

## NAG Toolbox

### nag\_lapack\_zstein (f08jx)

#### 1 Purpose

nag\_lapack\_zstein (f08jx) computes the eigenvectors of a real symmetric tridiagonal matrix corresponding to specified eigenvalues, by inverse iteration, storing the eigenvectors in a complex array.

#### 2 Syntax

```
[z, ifailv, info] = nag_lapack_zstein(d, e, m, w, iblock, isplit, 'n', n)
[z, ifailv, info] = f08jx(d, e, m, w, iblock, isplit, 'n', n)
```

#### 3 Description

nag\_lapack\_zstein (f08jx) computes the eigenvectors of a real symmetric tridiagonal matrix  $T$  corresponding to specified eigenvalues, by inverse iteration (see Jessup and Ipsen (1992)). It is designed to be used in particular after the specified eigenvalues have been computed by nag\_lapack\_dstebz (f08jj) with **order** = 'B', but may also be used when the eigenvalues have been computed by other functions in Chapters F02 or F08.

The eigenvectors of  $T$  are real, but are stored by this function in a complex array. If  $T$  has been formed by reduction of a full complex Hermitian matrix  $A$  to tridiagonal form, then eigenvectors of  $T$  may be transformed to (complex) eigenvectors of  $A$  by a call to nag\_lapack\_zunmtr (f08fu) or nag\_lapack\_zupmtr (f08gu).

nag\_lapack\_dstebz (f08jj) determines whether the matrix  $T$  splits into block diagonal form:

$$T = \begin{pmatrix} T_1 & & & & \\ & T_2 & & & \\ & & \cdot & & \\ & & & \cdot & \\ & & & & T_p \end{pmatrix}$$

and passes details of the block structure to this function in the arrays **iblock** and **isplit**. This function can then take advantage of the block structure by performing inverse iteration on each block  $T_i$  separately, which is more efficient than using the whole matrix.

#### 4 References

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

Jessup E and Ipsen I C F (1992) Improving the accuracy of inverse iteration *SIAM J. Sci. Statist. Comput.* **13** 550–572

#### 5 Parameters

##### 5.1 Compulsory Input Parameters

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

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

The diagonal elements of the tridiagonal matrix  $T$ .

- 2: **e**(:) – REAL (KIND=nag\_wp) array  
 The dimension of the array **e** must be at least  $\max(1, \mathbf{n} - 1)$   
 The off-diagonal elements of the tridiagonal matrix  $T$ .
- 3: **m** – INTEGER  
 $m$ , the number of eigenvectors to be returned.  
*Constraint:*  $0 \leq \mathbf{m} \leq \mathbf{n}$ .
- 4: **w**(:) – REAL (KIND=nag\_wp) array  
 The dimension of the array **w** must be at least  $\max(1, \mathbf{n})$   
 The eigenvalues of the tridiagonal matrix  $T$  stored in **w**(1) to **w**( $m$ ), as returned by nag\_lapack\_dstebz (f08jj) with **order** = 'B'. Eigenvalues associated with the first sub-matrix must be supplied first, in nondecreasing order; then those associated with the second sub-matrix, again in nondecreasing order; and so on.  
*Constraint:* if **iblock**( $i$ ) = **iblock**( $i + 1$ ), **w**( $i$ )  $\leq$  **w**( $i + 1$ ), for  $i = 1, 2, \dots, \mathbf{m} - 1$ .
- 5: **iblock**(:) – INTEGER array  
 The dimension of the array **iblock** must be at least  $\max(1, \mathbf{n})$   
 The first  $m$  elements must contain the sub-matrix indices associated with the specified eigenvalues, as returned by nag\_lapack\_dstebz (f08jj) with **order** = 'B'. If the eigenvalues were not computed by nag\_lapack\_dstebz (f08jj) with **order** = 'B', set **iblock**( $i$ ) to 1, for  $i = 1, 2, \dots, m$ .  
*Constraint:* **iblock**( $i$ )  $\leq$  **iblock**( $i + 1$ ), for  $i = 1, 2, \dots, \mathbf{m} - 1$ .
- 6: **isplit**(:) – INTEGER array  
 The dimension of the array **isplit** must be at least  $\max(1, \mathbf{n})$   
 The points at which  $T$  breaks up into sub-matrices, as returned by nag\_lapack\_dstebz (f08jj) with **order** = 'B'. If the eigenvalues were not computed by nag\_lapack\_dstebz (f08jj) with **order** = 'B', set **isplit**(1) to **n**.

## 5.2 Optional Input Parameters

- 1: **n** – INTEGER  
*Default:* the first dimension of the array **d** and the second dimension of the array **d**. (An error is raised if these dimensions are not equal.)  
 $n$ , the order of the matrix  $T$ .  
*Constraint:*  $\mathbf{n} \geq 0$ .

## 5.3 Output Parameters

- 1: **z**(ldz,:) – COMPLEX (KIND=nag\_wp) array  
 The first dimension of the array **z** will be  $\max(1, \mathbf{n})$ .  
 The second dimension of the array **z** will be  $\max(1, \mathbf{m})$ .  
 The  $m$  eigenvectors, stored as columns of  $Z$ ; the  $i$ th column corresponds to the  $i$ th specified eigenvalue, unless **info**  $>$  0 (in which case see Section 6).
- 2: **ifailv**(**m**) – INTEGER array  
 If **info** =  $i >$  0, the first  $i$  elements of **ifailv** contain the indices of any eigenvectors which have failed to converge. The rest of the first **m** elements of **ifailv** are set to 0.

- 3: **info** – INTEGER  
**info** = 0 unless the function detects an error (see Section 6).

## 6 Error Indicators and Warnings

**info** =  $-i$

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

1: **n**, 2: **d**, 3: **e**, 4: **m**, 5: **w**, 6: **iblock**, 7: **isplit**, 8: **z**, 9: **ldz**, 10: **work**, 11: **iwork**, 12: **ifailv**, 13: **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 (*warning*)

If **info** =  $i$ , then  $i$  eigenvectors (as indicated by the argument **ifailv** above) each failed to converge in five iterations. The current iterate after five iterations is stored in the corresponding column of **z**.

## 7 Accuracy

Each computed eigenvector  $z_i$  is the exact eigenvector of a nearby matrix  $A + E_i$ , such that

$$\|E_i\| = O(\epsilon)\|A\|,$$

where  $\epsilon$  is the *machine precision*. Hence the residual is small:

$$\|Az_i - \lambda_i z_i\| = O(\epsilon)\|A\|.$$

However, a set of eigenvectors computed by this function may not be orthogonal to so high a degree of accuracy as those computed by nag\_lapack\_zsteqr (f08js).

## 8 Further Comments

The real analogue of this function is nag\_lapack\_dstein (f08jk).

## 9 Example

See Section 10 in nag\_lapack\_zunmtr (f08fu).

### 9.1 Program Text

```
function f08jx_example

fprintf('f08jx example results\n\n');

% Lower triangular part of Hermitian matrix A
uplo = 'L';
a = [-2.28 + 0.00i, 0.00 + 0i, 0 + 0i, 0 + 0i;
     1.78 + 2.03i, -1.12 + 0i, 0 + 0i, 0 + 0i;
     2.26 - 0.10i, 0.01 - 0.43i, -0.37 + 0i, 0 + 0i;
     -0.12 - 2.53i, -1.07 - 0.86i, 2.31 + 0.92i, -0.73 + 0i];

% Reduce A to tridiagonal form
[aq, d, e, tau, info] = f08fs( ...
    uplo, a);

% Calculate two smallest eigenvalues
range = 'Indices';
order = 'Block';
vl = 0;
```

```

vu = 0;
il = nag_int(1);
iu = nag_int(2);
abstol = 0;
[m, nsplit, w, iblock, isplit, info] = ...
    f08jj( ...
        range, order, vl, vu, il, iu, abstol, d, e);

% Corresponding eigenvectors of T
[tz, ifailv, info] = f08jx( ...
    d, e, m, w, iblock, isplit);

% TTransform to eigenvalues of A (by premultiplying by Q)
trans = 'No transpose';
side = 'Left';
[z, info] = f08fu( ...
    side, uplo, trans, aq, tau, tz);

% Normalize vectors, largest element is real and positive.
for i = 1:m
    [~,k] = max(abs(real(z(:,i)))+abs(imag(z(:,i))));
    z(:,i) = z(:,i)*conj(z(k,i))/abs(z(k,i));
end

disp(' Selected eigenvalues of A:');
disp(w(1:m));
disp(' Corresponding eigenvectors:');
disp(z);

```

## 9.2 Program Results

f08jx example results

Selected eigenvalues of A:

```

-6.0002
-3.0030

```

Corresponding eigenvectors:

```

0.7299 + 0.0000i  -0.2120 + 0.1497i
-0.1663 - 0.2061i  0.7307 + 0.0000i
-0.4165 - 0.1417i  -0.3291 + 0.0479i
0.1743 + 0.4162i  0.5200 + 0.1329i

```

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