

D03PUF – NAG Fortran Library Routine Document

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

D03PUF calculates a numerical flux function using Roe's Approximate Riemann Solver for the Euler equations in conservative form. The routine is designed primarily for use with the upwind discretisation routines D03PFF, D03PLF or D03PSF, but may also be applicable to other conservative upwind schemes requiring numerical flux functions.

2 Specification

```
SUBROUTINE D03PUF(ULEFT, URIGHT, GAMMA, FLUX, IFAIL)
INTEGER          IFAIL
real             ULEFT(3), URIGHT(3), GAMMA, FLUX(3)
```

3 Description

D03PUF calculates a numerical flux function at a single spatial point using Roe's Approximate Riemann Solver [1] for the Euler equations (for a perfect gas) in conservative form. The user must supply the *left* and *right* solution values at the point where the numerical flux is required, i.e. the initial left and right states of the Riemann problem defined below. In the routines D03PFF, D03PLF and D03PSF, the left and right solution values are derived automatically from the solution values at adjacent spatial points and supplied to the subroutine argument NUMFLX from which the user may call D03PUF. The Euler equations for a perfect gas in conservative form are:

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x} = 0, \quad (1)$$

with

$$U = \begin{bmatrix} \rho \\ m \\ e \end{bmatrix} \quad \text{and} \quad F = \begin{bmatrix} m \\ \frac{m^2}{\rho} + (\gamma - 1) \left(e - \frac{m^2}{2\rho} \right) \\ \frac{me}{\rho} + \frac{m}{\rho} (\gamma - 1) \left(e - \frac{m^2}{2\rho} \right) \end{bmatrix}, \quad (2)$$

where ρ is the density, m is the momentum, e is the specific total energy, and γ is the (constant) ratio of specific heats. The pressure p is given by

$$p = (\gamma - 1) \left(e - \frac{\rho u^2}{2} \right), \quad (3)$$

where $u = m/\rho$ is the velocity.

The routine calculates the Roe approximation to the numerical flux function $F(U_L, U_R) = F(U^*(U_L, U_R))$, where $U = U_L$ and $U = U_R$ are the left and right solution values, and $U^*(U_L, U_R)$ is the intermediate state $\omega(0)$ arising from the similarity solution $U(y, t) = \omega(y/t)$ of the Riemann problem defined by

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial y} = 0, \quad (4)$$

with U and F as in (2), and initial piecewise constant values $U = U_L$ for $y < 0$ and $U = U_R$ for $y > 0$. The spatial domain is $-\infty < y < \infty$, where $y = 0$ is the point at which the numerical flux is required. This implementation of Roe's scheme for the Euler equations uses the so-called parameter-vector method described in [1].

4 References

- [1] Roe P L (1981) Approximate Riemann solvers, parameter vectors, and difference schemes *J. Comput. Phys.* **43** 357–372
- [2] LeVeque R J (1990) *Numerical Methods for Conservation Laws* Birkhäuser Verlag
- [3] Quirk J J (1994) A contribution to the great Riemann solver debate *Internat. J. Numer. Methods Fluids* **18** 555–574

5 Parameters

- 1: ULEFT(3) — *real* array *Input*
On entry: ULEFT(i) must contain the left value of the component U_i for $i = 1, 2, 3$. That is, ULEFT(1) must contain the left value of ρ , ULEFT(2) must contain the left value of m and ULEFT(3) must contain the left value of e .
- 2: URIGHT(3) — *real* array *Input*
On entry: URIGHT(i) must contain the right value of the component U_i for $i = 1, 2, 3$. That is, URIGHT(1) must contain the right value of ρ , URIGHT(2) must contain the right value of m and URIGHT(3) must contain the right value of e .
- 3: GAMMA — *real* *Input*
On entry: the ratio of specific heats γ .
Constraint: GAMMA > 0.0.
- 4: FLUX(3) — *real* array *Output*
On exit: FLUX(i) contains the numerical flux component \hat{F}_i for $i = 1, 2, 3$.
- 5: IFAIL — INTEGER *Input/Output*
On entry: IFAIL must be set to 0, -1 or 1. For users not familiar with this parameter (described in Chapter P01) the recommended value is 0.
On exit: IFAIL = 0 unless the routine detects an error (see Section 6).

Note. If the left and/or right values of ρ or p (from (3)) are found to be negative, then the routine will terminate with an error exit (IFAIL = 2). If the routine is being called from the user-supplied subroutine NUMFLX in D03PFF etc., then a **soft fail** option (IFAIL = 1 or -1) is recommended so that a recalculation of the current time step can be forced using the IRES parameter.

6 Error Indicators and Warnings

Errors detected by the routine:

IFAIL = 1

On entry, GAMMA \leq 0.0.

IFAIL = 2

On entry, the left and/or right density or pressure value is less than 0.0.

7 Accuracy

The routine performs an exact calculation of the Roe numerical flux function, and so the result will be accurate to machine precision.

8 Further Comments

The routine must only be used to calculate the numerical flux for the Euler equations in exactly the form given by (2), with $ULEFT(i)$ and $URIGHT(i)$ containing the left and right values of ρ, m and e for $i = 1, 2, 3$ respectively. It should be noted that Roe's scheme, in common with all Riemann solvers, may be unsuitable for some problems (see Quirk [3] for examples). In particular Roe's scheme does not satisfy an 'entropy condition' which guarantees that the approximate solution of the PDE converges to the correct physical solution, and hence it may admit non-physical solutions such as expansion shocks. The algorithm used in this routine does not detect or correct any entropy violation. The time taken by the routine is independent of the input parameters.

9 Example

See Section 9.2 of the document for D03PLF.
