

## NAG Toolbox

### nag\_lapack\_zpttrs (f07js)

#### 1 Purpose

nag\_lapack\_zpttrs (f07js) computes the solution to a complex system of linear equations  $AX = B$ , where  $A$  is an  $n$  by  $n$  Hermitian positive definite tridiagonal matrix and  $X$  and  $B$  are  $n$  by  $r$  matrices, using the  $LDL^H$  factorization returned by nag\_lapack\_zpttrf (f07jr).

#### 2 Syntax

```
[b, info] = nag_lapack_zpttrs(uplo, d, e, b, 'n', n, 'nrhs_p', nrhs_p)
```

```
[b, info] = f07js(uplo, d, e, b, 'n', n, 'nrhs_p', nrhs_p)
```

#### 3 Description

nag\_lapack\_zpttrs (f07js) should be preceded by a call to nag\_lapack\_zpttrf (f07jr), which computes a modified Cholesky factorization of the matrix  $A$  as

$$A = LDL^H,$$

where  $L$  is a unit lower bidiagonal matrix and  $D$  is a diagonal matrix, with positive diagonal elements. nag\_lapack\_zpttrs (f07js) then utilizes the factorization to solve the required equations. Note that the factorization may also be regarded as having the form  $U^H DU$ , where  $U$  is a unit upper bidiagonal matrix.

#### 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 <http://www.netlib.org/lapack/lug>

#### 5 Parameters

##### 5.1 Compulsory Input Parameters

1: **uplo** – CHARACTER(1)

Specifies the form of the factorization as follows:

**uplo** = 'U'

$$A = U^H DU.$$

**uplo** = 'L'

$$A = LDL^H.$$

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

2: **d**(:) – REAL (KIND=nag\_wp) array

The dimension of the array **d** must be at least  $\max(1, n)$

Must contain the  $n$  diagonal elements of the diagonal matrix  $D$  from the  $LDL^H$  or  $U^H DU$  factorization of  $A$ .

3: **e**(:) – COMPLEX (KIND=nag\_wp) array

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

If **uplo** = 'U', **e** must contain the  $(n - 1)$  superdiagonal elements of the unit upper bidiagonal matrix  $U$  from the  $U^H D U$  factorization of  $A$ .

If **uplo** = 'L', **e** must contain the  $(n - 1)$  subdiagonal elements of the unit lower bidiagonal matrix  $L$  from the  $LDL^H$  factorization of  $A$ .

4: **b**(ldb,:) – COMPLEX (KIND=nag\_wp) array

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

The second dimension of the array **b** must be at least  $\max(1, \mathbf{nrhs\_p})$ .

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

## 5.2 Optional Input Parameters

1: **n** – INTEGER

*Default:* the first dimension of the array **b** and the dimension of the array **d**.  
 $n$ , the order of the matrix  $A$ .

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

2: **nrhs\_p** – INTEGER

*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 Output Parameters

1: **b**(ldb,:) – COMPLEX (KIND=nag\_wp) array

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

The second dimension of the array **b** will be  $\max(1, \mathbf{nrhs\_p})$ .

The  $n$  by  $r$  solution matrix  $X$ .

2: **info** – INTEGER

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

## 6 Error Indicators and Warnings

**info** < 0

If **info** =  $-i$ , argument  $i$  had an illegal value. An explanatory message is output, and execution of the program is terminated.

## 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

$$\|E\|_1 = O(\epsilon)\|A\|_1$$

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

$$\frac{\|\hat{x} - x\|_1}{\|x\|_1} \leq \kappa(A) \frac{\|E\|_1}{\|A\|_1},$$

where  $\kappa(A) = \|A^{-1}\|_1 \|A\|_1$ , 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.

Following the use of this function `nag_lapack_zptcon` (f07ju) can be used to estimate the condition number of  $A$  and `nag_lapack_zptrfs` (f07jv) can be used to obtain approximate error bounds.

## 8 Further Comments

The total number of floating-point operations required to solve the equations  $AX = B$  is proportional to  $nr$ .

The real analogue of this function is `nag_lapack_dptrs` (f07je).

## 9 Example

This example solves the equations

$$AX = B,$$

where  $A$  is the Hermitian positive definite tridiagonal matrix

$$A = \begin{pmatrix} 16.0 & 16.0 - 16.0i & 0 & 0 \\ 16.0 + 16.0i & 41.0 & 18.0 + 9.0i & 0.0 \\ 0 & 18.0 - 9.0i & 46.0 & 1.0 + 4.0i \\ 0 & 0 & 1.0 - 4.0i & 21.0 \end{pmatrix}$$

and

$$B = \begin{pmatrix} 64.0 + 16.0i & -16.0 - 32.0i \\ 93.0 + 62.0i & 61.0 - 66.0i \\ 78.0 - 80.0i & 71.0 - 74.0i \\ 14.0 - 27.0i & 35.0 + 15.0i \end{pmatrix}.$$

### 9.1 Program Text

```
function f07js_example
fprintf('f07js example results\n\n');

% Hermitian tridiagonal A stored as two diagonals
d = [ 16          41          46          21];
e = [ 16 + 16i    18 - 9i     1 - 4i         ];

% Factorize
[df, ef, info] = f07jr( ...
    d, e);

% RHS
b = [ 64 + 16i,   -16 - 32i;
      93 + 62i,    61 - 66i;
      78 - 80i,    71 - 74i;
      14 - 27i,    35 + 15i];

% Solve
uplo = 'L';
[x, info] = f07js( ...
    uplo, df, ef, b);

disp('Solution(s)');
disp(x);
```

## 9.2 Program Results

f07js example results

Solution(s)

2.0000 + 1.0000i	-3.0000 - 2.0000i
1.0000 + 1.0000i	1.0000 + 1.0000i
1.0000 - 2.0000i	1.0000 - 2.0000i
1.0000 - 1.0000i	2.0000 + 1.0000i

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