# NAG Fortran Library Routine Document F08JJF (SSTEBZ/DSTEBZ)

Note: before using this routine, please read the Users' Note for your implementation to check the interpretation of **bold italicised** terms and other implementation-dependent details.

## 1 Purpose

F08JJF (SSTEBZ/DSTEBZ) computes some (or all) of the eigenvalues of a real symmetric tridiagonal matrix, by bisection.

# 2 Specification

```
SUBROUTINE FO8JJF(RANGE, ORDER, N, VL, VU, IL, IU, ABSTOL, D, E, M,

NSPLIT, W, IBLOCK, ISPLIT, WORK, IWORK, INFO)

ENTRY sstebz (RANGE, ORDER, N, VL, VU, IL, IU, ABSTOL, D, E, M,

NSPLIT, W, IBLOCK, ISPLIT, WORK, IWORK, INFO)

INTEGER N, IL, IU, M, NSPLIT, IBLOCK(*), ISPLIT(*), IWORK(*),

INFO

real VL, VU, ABSTOL, D(*), E(*), W(*), WORK(*)

CHARACTER*1 RANGE, ORDER
```

The ENTRY statement enables the routine to be called by its LAPACK name.

# 3 Description

This routine uses bisection to compute some or all of the eigenvalues of a real symmetric tridiagonal matrix T.

It searches for zero or negligible off-diagonal elements of T to see if the matrix splits into block diagonal form:

It performs bisection on each of the blocks  $T_i$  and returns the block index of each computed eigenvalue, so that a subsequent call to F08JKF (SSTEIN/DSTEIN) to compute eigenvectors can also take advantage of the block structure.

#### 4 References

Kahan W (1966) Accurate eigenvalues of a symmetric tridiagonal matrix Report CS41 Stanford University

#### 5 Parameters

1: RANGE – CHARACTER\*1

Input

On entry: indicates which eigenvalues are required as follows:

if RANGE = 'A', then all the eigenvalues are required;

if RANGE = 'V', then all the eigenvalues in the half-open interval (VL,VU] are required;

if RANGE = 'I', then eigenvalues with indices IL to IU are required.

Constraint: RANGE = 'A', 'V' or 'I'.

#### 2: ORDER - CHARACTER\*1

Input

On entry: indicates the order in which the eigenvalues and their block numbers are to be stored as follows:

if ORDER = 'B', then the eigenvalues are to be grouped by split-off block and ordered from smallest to largest within each block;

if ORDER = 'E', then the eigenvalues for the entire matrix are to be ordered from smallest to largest.

Constraint: ORDER = 'B' or 'E'.

#### 3: N - INTEGER

Input

On entry: n, the order of the matrix T.

Constraint: N > 0.

VL - real 4:

Input Input

VU - real

On entry: if RANGE = 'V', the lower and upper bounds, respectively, of the half-open interval (VL, VU] within which the required eigenvalues lie.

Not referenced if RANGE = 'A' or 'I'.

Constraint: VL < VU if RANGE = 'V'.

6: IL - INTEGER Input

IU - INTEGER

Input

On entry: if RANGE = 'I', the indices of the first and last eigenvalues, respectively, to be computed (assuming that the eigenvalues are in ascending order).

Not referenced if RANGE = 'A' or 'V'.

Constraint:  $1 \le IL \le IU \le N$  if RANGE = 'I'.

ABSTOL - real 8:

Input

On entry: the absolute tolerance to which each eigenvalue is required. An eigenvalue (or cluster) is considered to have converged if it lies in an interval of width  $\leq$  ABSTOL. If ABSTOL  $\leq$  0.0, then the tolerance is taken as **machine precision**  $\times ||T||_1$ .

9: D(\*) - real array Input

**Note:** the dimension of the array D must be at least max(1, N).

On entry: the diagonal elements of the tridiagonal matrix T.

10: E(\*) - real array Input

**Note:** the dimension of the array E must be at least max(1, N - 1).

On entry: the off-diagonal elements of the tridiagonal matrix T.

11: M - INTEGER Output

On exit: m, the actual number of eigenvalues found.

12: NSPLIT - INTEGER Output

On exit: the number of diagonal blocks which constitute the tridiagonal matrix T.

#### 13: W(\*) - real array

Output

**Note:** the dimension of the array W must be at least max(1, N).

On exit: the required eigenvalues of the tridiagonal matrix T stored in W(1) to W(m).

#### 14: IBLOCK(\*) – INTEGER array

Output

**Note:** the dimension of the array IBLOCK must be at least max(1, N).

On exit: at each row/column j where E(j) is zero or negligible, T is considered to split into a block diagonal matrix and IBLOCK(i) contains the block number of the eigenvalue stored in W(i), for  $i=1,2,\ldots,m$ . Note that IBLOCK(i)<0 for some i whenever INFO=1 or 3 (see Section 6) and RANGE=A or A.

#### 15: ISPLIT(\*) – INTEGER array

Output

**Note:** the dimension of the array ISPLIT must be at least max(1, N).

On exit: the leading NSPLIT elements contain the points at which T splits up into sub-matrices as follows. The first sub-matrix consists of rows/columns 1 to ISPLIT(1), the second sub-matrix consists of rows/columns ISPLIT(1) + 1 to ISPLIT(2), ..., and the NSPLIT(th) sub-matrix consists of rows/columns ISPLIT(NSPLIT - 1) + 1 to ISPLIT(NSPLIT) (= n).

#### 16: WORK(\*) - real array

Workspace

**Note:** the dimension of the array WORK must be at least max(1, 4 \* N).

#### 17: IWORK(\*) – INTEGER array

Workspace

**Note:** the dimension of the array IWORK must be at least max(1, 3 \* N).

#### 18: INFO – INTEGER

Output

On exit: INFO = 0 unless the routine detects an error (see Section 6).

### 6 Error Indicators and Warnings

INFO < 0

If INFO = -i, the *i*th parameter had an illegal value. An explanatory message is output, and execution of the program is terminated.

INFO = 1

If RANGE = 'A' or 'V', the algorithm failed to compute some (or all) of the required eigenvalues to the desired accuracy. More precisely, IBLOCK(i) < 0 indicates that the *i*th eigenvalue (stored in W(i)) failed to converge.

INFO = 2

If RANGE = 'I', the algorithm failed to compute some (or all) of the required eigenvalues. Try calling the routine again with RANGE = 'A'.

INFO = 3

If RANGE = 'I', see the description above for INFO = 2.

If RANGE = 'A' or 'V', see the description above for INFO = 1.

INFO = 4

No eigenvalues have been computed. The floating-point arithmetic on the computer is not behaving as expected.

If failures with INFO  $\geq 1$  are causing persistent trouble and the user has checked that the routine is being called correctly, please contact NAG.

# 7 Accuracy

The eigenvalues of T are computed to high relative accuracy which means that if they vary widely in magnitude, then any small eigenvalues will be computed more accurately than, for example, with the standard QR method. However, the reduction to tridiagonal form (prior to calling the routine) may exclude the possibility of obtaining high relative accuracy in the small eigenvalues of the original matrix if its eigenvalues vary widely in magnitude.

## **8** Further Comments

There is no complex analogue of this routine.

# 9 Example

See Section 9 of the document for F08FGF (SORMTR/DORMTR).