

# NAG Fortran Library Routine Document

## D03PZF

**Note:** before using this routine, please read the Users' Note for your implementation to check the interpretation of *bold italicised* terms and other implementation-dependent details.

### 1 Purpose

This routine interpolates in the spatial co-ordinate the solution and derivative of a system of partial differential equations (PDEs). The solution must first be computed using one of the finite difference scheme routines D03PCF/D03PCA, D03PHF/D03PHA or D03PPF/D03PPA, or one of the Keller box scheme routines D03PEF, D03PKF or D03PRF.

### 2 Specification

```

SUBROUTINE D03PZF(NPDE, M, U, NPTS, X, XP, INTPTS, ITYPE, UP, IFAIL)
  INTEGER          NPDE, M, NPTS, INTPTS, ITYPE, IFAIL
  real           U(NPDE,NPTS), X(NPTS), XP(INTPTS),
1                UP(NPDE,INTPTS,ITYPE)

```

### 3 Description

D03PZF is an interpolation routine for evaluating the solution of a system of partial differential equations (PDEs), at a set of user-specified points. The solution of the system of equations (possibly with coupled ordinary differential equations) must be computed using a finite difference scheme routine or a Keller box scheme routine on a set of mesh points. D03PZF can then be employed to compute the solution at a set of points anywhere in the range of the mesh. It can also evaluate the first spatial derivative of the solution. The routine uses linear interpolation for approximating the solution.

### 4 References

None.

### 5 Parameters

**Note:** the parameters X, M, U, NPTS and NPDE must be supplied unchanged from the PDE routine.

1: NPDE – INTEGER *Input*

*On entry:* the number of PDEs.

*Constraint:* NPDE  $\geq$  1.

2: M – INTEGER *Input*

*On entry:* the co-ordinate system used. If the call to D03PZF follows one of the finite difference routines then M must be the same parameter M as used by the finite difference routines. For the Keller box scheme routines only Cartesian co-ordinate systems are valid and so M **must** be set to zero. No check will be made by D03PZF in this case.

M = 0

indicates Cartesian co-ordinates

M = 1

indicates cylindrical polar co-ordinates

M = 2

indicates spherical polar co-ordinates

*Constraints:*

$0 \leq M \leq 2$  following a finite difference routine.  
 $M = 0$  following a Keller box scheme routine.

- 3: U(NPDE,NPTS) – *real* array *Input*  
*On entry:* the PDE part of the original solution returned in the parameter U by the PDE routine.  
*Constraint:* NPDE  $\geq$  1.
- 4: NPTS – INTEGER *Input*  
*On entry:* the number of mesh points.  
*Constraint:* NPTS  $\geq$  3.
- 5: X(NPTS) – *real* array *Input*  
*On entry:* X(*i*), for  $i = 1, 2, \dots, \text{NPTS}$ , must contain the mesh points as used by the PDE routine.
- 6: XP(INTPTS) – *real* array *Input*  
*On entry:* XP(*i*), for  $i = 1, 2, \dots, \text{INTPTS}$ , must contain the spatial interpolation points.  
*Constraint:* X(1)  $\leq$  XP(1)  $<$  XP(2)  $<$  ...  $<$  XP(INTPTS)  $\leq$  X(NPTS).
- 7: INTPTS – INTEGER *Input*  
*On entry:* the number of interpolation points.  
*Constraint:* INTPTS  $\geq$  1.
- 8: ITYPE – INTEGER *Input*  
*On entry:* specifies the interpolation to be performed.  
 If ITYPE = 1, the solutions at the interpolation points are computed. If ITYPE = 2, both the solutions and their first derivatives at the interpolation points are computed.  
*Constraint:* ITYPE = 1 or 2.
- 9: UP(NPDE,INTPTS,ITYPE) – *real* array *Output*  
*On exit:* if ITYPE = 1, UP(*i*, *j*, 1), contains the value of the solution  $U_i(x_j, t_{\text{out}})$ , at the interpolation points  $x_j = \text{XP}(j)$ , for  $j = 1, 2, \dots, \text{INTPTS}$ ;  $i = 1, 2, \dots, \text{NPDE}$ .  
 If ITYPE = 2, UP(*i*, *j*, 1) contains  $U_i(x_j, t_{\text{out}})$  and UP(*i*, *j*, 2) contains  $\frac{\partial U_i}{\partial x}$  at these points.
- 10: IFAIL – INTEGER *Input/Output*  
*On entry:* IFAIL must be set to 0, -1 or 1. Users who are unfamiliar with this parameter should refer to Chapter P01 for details.  
*On exit:* IFAIL = 0 unless the routine detects an error (see Section 6).  
 For environments where it might be inappropriate to halt program execution when an error is detected, the value -1 or 1 is recommended. If the output of error messages is undesirable, then the value 1 is recommended. Otherwise, for users not familiar with this parameter the recommended value is 0. **When the value -1 or 1 is used it is essential to test the value of IFAIL on exit.**

## 6 Error Indicators and Warnings

If on entry IFAIL = 0 or -1, explanatory error messages are output on the current error message unit (as defined by X04AAF).

Errors or warnings detected by the routine:

IFAIL = 1

On entry, ITYPE  $\neq$  1 or 2,  
or INTPTS  $<$  1,  
or NPDE  $<$  1,  
or NPTS  $<$  3,  
or  $M \neq$  0, 1 or 2,  
or the mesh points  $X(i)$ , for  $i = 1, 2, \dots, \text{NPTS}$ , are not in strictly increasing order.

IFAIL = 2

On entry, the interpolation points  $XP(i)$ , for  $i = 1, 2, \dots, \text{INTPTS}$ , are not in strictly increasing order.

IFAIL = 3

The user is attempting extrapolation, that is, one of the interpolation points  $XP(i)$ , for some  $i$ , lies outside the interval  $[X(1), X(\text{NPTS})]$ . Extrapolation is not permitted.

## 7 Accuracy

See the PDE routine documents.

## 8 Further Comments

None.

## 9 Example

See Section 9 of the documents for D03PCF/D03PCA, D03PPF/D03PPA and D03PRF.

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