

## NAG Toolbox

### nag\_lapack\_dsbevz (f08hb)

#### 1 Purpose

nag\_lapack\_dsbevz (f08hb) computes selected eigenvalues and, optionally, eigenvectors of a real  $n$  by  $n$  symmetric band matrix  $A$  of bandwidth  $(2k_d + 1)$ . Eigenvalues and eigenvectors can be selected by specifying either a range of values or a range of indices for the desired eigenvalues.

#### 2 Syntax

```
[ab, q, m, w, z, jfail, info] = nag_lapack_dsbevz(jobz, range, uplo, kd, ab, vl,
vu, il, iu, abstol, 'n', n)
```

```
[ab, q, m, w, z, jfail, info] = f08hb(jobz, range, uplo, kd, ab, vl, vu, il, iu,
abstol, 'n', n)
```

#### 3 Description

The symmetric band matrix  $A$  is first reduced to tridiagonal form, using orthogonal similarity transformations. The required eigenvalues and eigenvectors are then computed from the tridiagonal matrix; the method used depends upon whether all, or selected, eigenvalues and eigenvectors are required.

#### 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>

Demmel J W and Kahan W (1990) Accurate singular values of bidiagonal matrices *SIAM J. Sci. Statist. Comput.* **11** 873–912

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

#### 5 Parameters

##### 5.1 Compulsory Input Parameters

1: **jobz** – CHARACTER(1)

Indicates whether eigenvectors are computed.

**jobz** = 'N'

Only eigenvalues are computed.

**jobz** = 'V'

Eigenvalues and eigenvectors are computed.

*Constraint:* **jobz** = 'N' or 'V'.

2: **range** – CHARACTER(1)

If **range** = 'A', all eigenvalues will be found.

If **range** = 'V', all eigenvalues in the half-open interval  $(\mathbf{vl}, \mathbf{vu}]$  will be found.

If **range** = 'I', the **ilth** to **iuth** eigenvalues will be found.

*Constraint:* **range** = 'A', 'V' or 'I'.

3: **uplo** – CHARACTER(1)

If **uplo** = 'U', the upper triangular part of  $A$  is stored.

If **uplo** = 'L', the lower triangular part of  $A$  is stored.

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

4: **kd** – INTEGER

If **uplo** = 'U', the number of superdiagonals,  $k_d$ , of the matrix  $A$ .

If **uplo** = 'L', the number of subdiagonals,  $k_d$ , of the matrix  $A$ .

*Constraint:* **kd**  $\geq$  0.

5: **ab**(*ldab*,:) – REAL (KIND=nag\_wp) array

The first dimension of the array **ab** must be at least **kd** + 1.

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

The upper or lower triangle of the  $n$  by  $n$  symmetric band matrix  $A$ .

The matrix is stored in rows 1 to  $k_d + 1$ , more precisely,

if **uplo** = 'U', the elements of the upper triangle of  $A$  within the band must be stored with element  $A_{ij}$  in **ab**( $k_d + 1 + i - j, j$ ) for  $\max(1, j - k_d) \leq i \leq j$ ;

if **uplo** = 'L', the elements of the lower triangle of  $A$  within the band must be stored with element  $A_{ij}$  in **ab**( $1 + i - j, j$ ) for  $j \leq i \leq \min(n, j + k_d)$ .

6: **vl** – REAL (KIND=nag\_wp)

7: **vu** – REAL (KIND=nag\_wp)

If **range** = 'V', the lower and upper bounds of the interval to be searched for eigenvalues.

If **range** = 'A' or 'I', **vl** and **vu** are not referenced.

*Constraint:* if **range** = 'V', **vl** < **vu**.

8: **il** – INTEGER

9: **iu** – INTEGER

If **range** = 'I', the indices (in ascending order) of the smallest and largest eigenvalues to be returned.

If **range** = 'A' or 'V', **il** and **iu** are not referenced.

*Constraints:*

if **range** = 'I' and **n** = 0, **il** = 1 and **iu** = 0;

if **range** = 'I' and **n** > 0,  $1 \leq \mathbf{il} \leq \mathbf{iu} \leq \mathbf{n}$ .

10: **abstol** – REAL (KIND=nag\_wp)

The absolute error tolerance for the eigenvalues. An approximate eigenvalue is accepted as converged when it is determined to lie in an interval  $[a, b]$  of width less than or equal to

$$\mathbf{abstol} + \epsilon \max(|a|, |b|),$$

where  $\epsilon$  is the *machine precision*. If **abstol** is less than or equal to zero, then  $\epsilon \|T\|_1$  will be used in its place, where  $T$  is the tridiagonal matrix obtained by reducing  $A$  to tridiagonal form. Eigenvalues will be computed most accurately when **abstol** is set to twice the underflow threshold  $2 \times \text{x02am}()$ , not zero. If this function returns with **info** > 0, indicating that some

eigenvectors did not converge, try setting **abstol** to  $2 \times 10^{-24}$ . See Demmel and Kahan (1990).

## 5.2 Optional Input Parameters

1: **n** – INTEGER

*Default:* the second dimension of the array **ab**.

*n*, the order of the matrix *A*.

*Constraint:*  $n \geq 0$ .

## 5.3 Output Parameters

1: **ab**(*ldab*,:) – REAL (KIND=nag\_wp) array

The first dimension of the array **ab** will be **kd** + 1.

The second dimension of the array **ab** will be  $\max(1, n)$ .

**ab** stores values generated during the reduction to tridiagonal form.

The first superdiagonal or subdiagonal and the diagonal of the tridiagonal matrix *T* are returned in **ab** using the same storage format as described above.

2: **q**(*ldq*,:) – REAL (KIND=nag\_wp) array

The first dimension, *ldq*, of the array **q** will be

if **jobz** = 'V',  $ldq = \max(1, n)$ ;  
otherwise  $ldq = 1$ .

The second dimension of the array **q** will be  $\max(1, n)$  if **jobz** = 'V' and 1 otherwise.

If **jobz** = 'V', the *n* by *n* orthogonal matrix used in the reduction to tridiagonal form.

If **jobz** = 'N', **q** is not referenced.

3: **m** – INTEGER

The total number of eigenvalues found.  $0 \leq m \leq n$ .

If **range** = 'A', **m** = **n**.

If **range** = 'I', **m** = **iu** – **il** + 1.

4: **w**(**n**) – REAL (KIND=nag\_wp) array

The first **m** elements contain the selected eigenvalues in ascending order.

5: **z**(*ldz*,:) – REAL (KIND=nag\_wp) array

The first dimension, *ldz*, of the array **z** will be

if **jobz** = 'V',  $ldz = \max(1, n)$ ;  
otherwise  $ldz = 1$ .

The second dimension of the array **z** will be  $\max(1, m)$  if **jobz** = 'V' and 1 otherwise.

If **jobz** = 'V', then

if **info** = 0, the first **m** columns of *Z* contain the orthonormal eigenvectors of the matrix *A* corresponding to the selected eigenvalues, with the *i*th column of *Z* holding the eigenvector associated with **w**(*i*);

if an eigenvector fails to converge (**info** > 0), then that column of *Z* contains the latest approximation to the eigenvector, and the index of the eigenvector is returned in **jfail**.

If **jobz** = 'N', **z** is not referenced.

6: **jfail**(:) – INTEGER array

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

If **jobz** = 'V', then

if **info** = 0, the first **m** elements of **jfail** are zero;

if **info** > 0, **jfail** contains the indices of the eigenvectors that failed to converge.

If **jobz** = 'N', **jfail** is not referenced.

7: **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: **jobz**, 2: **range**, 3: **uplo**, 4: **n**, 5: **kd**, 6: **ab**, 7: **ldab**, 8: **q**, 9: **ldq**, 10: **vl**, 11: **vu**, 12: **il**, 13: **iu**,  
14: **abstol**, 15: **m**, 16: **w**, 17: **z**, 18: **ldz**, 19: **work**, 20: **iwork**, 21: **jfail**, 22: **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 failed to converge. Their indices are stored in array **jfail**. Please see **abstol**.

## 7 Accuracy

The computed eigenvalues and eigenvectors are exact for a nearby matrix  $(A + E)$ , where

$$\|E\|_2 = O(\epsilon)\|A\|_2,$$

and  $\epsilon$  is the *machine precision*. See Section 4.7 of Anderson *et al.* (1999) for further details.

## 8 Further Comments

The total number of floating-point operations is proportional to  $k_d n^2$  if **jobz** = 'N', and is proportional to  $n^3$  if **jobz** = 'V' and **range** = 'A', otherwise the number of floating-point operations will depend upon the number of computed eigenvectors.

The complex analogue of this function is nag\_lapack\_zhbevz (f08hp).

## 9 Example

This example finds the eigenvalues in the half-open interval  $(-3, 3]$ , and the corresponding eigenvectors, of the symmetric band matrix

$$A = \begin{pmatrix} 1 & 2 & 3 & 0 & 0 \\ 2 & 2 & 3 & 4 & 0 \\ 3 & 3 & 3 & 4 & 5 \\ 0 & 4 & 4 & 4 & 5 \\ 0 & 0 & 5 & 5 & 5 \end{pmatrix}.$$

## 9.1 Program Text

```
function f08hb_example

fprintf('f08hb example results\n\n');

% Symmetric band matrix A, stored on symmetric banded format
uplo = 'U';
kd = nag_int(2);
ab = [0, 0, 3, 4, 5;
      0, 2, 3, 4, 5;
      1, 2, 3, 4, 5];

% Eigenvalues in range [-3,3] and corresponding eigenvectors
jobz = 'Vectors';
range = 'Values in range';
vl = -3;
vu = 3;
il = nag_int(0);
iu = nag_int(0);
abstol = 0;
[abf, q, m, w, z, jfail, info] = ...
f08hb( ...
    jobz, range, uplo, kd, ab, vl, vu, il, iu, abstol);

% Normalize eigenvectors: largest element positive
for j = 1:m
    [k] = max(abs(z(:,j)));
    if z(k,j) < 0;
        z(:,j) = -z(:,j);
    end
end

fprintf('Number of eigenvalues in [-3,3] is %2d\n',m);
fprintf('\n Eigenvalues are:\n ');
fprintf(' %7.4f',w(1:m));
fprintf('\n\n');

[ifail] = x04ca( ...
    'General', ' ', z, 'Corresponding eigenvectors of A');
```

## 9.2 Program Results

```
f08hb example results

Number of eigenvalues in [-3,3] is  2

Eigenvalues are:
-2.6633  1.7511

Corresponding eigenvectors of A
      1      2
1  0.6238  0.5635
2 -0.2575 -0.3896
3 -0.5900  0.4008
4  0.4308 -0.5581
5  0.1039  0.2421
```

---