Func:** A user supplied function that returns the value of the estimator. An interface block of the following type should be defined.

```
Interface
  Function Func(X)
    USE NumTypes

    Real (kind=DP), Intent (in) :: X(:)
    Real (kind=DP) :: Func

  End Function Func
End Interface
```

**dmin:** Double precision real. Output. The lower limit of the confidence interval.

**dpls:** Double precision real. Output. The higher limit of the confidence interval.

### 9.21.3 Examples

Listing 9.21: Giving a confidence interval.

```
1 Program Tests
3   USE NumTypes
4   USE Error
5   USE Statistics
7   Integer, Parameter :: Nmax = 1000, Nb = 10000, Ndiv = 50
```

```

9   Real (kind=DP) :: Rdata(Nmax), Rmean(Nb), avg, Xmin, Xmax, h, Xac,&
      & err, dmin, dpls
11  Integer :: Id(Nmax), Ib(Nb, Nmax), Ntics(Ndiv)

13  Interface
      Function F(X)
      USE NumTypes

15      Real (kind=DP), Intent (in) :: X(:)
17      Real (kind=DP) :: F

19  End Function F
End Interface

21  CALL Normal(Rdata)

23  avg = Mean(Rdata)
25  ! Now create the resamples
27  CALL Bootstrap(Ib)

29  CALL EstBstrp(Rdata, Ib, F, avg, Err, Rmean)
Write(*,*)'#', avg, Err, Mean(Rdata)

31  CALL BstrpConfInt(Rdata, Ib, 0.1_DP, F, dmin, dpls)
Write(*,*)'Interval:', dmin, dpls
33  Write(*,*)Avg - Dmin, Dpls - Avg
Write(*,*)(dpls - dmin)/2.0_DP, 1.64485_DP/Sqrt(Real(Nmax,kind=DP))

35  Stop

37  End Program Tests

39  Function F(X)
41  USE NumTypes
43  USE Statistics

      Real (kind=DP), Intent (in) :: X(:)
45  Real (kind=DP) :: F

47  F = Mean(X)

49  Return
End Function F

```



# Ten

---

## MODULE Polynomial

---

This is the documentation of the `MODULE Polynomial`, a set of `FORTTRAN 90` routines to work with polynomials. This module make use of the `MODULE NumTypes`, `MODULE Constants`, `MODULE Error` and `MODULE Linear` so please read the documentation of these modules *before* reading this.

### 10.1 Type Pol

#### 10.1.1 Description

A new data type `Pol` is defined to work with polynomials. This type has two components: The coefficients of the polynomial, and the degree.

#### 10.1.2 Components

`Coef(:)`: Real double precision one dimensional array.

`dg`: Integer. The degree of the polynomial.

#### 10.1.3 Examples

A small example showing how to define a polynomial.

Listing 10.1: Defining a polynomial.

```
1 Program TestPoly
2
3   USE NumTypes
4   USE Error
5   USE Polynomial
6
7   Type (Pol) :: P1
8
9   Stop
10 End Program TestPoly
```

## 10.2 Type CmplxPol

### 10.2.1 Description

A new data type `CmplxPol` is defined to work with complex polynomials. This type has two components: The coefficients of the polynomial, and the degree.

All the routines, operators, etc... defined in this module works for complex as well as for real polynomials.

### 10.2.2 Components

`Coef(:)`: Complex double precision one dimensional array.

`dg`: Integer. The degree of the polynomial.

### 10.2.3 Examples

A small example showing how to define a polynomial of complex coefficients.

Listing 10.2: Defining a polynomial.

```

1 Program TestPoly
2
3   USE NumTypes
4   USE Error
5   USE Polynomial
6
7   Type (CmplxPol) :: P1
8
9   Stop
10 End Program TestPoly

```

## 10.3 Assignment

### 10.3.1 Description

You can directly assign one defined polynomial (complex or real) to another, or to an array of real numbers, that are interpreted as the coefficients.

### 10.3.2 Examples

Listing 10.3: Assigning polynomials.

```

1 Program TestPoly
2
3   USE NumTypes
4   USE Error
5   USE Polynomial
6
7   Integer, Parameter :: Deg = 4
8   Real (kind=DP) :: Hcoef(Deg+1)

```

```

10  Type (Pol) :: Hermite4
12  ! The fourth Hermite polynomial is  $x^4 - 6x^2 + 3$ , so
12  ! we first assign the values of the coefficients.
14  Hcoef = 0.0_DP
14  Hcoef(1) = 3.0_DP
16  Hcoef(3) = -6.0_DP
16  Hcoef(5) = 1.0_DP
18  Hermite4 = Hcoef
20  ! Now Show what we have in our data type:
22  Do I = 0, Hermite4%deg
22      Write(*, '(1I5,ES33.25)') I, Hermite4%Coef(I)
24  End Do
26  Stop
26  End Program TestPoly

```

## 10.4 Operator +

### 10.4.1 Description

You can naturally sum Pol or CmplxPol data types.

### 10.4.2 Examples

Listing 10.4: Adding polynomials.

```

Program TestPoly
2
3  USE NumTypes
4  USE Error
5  USE Polynomial
6
7  Integer, Parameter :: Deg = 4
8  Real (kind=DP) :: Hcoef(Deg+1)
9  Type (Pol) :: Hermite4, Hermite3, Sum
10
11  ! The Third Hermite polynomial is  $x^3 - 3x$ , so
12  ! we first assign the values of the coefficients.
14  Hcoef = 0.0_DP
14  Hcoef(2) = -3.0_DP
16  Hcoef(4) = 1.0_DP
18  Hermite3 = Hcoef(1:4)
20
21  ! The fourth Hermite polynomial is  $x^4 - 6x^2 + 3$ , so
22  ! we first assign the values of the coefficients.
24  Hcoef = 0.0_DP
24  Hcoef(1) = 3.0_DP

```

```

24   Hcoef(3) = -6.0_DP
      Hcoef(5) = 1.0_DP

26   Hermite4 = Hcoef

28   ! Now Add the two polynomials, and show the result.
      Sum = Hermite3 + Hermite4
30   Do I = 0, Sum%deg
      Write(*, '(1 I5, ES33.25) ') I, Sum%Coef(I)
32   End Do

34   Stop
End Program TestPoly

```

## 10.5 Operator -

### 10.5.1 Description

You can subtract Pol or CmplxPol data types.

### 10.5.2 Examples

Listing 10.5: Subtracting polynomials.

```

1  Program TestPoly

3   USE NumTypes
   USE Error
5   USE Polynomial

7   Integer, Parameter :: Deg = 4
   Real (kind=DP) :: Hcoef(Deg+1)
9   Type (Pol) :: Hermite4, Hermite3, Sum

11  ! The Third Hermite polynomial is  $x^3 - 3x$ , so
   ! we first assign the values of the coefficients.
13  Hcoef = 0.0_DP
   Hcoef(2) = -3.0_DP
15  Hcoef(4) = 1.0_DP

17  Hermite3 = Hcoef(1:4)

19  ! The fourth Hermite polynomial is  $x^4 - 6x^2 + 3$ , so
   ! we first assign the values of the coefficients.
21  Hcoef = 0.0_DP
   Hcoef(1) = 3.0_DP
23  Hcoef(3) = -6.0_DP
   Hcoef(5) = 1.0_DP

25  Hermite4 = Hcoef

27

```



```

29      ! Now Subtract the two polynomials, and show the result.
      Sum = Hermite3 - Hermite4
      Do I = 0, Sum%deg
31          Write(*, '(1 I5, ES33.25) ') I, Sum%Coef(I)
      End Do
33
      Stop
35 End Program TestPoly

```

## 10.6 Operator \*

### 10.6.1 Description

You can naturally multiply `Pol` data types, `Pol` data types with double precision real numbers, `CmplxPol` data types and `CmplxPol` data types with real or complex numbers.

### 10.6.2 Examples

Listing 10.6: Computing the product of two polynomials.

```

1 Program TestPoly
3     USE NumTypes
4     USE Error
5     USE Polynomial
7     Integer, Parameter :: Deg = 4
8     Real (kind=DP) :: Hcoef(Deg+1)
9     Type (Pol) :: Hermite4, Hermite3, Sum
11
12     ! The Third Hermite polynomial is  $x^3 - 3x$ , so
13     ! we first assign the values of the coefficients.
14     Hcoef = 0.0_DP
15     Hcoef(2) = -3.0_DP
16     Hcoef(4) = 1.0_DP
17
18     Hermite3 = Hcoef(1:4)
19
20     ! The fourth Hermite polynomial is  $x^4 - 6x^2 + 3$ , so
21     ! we first assign the values of the coefficients.
22     Hcoef = 0.0_DP
23     Hcoef(1) = 3.0_DP
24     Hcoef(3) = -6.0_DP
25     Hcoef(5) = 1.0_DP
26
27     Hermite4 = Hcoef
28
29     ! Now multiply the two polynomials, and show the result.
30     Sum = Hermite3 * Hermite4
31     Do I = 0, Sum%deg
        Write(*, '(1 I5, ES33.25) ') I, Sum%Coef(I)
    End Do

```

```

33     End Do
34     Stop
35 End Program TestPoly

```

## 10.7 Subroutine Init(P, Dgr)

### 10.7.1 Description

Allocate memory space for the coefficients of a `Pol` or a `CmplxPol` type.

### 10.7.2 Arguments

**P:** Type `Pol` or `CmplxPol`. The polynomial that you want to allocate space for.

**Dgr:** Integer. The degree of the polynomial.

### 10.7.3 Examples

Listing 10.7: Initialising a polynomial data type.

```

1 Program TestPoly
2
3     USE NumTypes
4     USE Error
5     USE Polynomial
6
7     Integer, Parameter :: Deg = 4
8     Real (kind=DP) :: Hcoef(Deg+1)
9     Type (Pol) :: Hermite4, Hermite3, Sum
10
11
12     ! An alternative way of setting the third Hermite
13     ! polynomial.
14     CALL Init(Hermite3, 3)
15     Hermite3%Coef(0) = 0.0_DP
16     Hermite3%Coef(1) = -3.0_DP
17     Hermite3%Coef(2) = 0.0_DP
18     Hermite3%Coef(3) = 1.0_DP
19     Hermite3%dg = 3
20
21
22     Stop
23 End Program TestPoly

```

## 10.8 Function Degree(P)

### 10.8.1 Description

Returns the degree of the polynomial P.

### 10.8.2 Arguments

P: Type Pol or CmplxPol. The polynomial whose degree we want to know.

### 10.8.3 Output

Integer. The degree of the polynomial P.

### 10.8.4 Examples

Listing 10.8: Returns the degree of a polynomial.

```

1 Program TestPoly
3   USE NumTypes
4   USE Error
5   USE Polynomial
7   Integer, Parameter :: Deg = 4
8   Real (kind=DP) :: Hcoef(Deg+1), X
9   Type (Pol) :: Hermite4, Hermite3, Sum
11
12   ! The Third Hermite polynomial is x^3 - 3x, so
13   ! we first assign the values of the coefficients.
14   Hcoef = 0.0_DP
15   Hcoef(2) = -3.0_DP
16   Hcoef(4) = 1.0_DP
17
18   Hermite3 = Hcoef(1:4)
19
20   ! The fourth Hermite polynomial is x^4 - 6x^2 + 3, so
21   ! we first assign the values of the coefficients.
22   Hcoef = 0.0_DP
23   Hcoef(1) = 3.0_DP
24   Hcoef(3) = -6.0_DP
25   Hcoef(5) = 1.0_DP
26
27   Hermite4 = Hcoef
28
29   ! Now Mutiply the two polynomials, and show the result.
30   Sum = Hermite3 * Hermite4
31
32   ! Show the degree of the product. It should be 4+3=7.
33   Write(*,*) Degree(Sum)
34
35   Stop
End Program TestPoly

```

## 10.9 Function Value(P, X)

### 10.9.1 Description

Computes the value of the polynomial P in the point X.

### 10.9.2 Arguments

**P:** Type `Pol` or `CmplxPol`. The polynomial.

**X:** Real double precision if P is of type `Pol` and Complex double precision if P is `CmplxPol`.  
The point in which you want to compute the value.

### 10.9.3 Output

Real double precision. The value of the polynomial P in the point X.

### 10.9.4 Examples

Listing 10.9: Computes the values of a polynomial at some points.

```

Program TestPoly
2
3   USE NumTypes
4   USE Error
5   USE Polynomial
6
7   Integer, Parameter :: Deg = 4
8   Real (kind=DP) :: Hcoef(Deg+1), X
9   Type (Pol) :: Hermite4, Hermite3, Sum
10
11   ! The Third Hermite polynomial is x^3 - 3x, so
12   ! we first assign the values of the coefficients.
13   Hcoef = 0.0_DP
14   Hcoef(2) = -3.0_DP
15   Hcoef(4) = 1.0_DP
16
17   Hermite3 = Hcoef(1:4)
18
19   ! The fourth Hermite polynomial is x^4 - 6x^2 + 3, so
20   ! we first assign the values of the coefficients.
21   Hcoef = 0.0_DP
22   Hcoef(1) = 3.0_DP
23   Hcoef(3) = -6.0_DP
24   Hcoef(5) = 1.0_DP
25
26   Hermite4 = Hcoef
27
28   ! Now Mutiply the two polynomials, and show the result.
29   Sum = Hermite3 * Hermite4
30
31   ! Compute the value of the product in some point in two

```

```

32  ! different ways.
    X = 9.34564_DP
34  Write(*, '(ES33.25) ') Value(Sum, X)
    Write(*, '(ES33.25) ') Value(Hermite3, X)*Value(Hermite4, X)
36
38  Stop
End Program TestPoly

```

## 10.10 Function Deriv(P)

### 10.10.1 Description

Computes the derivative of the polynomial P.

### 10.10.2 Arguments

P: Type Pol or CmplxPol. The polynomial whose derivative we want to compute.

### 10.10.3 Output

Type Pol. Another polynomial: the derivative of P.

### 10.10.4 Examples

Listing 10.10: Computing the derivative of a polynomial.

```

1  Program TestPoly
3
   USE NumTypes
   USE Error
   USE Polynomial
5
7  Integer, Parameter :: Deg = 4
   Real (kind=DP) :: Hcoef(Deg+1), X
9  Type (Pol) :: Hermite4, Hermite3, Res, Sum
11
   ! The Third Hermite polynomial is x^3 - 3x, so
   ! we first assign the values of the coefficients.
13  Hcoef = 0.0_DP
   Hcoef(2) = -3.0_DP
15  Hcoef(4) = 1.0_DP
17
   Hermite3 = Hcoef(1:4)
19
   ! The fourth Hermite polynomial is x^4 - 6x^2 + 3, so
   ! we first assign the values of the coefficients.
21  Hcoef = 0.0_DP
   Hcoef(1) = 3.0_DP
23  Hcoef(3) = -6.0_DP
   Hcoef(5) = 1.0_DP

```

```

25 Hermite4 = Hcoef
27
28 ! Now compute the derivative of Hermite4
29 Res = Deriv(Hermite4)
30
31 ! From the recursion relation of the Hermite polynomials
32 ! we should obtain twice the same number:
33 X = 7.346582_DP
34 Write(*, '(ES33.25) ') Value(Res, X)
35 Write(*, '(ES33.25) ') 4.0_DP*Value(Hermite3, X)
36
37 Stop
38 End Program TestPoly

```

## 10.11 Function Integra(P[, Cte])

### 10.11.1 Description

Computes the integral of the polynomial P. If Cte is present then it is used as *integration constant*.

### 10.11.2 Arguments

**P:** Type Pol or CmplxPol. The polynomial whose integral we want to compute.

**Cte:** Optional. Real double precision if P is of type Pol and Complex double precision if P is CmplxPol. If not present, the default value is 0.

### 10.11.3 Output

Type Pol. Another polynomial: the integral of P.

### 10.11.4 Examples

Listing 10.11: Computing the integral of a polynomial.

```

1 Program TestPoly
2
3   USE NumTypes
4   USE Error
5   USE Polynomial
6
7   Integer, Parameter :: Deg = 4
8   Real (kind=DP) :: Hcoef(Deg+1), X
9   Type (Pol) :: Hermite4, Hermite3, Res, Sum
10
11 ! The Third Hermite polynomial is x^3 - 3x, so
12 ! we first assign the values of the coefficients.
13 Hcoef = 0.0_DP

```

```

15   Hcoef(2) = -3.0_DP
    Hcoef(4) = 1.0_DP

17   Hermite3 = Hcoef(1:4)

19   ! The fourth Hermite polynomial is  $x^4 - 6x^2 + 3$ , so
    ! we first assign the values of the coefficients.
21   Hcoef = 0.0_DP
    Hcoef(1) = 3.0_DP
23   Hcoef(3) = -6.0_DP
    Hcoef(5) = 1.0_DP

25   Hermite4 = Hcoef

27   ! Now compute the derivative of Hermite4
29   Res = Integra(Hermite3, 3.0_DP/4.0_DP)

31   ! From the recursion relation of the Hermite polynomials
    ! we should obtain twice the same number:
33   X = 7.346582_DP
    Write(*, '(ES33.25) ') Value(Res, X)
35   Write(*, '(ES33.25) ') 0.25_DP*Value(Hermite4, X)

37
39   Stop
End Program TestPoly

```

## 10.12 Function InterpolValue(X, Y, Xo)

### 10.12.1 Description

Computes the value of the interpolation polynomial that pass trough  $(X(:), Y(:))$  in the point  $Xo$ .

### 10.12.2 Arguments

$X(:)$ ,  $Y(:)$ : Real double precision or Complex double precision one dimensional arrays. Specify the points at which the interpolation polynomial should pass.

$Xo$ : Same type as  $X(:)$  and  $Y(:)$ . The point at which you want to compute the interpolation polynomial.

### 10.12.3 Output

Real double precision if input is real and complex double precision if input is complex. The value of the interpolation polynomial in  $Xo$ .

### 10.12.4 Examples

Listing 10.12: Compute values of the Interpolation polynomial.

```

1  Program TestPoly
3      USE NumTypes
4      USE Error
5      USE Polynomial
7
8      Integer, Parameter :: Deg = 4, Np = 7
9      Real (kind=DP) :: Hcoef(Deg+1), X, Xp(Np), Yp(Np)
10     Type (Pol) :: Hermite4, Hermite3, Res, Sum
11
12     CALL Random_Number(Xp)
13     Yp = 3.347234_DP*Xp - 2.475875_DP*Xp**3 - 7.23467_DP*Xp**4 + &
14         & 1.47854_DP*Xp**6
15
16     ! Now we compute the value of the interpolation polynomial
17     ! at X, and compare it with the real value of the Polynomial
18     X = -1.23899843_DP
19     Write(*, '(ES33.25)') InterpolValue(Xp, Yp, X)
20     Write(*, '(ES33.25)') 3.347234_DP*X - 2.475875_DP*X**3 - &
21         & 7.23467_DP*X**4 + 1.47854_DP*X**6
22
23     Stop
24 End Program TestPoly

```

## 10.13 Function Interpol(X, Y)

Computes the interpolation polynomial that pass trough  $(X(:), Y(:))$ . **Note that using this function can be very unstable.**

### 10.13.1 Arguments

$X(:)$ ,  $Y(:)$ : Real double precision or Complex double precision one dimensional arrays. Specify the points at which the interpolation polynomial should pass.

### 10.13.2 Output

Type `Pol` if input is real, and `CmplxPol` if input is complex. The interpolation polynomial.

### 10.13.3 Examples

Listing 10.13: Computes the interpolation polynomial.

```

1  Program TestPoly
3      USE NumTypes
4      USE Error
5      USE Polynomial

```



```

7  Integer, Parameter :: Deg = 4, Np = 7
   Real (kind=DP) :: Hcoef(Deg+1), X, Xp(Np), Yp(Np)
9  Type (Pol) :: Hermite4, Hermite3, Res, Sum

11
   CALL Random_Number(Xp)
13  Yp = 3.347234_DP*Xp - 2.475875_DP*Xp**3 - 7.23467_DP*Xp**4 + &
      & 1.47854_DP*Xp**6

15
   ! Now we compute the interpolation polynomial
17  ! at X, and compare it with the real value of the Polynomial
   X = -1.23899843_DP
19  Res = Interpol(Xp,Yp)
   Write(*,'(ES33.25)') Value(Res, X)
21  Write(*,'(ES33.25)') 3.347234_DP*X - 2.475875_DP*X**3 - &
      & 7.23467_DP*X**4 + 1.47854_DP*X**6

23
25  Stop
End Program TestPoly

```

## 10.14 Subroutine Spline(X, Y, Ypp0, YppN, Pols)

### 10.14.1 Description

Compute the cubic spline interpolation polynomial that pass trough  $(X(:), Y(:))$ .

### 10.14.2 Arguments

$X(:), Y(:)$ : Real double precision one dimensional arrays. Specify the points at which the cubic spline interpolation polynomial should pass.

$Ypp0, YppN$ : The values of the second derivatives of the cubic spline interpolation polynomial in the first and last points.

$Pols(:)$ : Type Pol one dimensional array. Returns the N-1 cubic interpolation polynomials.

### 10.14.3 Examples

Listing 10.14: Computes the cubic spline interpolation polynomial.

```

Program TestPoly

2  USE NumTypes
   USE Error
4  USE Polynomial
   USE NonNumeric

6

8  Integer, Parameter :: Deg = 4, Np = 7
   Real (kind=DP) :: Hcoef(Deg+1), X, Xp(Np), Yp(Np)

```

```

10  Type (Pol) :: Hermite4, Hermite3, Res, Sum, Spl(Np-1)

12

13  CALL Random_Number(Xp)
14  ! Order Xp
15  CALL Qsort(Xp)
16  Yp = 3.347234_DP*Xp - 2.475875_DP*Xp**3 - 7.23467_DP*Xp**4 + &
      & 1.47854_DP*Xp**6

18

19  ! Now we compute the interpolation polynomial
20  ! at X, and compare it with the real value of the Polynomial, and
21  ! the value of the spline cubic interpolation polynomial.
22  X = 0.23899843_DP
23  Res = Interpol(Xp,Yp)
24  CALL Spline(Xp, Yp, 0.0_DP, 0.0_DP, Spl)
25  Write(*,'(ES33.25)') Value(Res, X)
26  Write(*,'(ES33.25)') Value(Spl(Locate(Xp, X)), X)
27  Write(*,'(ES33.25)') 3.347234_DP*X - 2.475875_DP*X**3 - &
28      & 7.23467_DP*X**4 + 1.47854_DP*X**6

30

31  Stop
32 End Program TestPoly

```

# Eleven

---

## MODULE Root

---

This is the documentation of the `MODULE Root`, a set of `FORTRAN 90` routines to compute roots of functions. This module make use of the `MODULE NumTypes`, `MODULE Constants` and `MODULE Error` so please read the documentation of these modules *before* reading this.

### 11.1 Subroutine `RootPol(a, b, [c, d], z1, z2, [z3, z4])`

#### 11.1.1 Description

Returns the complex roots of a polynomial of degree 2, 3 or 4.

#### 11.1.2 Arguments

**a, b, c, d:** The coefficients of the polynomial. The meaning of the coefficieents **a,b,c,d** depends on the degree of the polynomial:

$$\begin{aligned}P(x) &= x^2 + ax + b \\P(x) &= x^3 + ax^2 + bx + c \\P(x) &= x^4 + ax^3 + bx^2 + cx + d\end{aligned}$$

**z1,z2,z3,z4:** Complex simple or double precision. The roots of the polynomial.

#### 11.1.3 Examples

Listing 11.1: Computing roots of polynomials.

```
2 Program TestRoot
3
4   USE NumTypes
5   USE Error
6   USE Root
7
8   Real (kind=DP) :: a, b, c, d
9   Complex (kind=DPC) :: z1, z2, z3, z4, ac, bc, cc, dc
```

```

10  CALL Random_Number(a)
12  CALL Random_Number(b)
14  CALL Random_Number(c)
14  CALL Random_Number(d)
14  CALL RootPol(a,b,z1,z2)
16  Write(*,'(3ES20.12)') Z1, Abs(z1**2 + a*z1 + Cmplx(b,kind=DPC))
16  Write(*,'(3ES20.12)') Z2, Abs(z2**2 + a*z2 + Cmplx(b,kind=DPC))
18
18  CALL RootPol(a,b,c,z1,z2,z3)
20  Write(*,*)
20  Write(*,'(3ES20.12)') Z1, Abs(z1**3+a*z1**2+b*z1+Cmplx(c,kind=DPC))
22  Write(*,'(3ES20.12)') Z2, Abs(z2**3+a*z2**2+b*z2+Cmplx(c,kind=DPC))
22  Write(*,'(3ES20.12)') Z3, Abs(z3**3+a*z3**2+b*z3+Cmplx(c,kind=DPC))
24
24  ac = Cmplx(a,kind=DPC)
26  bc = Cmplx(b,a,kind=DPC)
26  cc = Cmplx(c,kind=DPC)
28  dc = Cmplx(d,kind=DPC)
28  CALL RootPol(ac,bc,z1,z2)
30  Write(*,*)
30  Write(*,'(3ES20.12)') Z1, Abs(z1**2 + ac*z1 + Cmplx(bc,kind=DPC))
32  Write(*,'(3ES20.12)') Z2, Abs(z2**2 + ac*z2 + Cmplx(bc,kind=DPC))
32  CALL RootPol(ac,bc,cc,dc,z1,z2,z3,z4)
34  Write(*,*)
34  Write(*,'(3ES20.12)') Z1, Abs(z1**4+ac*z1**3+bc*z1**2+cc*z1+dc)
36  Write(*,'(3ES20.12)') Z2, Abs(z2**4+ac*z2**3+bc*z2**2+cc*z2+dc)
36  Write(*,'(3ES20.12)') Z3, Abs(z3**4+ac*z3**3+bc*z3**2+cc*z3+dc)
38  Write(*,'(3ES20.12)') Z4, Abs(z4**4+ac*z4**3+bc*z4**2+cc*z4+dc)
40
40  Stop
42  End Program TestRoot

```

## 11.2 Function Newton( $X_0$ , $F_{new}$ , [Tol])

### 11.2.1 Description

Compute a root of the function defined by the routine **Fnew**.

### 11.2.2 Arguments

**Xo**: Real simple or double precision. An initial guess of the position of the root.

**Fnew**: The function whose root we want to compute. It is defined as a subroutine that returns the value of the function and of its derivative. If it is an external function, an interface block like this should be defined

```

Interface
  Subroutine FNew( $X_0$ , F, D)

```

```

USE NumTypes

Real (kind=DP), Intent (in) :: Xo
Real (kind=DP), Intent (out) :: F, D
End Subroutine FNew
End Interface

```

where F is the value of the function in Xo, and D the value of the derivative in Xo. If the arguments are of simple precision, a similar interface should be provided, where the arguments of Fnew are of single precision.

**Tol:** Real single or double precision. Optional. An estimation of the desired accuracy of the position of the root.

### 11.2.3 Output

Real single or double precision. The position of the root.

### 11.2.4 Examples

Listing 11.2: Computing roots of non-linear functions with the Newton method.

```

Program TestRoot
2
  USE NumTypes
4  USE Error
  USE Root
6
  Real (kind=DP) :: a, b, c, d, X
8  Complex (kind=DPC) :: z1, z2, z3, z4, ac, bc, cc, dc
10
11  Interface
12    Subroutine FNew(Xo, F, D)
14      USE NumTypes
16      Real (kind=DP), Intent (in) :: Xo
17      Real (kind=DP), Intent (out) :: F, D
18    End Subroutine FNew
19  End Interface
20
21  ! Compute the value such that cos(x) = x
22  X = Newton(0.0_DP, Fnew, 1.0E-10_DP)
24  Write(*, '(1A,ES33.25)') 'Point: ', X
25  Write(*, '(1A,ES33.25)') 'Value of Cos: ', Cos(X)
26
28  Stop

```

```

End Program TestRoot
30
! *****
32 ! *
Subroutine FNew(Xo, F, D)
34 ! *
! *****
36
USE NumTypes
38
Real (kind=DP), Intent (in) :: Xo
40 Real (kind=DP), Intent (out) :: F, D
42
F = Xo - Cos(Xo)
44 D = 1.0_DP + Sin(Xo)
46
Return
End Subroutine FNew

```

## 11.3 Function Bisec(a, b, Fbis, [Tol])

### 11.3.1 Description

Compute the root of the function defined by **Fbis**.

### 11.3.2 Arguments

**a, b:** Real single or double precision. Initial points, such that  $\text{Fbis}(a)\text{Fbis}(b) < 0$ .

**Fbis:** The function whose root we want to compute. It is defined as a function that returns the value of the function. If it is an external function, an interface block like this should be defined

```

Interface
  Function F(X)

      USE NumTypes

      Real (kind=DP), Intent (in) :: X
      Real (kind=DP) :: F
  End Function F
End Interface

```

where **F** is the value of the function in **X**. If the arguments are of simple precision, a similar interface should be provided, where the arguments of **F** are of single precision.

**Tol:** Real single or double precision. Optional. An estimation of the desired accuracy of the position of the root.

### 11.3.3 Output

Real single or double precision. The position of the root of Fbis.

### 11.3.4 Examples

Listing 11.3: Computing roots with the bisection method.

```

1 Program TestRoot
3   USE NumTypes
4   USE Error
5   USE Root
7   Real (kind=DP) :: a, b, c, d, X
8   Complex (kind=DPC) :: z1, z2, z3, z4, ac, bc, cc, dc
9
10  Interface
11    Function Fbis(X)
12
13      USE NumTypes
14
15      Real (kind=DP), Intent (in) :: X
16      Real (kind=DP) :: Fbis
17    End Function Fbis
18  End Interface
19
20  ! Compute the value such that cos(x) = x
21  X = Bisec(0.0_DP, 1.1_DP, Fbis, 1.0E-10_DP)
22  Write(*, '(1A,ES33.25)') 'Point: ', X
23  Write(*, '(1A,ES33.25)') 'Value of Cos: ', Cos(X)
24
25
26  Stop
27 End Program TestRoot
28
29 ! *****
30 ! *
31 Function FBis(X)
32 ! *
33 ! *****
34
35  USE NumTypes
36
37  Real (kind=DP), Intent (in) :: X
38  Real (kind=DP) :: FBis
39
40
41  FBis = X - Cos(X)
42
43  Return
44 End Function FBis

```





# Twelve

---

## MODULE Fourier

---

This is the documentation of the `MODULE Fourier`, a set of `FORTRAN 90` routines to work with Fourier series. This module make use of the `MODULE NumTypes` and the `MODULE Constants` so please read the documentation of these modules *before* reading this.

### 12.1 Type `Fourier_Serie`

#### 12.1.1 Description

A new data type `Fourier_Serie` is defined to work with Fourier series. This type has two components: The modes, and the number of modes.

#### 12.1.2 Components

`Coef(:)`: Complex double precision one dimensional array. The modes.

`Nterm`: Integer. The number of terms of the Fourier series.

#### 12.1.3 Examples

A small example showing how to define a Fourier serie.

Listing 12.1: Defining a Fourier serie.

```
2  Program TestFourier
4      USE NumTypes
4      USE Constants
4      USE Fourier
6
6      Type (Fourier_Serie) :: Ff
8
8      Stop
10 End Program TestPoly
```

## 12.2 Type Fourier\_Serie\_2D

### 12.2.1 Description

A new data type `Fourier_Serie_2D` is defined to work with two dimensional Fourier series. This type has two components: The modes, and the number of modes.

### 12.2.2 Components

`Coef(:, :)`: Complex double precision two dimensional array. The modes.

`Nterm`: Integer. The number of terms of the Fourier series.

### 12.2.3 Examples

A small example showing how to define a polynomial.

Listing 12.2: Defining a two-dimensional Fourier serie.

```

1 Program TestFourier
2
3     USE NumTypes
4     USE Constants
5     USE Fourier
6
7     Type (Fourier_Serie_2D) :: Ff
8
9     Stop
10 End Program TestPoly

```

## 12.3 Assignment

### 12.3.1 Description

You can directly assign one defined Fourier series (one or two dimensional) to another.

### 12.3.2 Examples

This example uses the `Init_Serie` subroutine. For details of the usage of this function look at the section (12.8), page (100).

Listing 12.3: Assigning Fourier series.

```

1 Program TestFourier
2
3     USE NumTypes
4     USE Constants
5     USE Fourier
6
7     Type (Fourier_Serie) :: FS1, FS2
8
9     CALL Init_Serie(FS1, 20)
10    CALL Init_Serie(FS2, 20)

```

```

12  FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
    FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
14
16  FS2 = FS1
18  Write(*, '(2ES33.25) ') FS2%Coef( 1)
    Write(*, '(2ES33.25) ') FS2%Coef(-1)
20
    Stop
End Program TestFourier

```

## 12.4 Operator +

### 12.4.1 Description

You can naturally sum one or two dimensional Fourier series. If they have different sizes, it is assumed that the non defined modes of the short Fourier Series are zero.

### 12.4.2 Examples

This example uses the `Init_Serie` subroutine. For details of the usage of this function look at the section (12.8), page (100).

Listing 12.4: Adding Fourier series.

```

1  Program TestFourier
3
4  USE NumTypes
5  USE Constants
6  USE Fourier
7
8  Type (Fourier_Serie_2D) :: FS1, FS2, FS3
9  Integer :: Nt
10
11 Nt = 4
12 CALL Init_Serie(FS1, Nt)
13 CALL Init_Serie(FS2, Nt)
14
15 FS1%Coef( 1,1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
16 FS1%Coef(-1,1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
17
18 FS2%Coef( 1,1) = Cmplx(-1.0_DP, 4.5_DP, kind=DPC)
19 FS2%Coef(-1,1) = Cmplx(-1.0_DP, -6.78745_DP, kind=DPC)
20
21 FS3 = FS1 + FS2
22 Write(*, '(2ES33.25) ') FS3%Coef( 1,1)
23 Write(*, '(2ES33.25) ') FS3%Coef(-1,1)
24
25 Stop
End Program TestFourier

```

## 12.5 Operator -

### 12.5.1 Description

You can naturally subtract one or two dimensional Fourier series. If they have different sizes, it is assumed that the non defined modes of the short Fourier Series are zero.

### 12.5.2 Examples

Listing 12.5: Subtracting Fourier series.

```

Program TestFourier
2
  USE NumTypes
4  USE Constants
  USE Fourier
6
  Type (Fourier_Serie) :: FS1, FS2, FS3
8  Integer :: Nt
10
  Nt = 4
  CALL Init_Serie(FS1, Nt)
12
  FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
14  FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
16
  FS2 = FS1
18
  FS3 = FS1 - FS2
  Write(*, '(2ES33.25)') FS3%Coef( 1)
20  Write(*, '(2ES33.25)') FS3%Coef(-1)
22
  Stop
End Program TestFourier

```

## 12.6 Operator \*

### 12.6.1 Description

You can naturally multiply one or two dimensional Fourier series, in which case the convolution of the Fourier Modes is performed. If they have different sizes, it is assumed that the non defined modes of the short Fourier Series are zero.

### 12.6.2 Examples

Listing 12.6: Computing the convolution of Fourier series.

```

1 Program TestFourier
3
  USE NumTypes
  USE Constants

```

```

5  USE Fourier
7  Type (Fourier_Serie) :: FS1, FS2, FS3
   Integer :: Nt
9
   Nt = 4
11 CALL Init_Serie(FS1, Nt)
13 FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
   FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
15
   FS2 = FS1
17
   FS3 = FS1 * FS2
19 Write(*, '(2ES33.25)') FS3%Coef( 0)
21
   Stop
End Program TestFourier

```

## 12.7 Operator \*\*

### 12.7.1 Description

You can naturally compute the integer power of a one or two dimensional Fourier series, in which case the convolution of the Fourier modes with themselves are performed a certain number of times.

### 12.7.2 Examples

Listing 12.7: "Exponentiating" Fourier series.

```

Program TestFourier
2
   USE NumTypes
   USE Constants
   USE Fourier
4
6
   Type (Fourier_Serie) :: FS1, FS2, FS3
   Integer :: Nt
8
   Nt = 4
   CALL Init_Serie(FS1, Nt)
12  CALL Init_Serie(FS2, Nt)
   CALL Init_Serie(FS3, Nt)
14
   FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
16  FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
18
   FS3%Coef(0) = Cmplx(1.0_DP, 0.0_DP, kind=DPC)
20
   FS2 = FS1**8

```

```

22      Do I = 1, 8
          FS3 = FS3 * FS1
24      End Do

26      Write(*, '(2ES33.25) ') FS2%Coef( 0)
      Write(*, '(2ES33.25) ') FS3%Coef( 0)

28      Stop
End Program TestFourier

```

## 12.8 Subroutine Init\_Serie(FS,Ns)

### 12.8.1 Description

Allocate memory space for the modes of a one or two dimensional Fourier series.

### 12.8.2 Arguments

**FS:** Type `Fourier_Serie` or type `Fourier_Serie_2D`. The Fourier series that you want to allocate space for.

**Ns:** Integer. The number of modes.

### 12.8.3 Examples

Any of the examples of some of the previous sections are also good examples of the use of the `Init_Serie` subroutine. Here we simply repeat one of them.

Listing 12.8: Initialising a Fourier series.

```

1  Program TestFourier

3      USE NumTypes
      USE Constants
5      USE Fourier

7      Type (Fourier_Serie) :: FS1, FS2, FS3
      Integer :: Nt

9      Nt = 4
11     CALL Init_Serie(FS1, Nt)

13     FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
      FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
15

17     FS2 = FS1

      FS3 = FS1 * FS2
19     Write(*, '(2ES33.25) ') FS3%Coef( 0)

21     Stop
End Program TestFourier

```

## 12.9 Function Eval\_Serie(FS, X, [Y], Tx, [Ty])

### 12.9.1 Description

Compute the value of the Fourier series FS with periods Tx,Ty at the point X,Y.

### 12.9.2 Arguments

**FS:** Type `Fourier_Serie` or type `Fourier_Serie_2D`. The Fourier series that you want to evaluate.

**X,Y:** Real double precision. The point in which you want to evaluate the Fourier series. If FS is a two dimensional Fourier series, then Y must be present.

**Tx,Ty:** Real double precision. The period(s). If FS is a two dimensional Fourier series, then Ty must be present.

### 12.9.3 Output

Real double precision. The value of the function defined by the modes in FS at the point (X[,Y]).

### 12.9.4 Examples

Listing 12.9: Evaluating a Fourier series at a point.

```

1 Program TestFourier
2
3   USE NumTypes
4   USE Constants
5   USE Fourier
6
7   Type (Fourier_Serie) :: FS1, FS2, FS3
8   Integer :: Nt
9
10  Nt = 4
11  CALL Init_Serie(FS1, Nt)
12  CALL Init_Serie(FS2, Nt)
13  CALL Init_Serie(FS3, Nt)
14
15
16  FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
17  FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
18
19  FS2 = FS1**2
20
21  FS3 = FS1*FS2
22
23  Write(*, '(2 ES33.25) ') Eval_Serie(FS1,0.12_DP,1.0_DP) * &
24                               & Eval_Serie(FS2,0.12_DP,1.0_DP)
25  Write(*, '(2 ES33.25) ') Eval_Serie(FS3,0.12_DP,1.0_DP)
26

```

```

28      Stop
      End Program TestFourier

```

## 12.10 Function Unit(FS, Ns)

### 12.10.1 Description

Allocate memory space for the modes of a one or two dimensional Fourier series and sets the zero mode equal to 1.

### 12.10.2 Arguments

**FS:** Type `Fourier_Serie` or type `Fourier_Serie_2D`. The Fourier series that you want to allocate space for.

**Ns:** Integer. The number of modes.

### 12.10.3 Examples

Listing 12.10: Obtaining a constant Fourier series.

```

Program TestFourier
2
  USE NumTypes
4  USE Constants
  USE Fourier
6
  Type (Fourier_Serie) :: FS1, FS2, FS3
8  Integer :: Nt
10
  Nt = 4
  CALL Init_Serie(FS1, Nt)
12  CALL Init_Serie(FS2, Nt)
  CALL Init_Serie(FS3, Nt)
14
16  FS1%Coef( 1) = Cmplx(1.0_DP, 0.5_DP, kind=DPC)
  FS1%Coef(-1) = Cmplx(1.0_DP, 0.7_DP, kind=DPC)
18
  CALL Unit(FS2, Nt)
20
  FS3 = FS1*FS2
22
  Write(*, '(2ES33.25)') Eval_Serie(FS1, 0.12_DP, 1.0_DP)
24  Write(*, '(2ES33.25)') Eval_Serie(FS3, 0.12_DP, 1.0_DP)
26
  Stop
End Program TestFourier

```



## 12.11 Function DFT(Data, Is)

### 12.11.1 Description

Compute the Discrete Fourier Transform of the values stored in the complex array **Data**. If **Is** is present and is set to -1, the inverse Discrete Fourier Transform is performed. The direct Fourier transform is defined as

$$\tilde{f}(k) = \sum_{n=0}^N f_n e^{\frac{2\pi i n k}{N}} \quad \forall k \in \left[-\frac{N}{2}, \frac{N}{2}\right]$$

the inverse one is defined as

$$\tilde{f}(k) = \frac{1}{N} \sum_{n=0}^N f_n e^{\frac{-2\pi i n k}{N}} \quad \forall k \in \left[-\frac{N}{2}, \frac{N}{2}\right]$$

### 12.11.2 Arguments

**Data(:, :)**: One or two dimensional double precision complex array. The data whose Discrete Fourier Transform we want to compute.

**Is**: Integer. Optional. A flag to tell if we want to compute the direct or the inverse Fourier transform.

### 12.11.3 Output

Type **Fourier\_Serie** if **Data(:)** is one dimensional, and type **Fourier\_Serie\_2D** if **Data(:, :)** is two dimensional.

### 12.11.4 Examples

This example compute the discrete Fourier transform of  $f(x_i) = \sin(x_i)$ .

Listing 12.11: Computing the Discrete Fourier Transform.

```

1 Program TestFourier
3   USE NumTypes
3   USE Constants
5   USE Fourier
7   Integer, Parameter :: Nmax=20
7   Type (Fourier_Serie) :: FS1, FS2, FS3
9   Complex (kind=DPC) :: Data(Nmax), X
9   Integer :: Nt
11
11  Do I = 1, Nmax
13    X = Cmplx(TWOPLDP*I/Nmax)
13    Data(I) = Sin(X)
15  End Do
17  FS1 = DFT(Data)

```

```

19  Write(*, '(1A,2 ES33.25) ') 'Mode k= 1: ', FS1%Coef( 1)
21  Write(*, '(1A,2 ES33.25) ') 'Mode k=-1: ', FS1%Coef(-1)
21  Write(*, '(ES33.25) ') Sum(Abs(FS1%Coef(:)))
23  Stop
End Program TestFourier

```

## 12.12 Function Conjg(FS)

### 12.12.1 Description

Computes the Fourier modes that correspond to the conjugate function. This means: If the modes of FS are  $\tilde{f}(k)$ , this function returns a Fourier series with modes  $\tilde{f}(-k)$ .

### 12.12.2 Arguments

**FS:** Type `Fourier_Serie` or type `Fourier_Serie_2D`. The Fourier series whose conjugate you want to compute.

### 12.12.3 Output

Type `Fourier_Serie` if FS is of type `Fourier_Serie`, and type `Fourier_Serie_2D` if FS is of Type `Fourier_Serie_2D`.

### 12.12.4 Examples

Listing 12.12: Computing the Conjugate Fourier Series.

```

Program TestFourier
2
  USE NumTypes
4  USE Constants
  USE Fourier
6
  Integer, Parameter :: Nmax=20
8  Type (Fourier_Serie) :: FS1, FS2, FS3
  Complex (kind=DPC) :: Data(Nmax), X
10 Integer :: Nt
12
  Do I = 1, Nmax
    X = Cmplx(TWOPLDP*I/Nmax, kind=DPC)
14    Data(I) = Sin(X) + Cmplx(0.0_DP, I*2.0_DP, kind=DPC)
  End Do
16
  FS1 = DFT(Data)
18
  Write(*, '(2 ES33.25) ') Eval_Serie(FS1, 0.23_DP, 1.0_DP)
20  Write(*, '(2 ES33.25) ') Eval_Serie(Conjg(FS1), 0.23_DP, 1.0_DP)

```

```

22      Stop
24 End Program TestFourier

```

## 12.13 Subroutine Save\_Serie(FS, File)

### 12.13.1 Description

Write the Fourier series FS to the file File.

### 12.13.2 Arguments

**FS:** Type Fourier\_Serie or type Fourier\_Serie\_2D. The Fourier series that you want to store in a file.

**File:** Character string of arbitrary length. The name of the file in which you want to save FS.

### 12.13.3 Examples

Listing 12.13: Saving a Fourier Serie in a file.

```

Program TestFourier
2
  USE NumTypes
4  USE Constants
  USE Fourier
6
  Integer, Parameter :: Nmax=20
8  Type (Fourier_Serie) :: FS1, FS2, FS3
  Complex (kind=DPC) :: Data(Nmax), X
10 Integer :: Nt
12
  Do I = 1, Nmax
    X = Cmplx(TWOPLDP*I/Nmax, kind=DPC)
14    Data(I) = Sin(X) + Cmplx(0.0_DP, I*2.0_DP, kind=DPC)
  End Do
16
  FS1 = DFT(Data)
18
  CALL Save(FS1, 'datamodes.dat')
20
  Stop
22 End Program TestFourier

```

## 12.14 Subroutine Read\_Serie(FS, File)

### 12.14.1 Description

Reads the Fourier series FS stored in the file File.

### 12.14.2 Arguments

**FS:** Type `Fourier_Serie` or type `Fourier_Serie_2D`. The name of the Fourier series data type in which you want to store that data.

**File:** Character string of arbitrary length. The name of the file in which the saved series is.

### 12.14.3 Examples

Listing 12.14: Reading a Fourier series from a file.

```

Program TestFourier
2
  USE NumTypes
4  USE Constants
  USE Fourier
6
  Integer, Parameter :: Nmax=20
8  Type (Fourier_Serie) :: FS1, FS2, FS3
  Complex (kind=DPC) :: Data(Nmax), X
10 Integer :: Nt
12
  Do I = 1, Nmax
    X = Cmplx(TWOPLDP*I/Nmax, kind=DPC)
14    Data(I) = Sin(X) + Cmplx(0.0_DP, I*2.0_DP, kind=DPC)
  End Do
16
  FS1 = DFT(Data)
18
  CALL Save_Serie(FS1, 'datamodes.dat')
20  CALL Read_Serie(FS2, 'datamodes.dat')
22
  Write(*, '(ES33.25)') Sum(Abs(FS1%Coef(:) - FS2%Coef(:)))
24
  Stop
26 End Program TestFourier

```

# Thirteen

---

## MODULE Time

---

The `MODULE Time` is a module to provide access to date and time properties.

### 13.1 Type `tm`

#### 13.1.1 Description

A new data type, called `tm` is defined. It has some properties common with the same derived type defined in the C standard library. The components of the type specify a time: Day, year, month, hour, etc. . .

#### 13.1.2 Components

`hour`: Integer. Hour of the day [0-23].  
`min`: Integer, Minutes after the hour [0-59].  
`sec`: Integer. Seconds after the minute [0-59].  
`msec`: Integer. Milliseconds after the second [0-999].  
`year`: Integer. Year.  
`month`: Integer. Month of the year [0-11].  
`mday`: Integer. Day of the month [1-31].  
`wday`: Integer. Day of the week since Sunday [0-6].

#### 13.1.3 Example

A small example defining a `tm` data type.

Listing 13.1: Defining a Time data type.

```
2 Program Test
4   USE NumTypes
   USE Time
```

```
6   Type (tm) :: Oneday
8   OneDay%hour = 12
9   OneDay%min  = 0
10  OneDay%sec  = 0
11  OneDay%mday = 10
12  OneDay%mon  = 0
13  OneDay%year = 2007
14  OneDay%yday = 3
16  Stop
End Program Test
```

## 13.2 Function `gettime()`

### 13.2.1 Description

The function `gettime()` returns the current time and date in a `type tm` data type.

### 13.2.2 Arguments

This function has no arguments.

### 13.2.3 Output

Type `tm`, containing all the information about the date and time.

### 13.2.4 Example

A small program that prints the current year.

Listing 13.2: Obtaining the current date and time.

```
1 Program Test
3   USE NumTypes
4   USE Time
5
6   Type (tm) :: Oneday
7
8   Oneday = gettime()
9
10  Write(*,*) 'Current year: ', Oneday%year
11
12  Stop
13 End Program Test
```

## 13.3 Function isleap(Nyr)

### 13.3.1 Description

The function `isleap(Nyr)` returns `.true.` if `Nyr` is a leap year, and `.false.` otherwise. Note that the leap years are different in the Julian and Gregorian calendars. In this code the Gregorian calendar is supposed valid *after* 1582<sup>1</sup>.

### 13.3.2 Arguments

**Nyr:** Integer. The year.

### 13.3.3 Output

Logical. `.true.` if `Nyr` is a leap year, and `.false.` otherwise.

### 13.3.4 Example

A small program that tell us if the current year is leap.

Listing 13.3: Are we in a leap year?.

```

1 Program Test
3   USE NumTypes
4   USE Time
5
6   Type (tm) :: Oneday
7
8   Oneday = gettime()
9
10  If (isleap(Oneday%year)) Then
11    Write(*,*) 'We are in a leap year.'
12  Else
13    Write(*,*) 'We are not in a leap year.'
14  End If
15
16  Stop
17 End Program Test

```

## 13.4 Function asctime(t)

### 13.4.1 Description

The function `asctime`, returns a 24 length character string from a type `tm` data type, containing the date and time, in a similar way that the function `asctime` of the C standard library, for example:

<sup>1</sup>For more details, take a look at

[http://en.wikipedia.org/wiki/Gregorian\\_calendar](http://en.wikipedia.org/wiki/Gregorian_calendar)

Wed Jan 10 19:15:49 2007

### 13.4.2 Arguments

**t:** Type `tm`. A Type `tm` data type containing the date and time.

### 13.4.3 Output

Character (len=24). A 24 length character string with the format `Www Mmm dd hh:mm:ss yyyy`, where `Www` is the weekday, `Mmm` the month in letters, `dd` the day of the month, `hh:mm:ss` the time, and `yyyy` the year.

### 13.4.4 Example

A small program that prints the current time.

Listing 13.4: Printing current date/time.

```
1 Program Test
3     USE NumTypes
4     USE Time
5
7     Write(*, '(1A)') asctime(gettime())
9
11    Stop
11 End Program Test
```

## 13.5 Function Day\_of\_Week(Day, Month, Year)

### 13.5.1 Description

The function `Day_of_Week(Day, Month, Year)`, returns the day of the week since sunday (sunday is 0), of the date that correspond to the input `Day, Month, Year`.

### 13.5.2 Arguments

**Day:** Integer. The day of the month [1-31].

**Month:** Integer. The month of the year [0-11].

**Year:** Integer. The year.

### 13.5.3 Output

Integer. The day of the week since sunday, thus a number between 0 and 6, with 0 corresponding to sunday.



### 13.5.4 Example

A small program that prints the date and time of the first of january of 1900.

Listing 13.5: Day of week of the first of January 1900.

```
1 Program Test
3   USE NumTypes
   USE Time
5
   Type (tm) :: Oneday
7
   Oneday%hour = 12
   Oneday%min  = 0
   Oneday%sec  = 0
11  Oneday%nday = 1
   Oneday%mon  = 0
13  Oneday%year = 1900
15
   Oneday%wday = Day_of_Week(Oneday%nday, Oneday%mon, Oneday%year)
17
   Write(*,*) asctime(Oneday)
19
   Stop
End Program Test
```



---

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