NAG Toolbox for MATLAB

f07fp

1 Purpose

f07fp uses the Cholesky factorization

$$A = U^{\mathrm{H}}U \qquad \text{or} \qquad A = LL^{\mathrm{H}}$$

to compute the solution to a complex system of linear equations

$$AX = B,$$

where A is an n by n Hermitian positive-definite matrix and X and B are n by r matrices. Error bounds on the solution and a condition estimate are also provided.

2 Syntax

```
[a, af, equed, s, b, x, rcond, ferr, berr, info] = f07fp(fact, uplo, a, af, equed, s, b, 'n', n, 'nrhs_p', nrhs_p)
```

3 Description

f07fp performs the following steps:

1. If **fact** = 'E', real diagonal scaling factors, D_S, are computed to equilibrate the system:

   $$(D_S A D_S)\left(D_S^{-1}X\right) = D_S B.$$

   Whether or not the system will be equilibrated depends on the scaling of the matrix A, but if equilibration is used, A is overwritten by $D_S A D_S$ and B by $D_S B$.
2. If **fact** = 'N' or 'E', the Cholesky decomposition is used to factor the matrix A (after equilibration if **fact** = 'E') as $A = U^{\mathrm{H}}U$ if **uplo** = 'U' or $A = LL^{\mathrm{H}}$ if **uplo** = 'L', where U is an upper triangular matrix and L is a lower triangular matrix.
3. If the leading i by i principal minor is not positive-definite, then the function returns with **info** = i. Otherwise, the factored form of A is used to estimate the condition number of the matrix A. If the reciprocal of the condition number is less than ***machine precision***, **info** $\geq N + 1$ is returned as a warning, but the function still goes on to solve for X and compute error bounds as described below.
4. The system of equations is solved for X using the factored form of A.
5. Iterative refinement is applied to improve the computed solution matrix and to calculate error bounds and backward error estimates for it.
6. If equilibration was used, the matrix X is premultiplied by D_S so that it solves the original system before equilibration.

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

Golub G H and Van Loan C F 1996 *Matrix Computations* (3rd Edition) Johns Hopkins University Press, Baltimore

Higham N J 2002 *Accuracy and Stability of Numerical Algorithms* (2nd Edition) SIAM, Philadelphia

5 Parameters

5.1 Compulsory Input Parameters

1: **fact – string**

Specifies whether or not the factorized form of the matrix A is supplied on entry, and if not, whether the matrix A should be equilibrated before it is factorized.

fact = 'F'

af contains the factorized form of A. If **equed** = 'Y', the matrix A has been equilibrated with scaling factors given by **s**. **af** will not be modified.

fact = 'N'

The matrix A will be copied to **af** and factorized.

fact = 'E'

The matrix A will be equilibrated if necessary, then copied to **af** and factorized.

Constraint: **fact** = 'F', 'N' or 'E'.

2: **uplo – string**

If **uplo** = 'U', the upper triangle of A is stored.

If **uplo** = 'L', the lower triangle of A is stored.

Constraint: **uplo** = 'U' or 'L'.

3: **a**(**lda**,$*$) **– complex array**

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

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

The n by n Hermitian matrix A.

If **fact** = 'F' and **equed** = 'Y', **a** must have been equilibrated by the scaling factor in **s** as $D_S A D_S$.

If **uplo** = 'U', the upper triangular part of A must be stored and the elements of the array below the diagonal are not referenced.

If **uplo** = 'L', the lower triangular part of A must be stored and the elements of the array above the diagonal are not referenced.

4: **af**(**ldaf**,$*$) **– complex array**

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

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

If **fact** = 'F', **af** contains the triangular factor U or L from the Cholesky factorization $A = U^{\mathrm{H}}U$ or $A = LL^{\mathrm{H}}$, in the same storage format as **a**. If **equed** $\neq$ 'N', **af** is the factorized form of the equilibrated matrix $D_S A D_S$.

5: **equed – string**

If **fact** = 'N' or 'E', **equed** need not be set.

If **fact** = 'F', **equed** must specify the form of the equilibration that was performed as follows:

if **equed** = 'N', no equilibration;
if **equed** = 'Y', equilibration was performed, i.e., A has been replaced by $D_S A D_S$.

Constraint: if **fact** = 'F', **equed** = 'N' or 'Y'.

6: **s**$(*)$ **– double array**

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

If **fact** = 'N' or 'E', **s** need not be set.

If **fact** = 'F' and **equed** = 'Y', **s** must contain the scale factors, D_S, for A; each element of **s** must be positive.

7: **b**(**ldb,***) **– complex 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 right-hand side matrix B.

5.2 Optional Input Parameters

1: **n – int32 scalar**

Default: The second dimension of the array **a** The second dimension of the array **af** The dimension of the array **s**.

n, the number of linear equations, i.e., the order of the matrix A.

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

2: **nrhs_p – int32 scalar**

Default: The second dimension of the array **b**.

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

lda, ldaf, ldb, ldx, work, rwork

5.4 Output Parameters

1: **a**(**lda,***) **– complex array**

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

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

If **fact** = 'F' or 'N', or if **fact** = 'E' and **equed** = 'N', **a** is not modified.

If **fact** = 'E' and **equed** = 'Y', **a** contains $D_S A D_S$.

2: **af**(**ldaf,***) **– complex array**

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

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

If **fact** = 'N', **af** returns the triangular factor U or L from the Cholesky factorization $A = U^{\mathrm{H}}U$ or $A = LL^{\mathrm{H}}$ of the original matrix A.

If **fact** = 'E', **af** returns the triangular factor U or L from the Cholesky factorization $A = U^{\mathrm{H}}U$ or $A = LL^{\mathrm{H}}$ of the equilibrated matrix A (see the description of **a** for the form of the equilibrated matrix).

3: **equed – string**

If **fact** = 'F', **equed** is unchanged from entry.

Otherwise, if **info** ≥ 0, **equed** specifies the form of the equilibration that was performed as specified above.

4: **s**$(*)$ **– double array**

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

If **fact** = 'F', **s** is unchanged from entry.

Otherwise, if **info** ≥ 0 and **equed** = 'Y', **s** contains the scale factors, D_S, for A; each element of **s** is positive.

5: **b**(**ldb**,$*$) **– complex 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 $\max(1, \mathbf{nrhs\_p})$

If **equed** = 'N', **b** is not modified.

If **equed** = 'Y', **b** contains $D_S B$.

6: **x**(**ldx**,$*$) **– complex 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 $\max(1, \mathbf{nrhs\_p})$

If **info** $= 0$ or **info** $\geq N + 1$, the n by r solution matrix X to the original system of equations. Note that the arrays A and B are modified on exit if **equed** = 'Y', and the solution to the equilibrated system is $D_S^{-1} X$.

7: **rcond – double scalar**

If **info** ≥ 0, an estimate of the reciprocal condition number of the matrix A (after equilibration if that is performed), computed as **rcond** $= 1/\left(\|A\|_1 \left\|A^{-1}\right\|_1\right)$.

8: **ferr**$(*)$ **– double array**

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

If **info** $= 0$ or **info** $\geq N + 1$, an estimate of the forward error bound for each computed solution vector, such that $\left\|\hat{x}_j - x_j\right\|_\infty / \left\|x_j\right\|_\infty \leq$ **ferr**(j) where $\hat{x}_j$ is the jth column of the computed solution returned in the array **x** and x_j is the corresponding column of the exact solution X. The estimate is as reliable as the estimate for **rcond**, and is almost always a slight overestimate of the true error.

9: **berr**$(*)$ **– double array**

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

If **info** $= 0$ or **info** $\geq N + 1$, an 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).

10: **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: **fact**, 2: **uplo**, 3: **n**, 4: **nrhs_p**, 5: **a**, 6: **lda**, 7: **af**, 8: **ldaf**, 9: **equed**, 10: **s**, 11: **b**, 12: **ldb**, 13: **x**, 14: **ldx**, 15: **rcond**, 16: **ferr**, 17: **berr**, 18: **work**, 19: **rwork**, 20: **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.

info > 0 and **info** $\leq N$

If **info** $= i$ and $i \leq$ **n**, the leading minor of order i of A is not positive-definite, so the factorization could not be completed, and the solution has not been computed. **rcond** $= 0$ is returned.

info $= N + 1$

U is nonsingular, but **rcond** is less than ***machine precision***, meaning that the matrix is singular to working precision. Nevertheless, the solution and error bounds are computed because there are a number of situations where the computed solution can be more accurate than the value of **rcond** would suggest.

7 Accuracy

For each right-hand side vector b, the computed solution x is the exact solution of a perturbed system of equations $(A + E)x = b$, where

$$|E| \leq c(n)\epsilon \left|U^{\mathrm{T}}\right||U|,$$

$c(n)$ is a modest linear function of n, and ϵ is the ***machine precision***. See Section 10.1 of Higham 2002 for further details.

If $\hat{x}$ is the true solution, then the computed solution x satisfies a forward error bound of the form

$$\frac{\|x - \hat{x}\|_\infty}{\|\hat{x}\|_\infty} \leq w_c \operatorname{cond}(A, \hat{x}, b)$$

where $\operatorname{cond}(A, \hat{x}, b) = \left\| \left|A^{-1}\right|(|A||\hat{x}| + |b|)\right\|_\infty / \|\hat{x}\|_\infty \leq \operatorname{cond}(A) = \left\| \left|A^{-1}\right||A| \right\|_\infty \leq \kappa_\infty(A)$. If $\hat{x}$ is the jth column of X, then w_c is returned in **berr**(j) and a bound on $\|x - \hat{x}\|_\infty / \|\hat{x}\|_\infty$ is returned in **ferr**(j). See Section 4.4 of Anderson *et al.* 1999 for further details.

8 Further Comments

The factorization of A requires approximately $\frac{4}{3}n^3$ floating-point operations.

For each right-hand side, computation of the backward error involves a minimum of $16n^2$ floating-point operations. Each step of iterative refinement involves an additional $24n^2$ operations. At most five steps of iterative refinement are performed, but usually only one or two steps are required. Estimating the forward error involves solving a number of systems of equations of the form $Ax = b$; the number is usually 4 or 5 and never more than 11. Each solution involves approximately $8n^2$ operations.

The real analogue of this function is f07fb.

9 Example

```
fact = 'Equilibration';
uplo = 'Upper';
a  =  [complex(3.23,  +0),  complex(1.51,  -1.92),  complex(1.9,  +0.84),
complex(0.42, +2.5);
      complex(0, 0), complex(3.58, +0), complex(-0.23, +1.11), complex(-
1.18, +1.37);
       complex(0, 0), complex(0, 0), complex(4.09, +0), complex(2.33, -
0.14);
     complex(0, 0), complex(0, 0), complex(0, +0), complex(4.29, +0)];
```

```
af = complex(zeros(4, 4));
equed = ' ';
s = zeros(4, 1);
b = [complex(3.93, -6.14), complex(1.48, +6.58);
     complex(6.17, +9.42), complex(4.65, -4.75);
     complex(-7.17, -21.83), complex(-4.91, +2.29);
     complex(1.99, -14.38), complex(7.64, -10.79)];
[aOut, afOut, equedOut, sOut, bOut, x, rcond, ferr, berr, info] =
f07fp(fact, uplo, a, af, equed, s, b)
```

```
aOut =
    3.2300                1.5100 - 1.9200i    1.9000 + 0.8400i    0.4200 +
2.5000i
         0                3.5800              -0.2300 + 1.1100i   -1.1800 +
1.3700i
         0                     0               4.0900                2.3300 -
0.1400i
         0                     0                    0             4.2900
afOut =
    1.7972                0.8402 - 1.0683i    1.0572 + 0.4674i    0.2337 +
1.3910i
         0                1.3164              -0.4702 - 0.3131i    0.0834 -
0.0368i
         0                     0               1.5604                0.9360 -
0.9900i
         0                     0                    0             0.6603
equedOut =
N
sOut =
    0.5564
    0.5285
    0.4945
    0.4828
bOut =
   3.9300 - 6.1400i   1.4800 + 6.5800i
   6.1700 + 9.4200i   4.6500 - 4.7500i
  -7.1700 -21.8300i  -4.9100 + 2.2900i
   1.9900 -14.3800i   7.6400 -10.7900i
x =
   1.0000 - 1.0000i  -1.0000 + 2.0000i
  -0.0000 + 3.0000i   3.0000 - 4.0000i
  -4.0000 - 5.0000i  -2.0000 + 3.0000i
   2.0000 + 1.0000i   4.0000 - 5.0000i
rcond =
    0.0066
ferr =
   1.0e-13 *
    0.5965
    0.7494
berr =
   1.0e-16 *
    0.8204
    0.9252
info =
           0
```