

NAG Toolbox for MATLAB

f07jh

1 Purpose

f07jh computes error bounds and refines the solution to a real system of linear equations $AX = B$, where A is an n by n symmetric positive-definite tridiagonal matrix and X and B are n by r matrices, using the modified Cholesky factorization returned by f07jd and an initial solution returned by f07je. Iterative refinement is used to reduce the backward error as much as possible.

2 Syntax

```
[x, ferr, berr, info] = f07jh(d, e, df, ef, b, x, 'n', n, 'nrhs_p',  
nrhs_p)
```

3 Description

f07jh should normally be preceded by calls to f07jd and f07je. f07jd computes a modified Cholesky factorization of the matrix A as

$$A = LDL^T,$$

where L is a unit lower bidiagonal matrix and D is a diagonal matrix, with positive diagonal elements. f07je then utilizes the factorization to compute a solution, \hat{X} , to the required equations. Letting \hat{x} denote a column of \hat{X} , f07jh computes a *component-wise backward error*, β , the smallest relative perturbation in each element of A and b such that \hat{x} is the exact solution of a perturbed system

$$(A + E)\hat{x} = b + f, \quad \text{with} \quad |e_{ij}| \leq \beta |a_{ij}|, \quad \text{and} \quad |f_j| \leq \beta |b_j|.$$

The function also estimates a bound for the *component-wise forward error* in the computed solution defined by $\max |x_i - \hat{x}_i| / \max |\hat{x}_i|$, where x is the corresponding column of the exact solution, X .

Note that the modified Cholesky factorization of A can also be expressed as

$$A = U^T D U,$$

where U is unit upper bidiagonal.

4 References

Anderson E, Bai Z, Bischof C, Blackford S, Demmel J, Dongarra J J, Du Croz J J, Greenbaum A, Hammarling S, McKenney A and Sorensen D 1999 *LAPACK Users' Guide* (3rd Edition) SIAM, Philadelphia URL: <http://www.netlib.org/lapack/lug>

5 Parameters

5.1 Compulsory Input Parameters

1: **d(*)** – double array

Note: the dimension of the array **d** must be at least $\max(1, \mathbf{n})$.

Must contain the n diagonal elements of the matrix of A .

2: **e(*)** – double array

Note: the dimension of the array **e** must be at least $\max(1, \mathbf{n} - 1)$.

Must contain the $(n - 1)$ subdiagonal elements of the matrix A .

3: **df(*) – double array**

Note: the dimension of the array **df** must be at least $\max(1, \mathbf{n})$.

Must contain the n diagonal elements of the diagonal matrix D from the LDL^T factorization of A .

4: **ef(*) – double array**

Note: the dimension of the array **ef** must be at least $\max(1, \mathbf{n})$.

Must contain the $(n - 1)$ subdiagonal elements of the unit bidiagonal matrix L from the LDL^T factorization of A .

5: **b(lb,*) – double array**

The first dimension of the array **b** must be at least $\max(1, \mathbf{n})$

The second dimension of the array must be at least $\max(1, \mathbf{nrhs_p})$

The n by r matrix of right-hand sides B .

6: **x(ldx,*) – double array**

The first dimension of the array **x** must be at least $\max(1, \mathbf{n})$

The second dimension of the array must be at least $\max(1, \mathbf{nrhs_p})$

The n by r initial solution matrix X .

5.2 Optional Input Parameters

1: **n – int32 scalar**

Default: The dimension of the array **d** The dimension of the array **df** The dimension of the array **ef**.
 n , the order of the matrix A .

Constraint: $\mathbf{n} \geq 0$.

2: **nrhs_p – int32 scalar**

Default: The second dimension of the array **b** The second dimension of the array **x**.
 r , the number of right-hand sides, i.e., the number of columns of the matrix B .

Constraint: $\mathbf{nrhs_p} \geq 0$.

5.3 Input Parameters Omitted from the MATLAB Interface

ldb, ldx, work

5.4 Output Parameters

1: **x(ldx,*) – double array**

The first dimension of the array **x** must be at least $\max(1, \mathbf{n})$

The second dimension of the array must be at least $\max(1, \mathbf{nrhs_p})$

The n by r refined solution matrix X .

2: **ferr(*) – double array**

Note: the dimension of the array **ferr** must be at least $\max(1, \mathbf{nrhs_p})$.

Estimate of the forward error bound for each computed solution vector, such that $\|\hat{x}_j - x_j\|_\infty / \|x_j\|_\infty \leq \mathbf{ferr}(j)$, where \hat{x}_j is the j th column of the computed solution returned in the

array \mathbf{x} and x_j is the corresponding column of the exact solution X . The estimate is almost always a slight overestimate of the true error.

3: **berr**(*) – double array

Note: the dimension of the array **berr** must be at least $\max(1, \text{nrhs_p})$.

Estimate of the component-wise relative backward error of each computed solution vector \hat{x}_j (i.e., the smallest relative change in any element of A or B that makes \hat{x}_j an exact solution).

4: **info** – int32 scalar

info = 0 unless the function detects an error (see Section 6).

6 Error Indicators and Warnings

Errors or warnings detected by the function:

info = $-i$

If **info** = $-i$, parameter i had an illegal value on entry. The parameters are numbered as follows:

1: **n**, 2: **nrhs_p**, 3: **d**, 4: **e**, 5: **df**, 6: **ef**, 7: **b**, 8: **ldb**, 9: **x**, 10: **ldx**, 11: **ferr**, 12: **berr**, 13: **work**, 14: **info**.

It is possible that **info** refers to a parameter that is omitted from the MATLAB interface. This usually indicates that an error in one of the other input parameters has caused an incorrect value to be inferred.

7 Accuracy

The computed solution for a single right-hand side, \hat{x} , satisfies an equation of the form

$$(A + E)\hat{x} = b,$$

where

$$\|E\|_{\infty} = O(\epsilon)\|A\|_{\infty}$$

and ϵ is the *machine precision*. An approximate error bound for the computed solution is given by

$$\frac{\|\hat{x} - x\|_{\infty}}{\|x\|_{\infty}} \leq \kappa(A) \frac{\|E\|_{\infty}}{\|A\|_{\infty}},$$

where $\kappa(A) = \|A^{-1}\|_{\infty}\|A\|_{\infty}$, the condition number of A with respect to the solution of the linear equations. See Section 4.4 of Anderson *et al.* 1999 for further details. Function f07jg can be used to compute the condition number of A .

8 Further Comments

The total number of floating-point operations required to solve the equations $AX = B$ is proportional to nr . At most five steps of iterative refinement are performed, but usually only one or two steps are required.

The complex analogue of this function is f07jv.

9 Example

```
d = [4;
     10;
     29;
     25;
     5];
```

```
e = [-2;  
      -6;  
      15;  
      8];  
df = [4;  
      9;  
      25;  
      16;  
      1];  
ef = [-0.5;  
      -0.6666666666666666;  
      0.6;  
      0.5;  
      0];  
b = [6, 10;  
      9, 4;  
      2, 9;  
      14, 65;  
      7, 23];  
x = [2.5, 2;  
      2, -0.9999999999999999;  
      1, -3;  
      -1, 6;  
      3, -5];  
[xOut, ferr, berr, info] = f07jh(d, e, df, ef, b, x)
```

```
xOut =  
    2.5000    2.0000  
    2.0000   -1.0000  
    1.0000   -3.0000  
   -1.0000    6.0000  
    3.0000   -5.0000  
ferr =  
    1.0e-13 *  
    0.2425  
    0.4663  
berr =  
    1.0e-16 *  
         0  
    0.1110  
info =  
         0
```