

NAG Toolbox

nag_lapack_dsyevx (f08fb)

1 Purpose

nag_lapack_dsyevx (f08fb) computes selected eigenvalues and, optionally, eigenvectors of a real n by n symmetric matrix A . Eigenvalues and eigenvectors can be selected by specifying either a range of values or a range of indices for the desired eigenvalues.

2 Syntax

```
[a, m, w, z, jfail, info] = nag_lapack_dsyevx(jobz, range, uplo, a, vl, vu, il, iu, abstol, 'n', n)
```

```
[a, m, w, z, jfail, info] = f08fb(jobz, range, uplo, a, vl, vu, il, iu, abstol, 'n', n)
```

3 Description

The symmetric 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: **a**(*lda*,:) – REAL (KIND=nag_wp) array

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

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

The n by n symmetric matrix A .

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.

5: **vl** – REAL (KIND=nag_wp)

6: **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', $\mathbf{vl} < \mathbf{vu}$.

7: **il** – INTEGER

8: **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 $\mathbf{n} = 0$, $\mathbf{il} = 1$ and $\mathbf{iu} = 0$;

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

9: **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 ϵ 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 \text{x02am}(\)$. See Demmel and Kahan (1990).

5.2 Optional Input Parameters

1: **n** – INTEGER

Default: the first dimension of the array **a** and the second dimension of the array **a**. (An error is raised if these dimensions are not equal.)

n , the order of the matrix A .

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

5.3 Output Parameters

1: $\mathbf{a}(\mathit{lda}, :)$ – REAL (KIND=nag_wp) array

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

The second dimension of the array \mathbf{a} will be $\max(1, \mathbf{n})$.

The lower triangle (if $\mathbf{uplo} = 'L'$) or the upper triangle (if $\mathbf{uplo} = 'U'$) of \mathbf{a} , including the diagonal, is overwritten.

2: \mathbf{m} – INTEGER

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

If $\mathbf{range} = 'A'$, $\mathbf{m} = \mathbf{n}$.

If $\mathbf{range} = 'I'$, $\mathbf{m} = \mathbf{i}u - \mathbf{i}l + 1$.

3: $\mathbf{w}(:)$ – REAL (KIND=nag_wp) array

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

The first \mathbf{m} elements contain the selected eigenvalues in ascending order.

4: $\mathbf{z}(\mathit{ldz}, :)$ – REAL (KIND=nag_wp) array

The first dimension, ldz , of the array \mathbf{z} will be

if $\mathbf{jobz} = 'V'$, $\mathit{ldz} = \max(1, \mathbf{n})$;
otherwise $\mathit{ldz} = 1$.

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

If $\mathbf{jobz} = 'V'$, then

if $\mathbf{info} = 0$, the first \mathbf{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 $\mathbf{w}(i)$;

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

If $\mathbf{jobz} = 'N'$, \mathbf{z} is not referenced.

5: $\mathbf{jfail}(:)$ – INTEGER array

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

If $\mathbf{jobz} = 'V'$, then

if $\mathbf{info} = 0$, the first \mathbf{m} elements of \mathbf{jfail} are zero;

if $\mathbf{info} > 0$, \mathbf{jfail} contains the indices of the eigenvectors that failed to converge.

If $\mathbf{jobz} = 'N'$, \mathbf{jfail} is not referenced.

6: \mathbf{info} – INTEGER

$\mathbf{info} = 0$ 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.

info > 0 (*warning*)

The algorithm failed to converge; $\langle value \rangle$ eigenvectors did not converge. Their indices are stored in array **jfail**.

7 Accuracy

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

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

and ϵ 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 n^3 .

The complex analogue of this function is `nag_lapack_zheevx` (f08fp).

9 Example

This example finds the eigenvalues in the half-open interval $(-1, 1]$, and the corresponding eigenvectors, of the symmetric matrix

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

9.1 Program Text

```
function f08fb_example

fprintf('f08fb example results\n\n');

% Eigenvalues between -1 and 1 of A, and corresponding eigenvectors.
a = [1, 2, 3, 4;
     0, 2, 3, 4;
     0, 0, 3, 4;
     0, 0, 0, 4];

jobz = 'Vectors';
range = 'Values in range';
uplo = 'Upper';
vl = -1;
vu = 1;
il = nag_int(0);
iu = nag_int(0);
abstol = 0;
[~, m, w, z, jfail, info] = ...
    f08fb(...
        jobz, range, uplo, a, vl, vu, il, iu, abstol);

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

```
fprintf(' %7.4f',w(1:m));  
fprintf('\n\n');  
[ifail] = x04ca( ...  
    'General', ' ', z, 'Corresponding eigenvectors of A');
```

9.2 Program Results

f08fb example results

Number of eigenvalues in [-1,1] is 2

Eigenvalues are:

-0.5146 -0.2943

Corresponding eigenvectors of A

	1	2
1	-0.5144	0.2767
2	0.4851	-0.6634
3	0.5420	0.6504
4	-0.4543	-0.2457
