



f2py

Python for computational science

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CINECA

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f2py (1)



- f2py allows to build an extension module that interfaces Python to routines Fortran 77/90/95
- Let's write a Fortran function to plot a Mandelbrot fractal
- This fractal is defined by the iteration

$$z \leftarrow z^2 + c$$

where z and c are complex variables. This expression is iterated; if z stays finite, c belongs to the Mandelbrot set

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```
subroutine single_point_mandelbrot(z,c,zout)
!
! ****
!
! *      compute single point mandlebrot      *
!
! ****
implicit none
integer :: I
complex*16 :: z, c, zout
do i = 1,100
    z = z*z + c
    if ((real(z)**2 + aimag(z)**2) .gt. 1000.0) then
        exit
    endif
enddo
zout = z
end subroutine
```

(mandel.f90)

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- Let us get the new module a more pythonic interface

```
$ f2py -h mandel.pyf -m mandel mandel.f90
```

(mandel.pyf)

```
! Note: the context of this file is case sensitive.

python module mandel ! in

    interface ! in :mandel

        subroutine single_point_mandelbrot(z,c,zout) ! in :mandel:mandel.f90
            complex*16 :: z
            complex*16 :: c
            complex*16 :: zout
        end subroutine single_point_mandelbrot

    end interface

end python module mandel

! This file was auto-generated with f2py (version:2).
```

f2py (3)



- Let us get the new module a more pythonic interface

```
$ f2py -h mandel.pyf -m mandel mandel.f90
```

(mandel.pyf)

```
! Note: the context of this file is case sensitive.

python module mandel ! in

    interface ! in :mandel

        subroutine single_point_mandelbrot(z,c,zout) ! in :mandel:mandel.f90
            complex*16 :: z
            complex*16 :: c
            complex*16, intent(out) :: zout
        end subroutine single_point_mandelbrot

    end interface

end python module mandel

! This file was auto-generated with f2py (version:2).
```

f2py (4)



- Let's test the module (and the interface)

```
$ rm mandel.so  
$ f2py -c -m mandel mandel.pyf mandel.f90  
$ python -c 'import mandel'
```

```
>>> import mandel  
>>> print mandel.single_point_mandelbrot.__doc__  
single_point_mandelbrot - Function signature:  
    zout = single_point_mandelbrot(z,c)
```

Required arguments:

```
    z : input complex  
    c : input complex
```

Return objects:

```
    zout : complex
```

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- Let's test the module (and the interface)

```
In [1]: z = complex(1,2)
```

```
In [2]: c = complex(1,2)
```

```
In [3]: zout = mandel.single_point_mandelbrot(z,c)
```

```
In [4]: zout
```

```
Out[4]: (-31-22j)
```

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- Now we can employ our function
- First, we need to wrap our function

```
def myfunc(a,b):  
    return mandel.single_point_mandelbrot(a,b)
```

- And then vectorize it

```
vfunc = np.vectorize(myfunc)
```

- Let's build the input points ...

```
x = np.linspace(-1.7, 0.6, 1000)  
y = np.linspace(-1.4, 1.4, 1000)  
xx, yy = np.meshgrid(x, y)
```

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- ... and then get the complex numbers

```
def f(aa, bb):  
  
    return aa + 1j*bb
```

```
ff = f(xx, yy)      # build a 2d array of complex numbers
```

- Now we can use our new function and plot the results

```
z = vfunc(ff,ff)  # just used our single_point_mandelbrot func  
  
import matplotlib.pyplot as plt  
  
plt.imshow(abs(z)**2 < 1000, extent=[-1.7, 0.6, -1.4, 1.4])  
  
plt.gray()  
  
plt.show()
```

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- Another example: matmul

```
(mult.f90)

subroutine fmult(a,b,c,n)
implicit none
real*8 :: a(*)
real*8 :: b(*)
real*8 :: c(*)
integer :: n, i
do i =1,n
    c(i) = a(i) * b(i)
enddo
end
```

```
$ f2py -c -m mult mult.f90
$ python -c 'import mult'

>>> import mult
>>> print mult.fmult.__doc__
fmult - Function signature:
fmult(a,b,c,n)

Required arguments:
a : input rank-1 array('d') with bounds (100)
b : input rank-1 array('d') with bounds (100)
c : input rank-1 array('d') with bounds (100)
n : input int

>>> a = np.ones(100)+ 3; c = a.copy()
>>> b = np.ones(100)+ 1.5
>>> mult.fmult(a,b,c,100)
```

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- One can improve the interface automatically built

```
$ f2py -h mult.pyf -m mult mult.f90
```

```
(mult.f90)

subroutine fmultiply(a,b,c,n)
    implicit none
    real*8 :: a(*)
    real*8 :: b(*)
    real*8 :: c(*)
    integer :: n, i
    do i =1,n
        c(i) = a(i) * b(i)
    enddo
end
```

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- One can improve the interface automatically built

```
$ f2py -h mult.pyf -m mult mult.f90
```

```
(mult.pyf)
```

```
python module mult

interface

    subroutine fmult(a,b,c,n)
        real*8 dimension(*) :: a
        real*8 dimension(*) :: b
        real*8 dimension(*) :: c
        integer :: n
    end subroutine fmult

    end interface

end python module mult
```

```
(mult.pyf)
```

```
python module mult

interface

    subroutine fmult(a,b,c,n)
        real*8 dimension(n) :: a
        real*8 dimension(n) :: b
        real*8 intent(out),
        dimension(n) :: c
        integer intent(hide), depend(a) :: n=len(a)
    end subroutine fmult

    end interface

end python module mult
```

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- Let us check the new interface

```
$ rm mult.so  
$ f2py -c -m mult mult.pyf mult.f90  
$ python -c 'import mult'
```

```
>>> import mult  
>>> print mult.fmult.__doc__  
fmult - Function signature:  
    c = fmult(a,b)
```

Required arguments:

a : input rank-1 array('d') with bounds (n)
b : input rank-1 array('d') with bounds (n)

Return objects:

c : rank-1 array('d') with bounds (n)

```
>>> import numpy as np  
>>> a = np.array([1, 3, 4])  
>>> b = np.array([2, 5, 1.5])  
>>> c = mult.fmult(a, b)
```

```
>>> f = mult.fmult([3,4], [1.2,1.4])  
>>> f  
array([ 3.6,  5.6])
```

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- Inserting directives in fortran source

```
subroutine fmult(a,b,c,n)
implicit none
integer, parameter :: dim=1000
real*8 :: a(n)
real*8 :: b(n)
real*8 :: c(n)
integer :: n, i
!f2py intent(hide), depend(a) :: n=len(a)
!f2py real*8          :: a(n)
!f2py real*8          :: b(n)
!f2py real*8, intent(out) :: c(n)
do i =1,n
    c(i) = a(i) * b(i)
enddo
end
```

```
$ f2py -c -m mult2 mult2.f90
```

```
>>> import numpy as np
>>> import mult2
>>> a = np.array([1, 3, 4])
>>> b = np.array([2, 5, 1.5])
>>> c = mult2.fmult(a, b)

>>> f = mult2.fmult([3,4], [1.2,1.4])
>>> f
array([ 3.6,  5.6])
```