

# F08JJFP (PDSTEBZ)

## NAG Parallel Library Routine Document

**Note:** before using this routine, please read the Users' Note for your implementation to check for implementation-dependent details. You are advised to enclose any calls to NAG Parallel Library routines between calls to Z01AAFP and Z01ABFP.

### 1 Description

F08JJFP (PDSTEBZ) computes eigenvalues of a real symmetric tridiagonal matrix  $T$  in parallel by bisection. There are options to request all eigenvalues, all eigenvalues in the interval  $(v_l, v_u]$ , or the eigenvalues indexed  $i_l$  through  $i_u$ .

Each logical processor must contain the vectors  $d$  and  $e$  which are the diagonal and off-diagonal elements of  $T$ , respectively. A static partitioning of the workload is performed at the beginning of this routine which results in all processes finding an (approximately) equal number of eigenvalues.

### 2 Specification

```

SUBROUTINE F08JJFP(ICNTXT, RANGE, ORDER, N, VL, VU, IL, IU,
1                ABSTOL, D, E, M, NSPLIT, W, IBLOCK, ISPLIT,
2                WORK, LWORK, IWORK, LIWORK, INFO)
ENTRY          PDSTEBZ(ICNTXT, RANGE, ORDER, N, VL, VU, IL, IU,
1                ABSTOL, D, E, M, NSPLIT, W, IBLOCK, ISPLIT,
2                WORK, LWORK, IWORK, LIWORK, INFO)
DOUBLE PRECISION VL, VU, ABSTOL, D(*), E(*), W(*), WORK(*)
INTEGER          ICNTXT, N, IL, IU, M, NSPLIT, IBLOCK(*),
1                ISPLIT(*), LWORK, IWORK(*), LIWORK, INFO
CHARACTER*1      RANGE, ORDER

```

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

### 3 Usage

#### 3.1 Definitions

None.

#### 3.2 Global and Local Arguments

The following global **input** arguments must have the same value on entry to the routine on each processor and the global **output** arguments will have the same value on exit from the routine on each processor:

Global input arguments: RANGE, ORDER, N, VL, VU, IL, IU, ABSTOL, D and E

Global output arguments: M, NSPLIT, W, IBLOCK, ISPLIT and INFO

The remaining arguments are local.

#### 3.3 Distribution Strategy

Each logical processor (defined by the Library context ICNTXT) must contain the vectors  $d$  and  $e$  which are the diagonal and off-diagonal entries of  $T$ , respectively.

#### 3.4 Related Routines

The Library provides many support routines for the generation/distribution and input/output of data. The following routines may be used in conjunction with F08JJFP (PDSTEBZ):

Real vector gather: F01ZFPF

## 4 Arguments

- 1:** ICNTXT — INTEGER *Local Input*  
*On entry:* the Library context, usually returned by a call to the Library Grid initialisation routine Z01AAFP.  
**Note:** the value of ICNTXT **must not** be changed.
- 2:** RANGE — CHARACTER\*1 *Global 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  $(v_l, v_u]$  are required;  
 if RANGE = 'I', then eigenvalues with indices  $i_l$  to  $i_u$  are required.  
*Constraint:* RANGE = 'A', 'V' or 'I'.
- 3:** ORDER — CHARACTER\*1 *Global 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'.
- 4:** N — INTEGER *Global Input*  
*On entry:*  $n$ , the order of the matrix  $T$ .  
*Constraint:*  $N \geq 0$ .
- 5:** VL — DOUBLE PRECISION *Global Input*  
**6:** VU — DOUBLE PRECISION *Global Input*  
*On entry:* if RANGE = 'V', the lower and the upper bounds, respectively, of the half-open interval  $(v_l, v_u]$  within which the required eigenvalues lie.  
 Not referenced if RANGE = 'A' or 'I'.  
*Constraint:*  $VL < VU$  if RANGE = 'V'.
- 7:** IL — INTEGER *Global Input*  
**8:** IU — INTEGER *Global Input*  
*On entry:* if RANGE = 'I',  $i_l$  and  $i_u$ , the indices of the first and the last eigenvalues, respectively, to be computed (assuming that the eigenvalues are in non-decreasing order).  
 Not referenced if RANGE = 'A' or 'V'.  
*Constraint:*  $1 \leq IL \leq IU \leq N$  if RANGE = 'I'.
- 9:** ABSTOL — DOUBLE PRECISION *Global 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*  $\times \|T\|_1$ .
- 10:** D(\*) — DOUBLE PRECISION array *Global Input*  
**Note:** the dimension of the array D must be at least  $\max(1, N)$ .  
*On entry:*  $d$ , the diagonal elements of the tridiagonal matrix  $T$ .

- 11:** E(\*) — DOUBLE PRECISION array *Global Input*  
**Note:** the dimension of the array E must be at least  $\max(1, N - 1)$ .  
*On entry:*  $e$ , the off-diagonal elements of the tridiagonal matrix  $T$ .
- 12:** M — INTEGER *Global Output*  
*On exit:* the actual number of eigenvalues found.
- 13:** NSPLIT — INTEGER *Global Output*  
*On exit:* the number of diagonal blocks which constitute the tridiagonal matrix  $T$ .
- 14:** W(\*) — DOUBLE PRECISION array *Global 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).
- 15:** IBLOCK(\*) — INTEGER array *Global 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, the matrix  $T$  is considered to split into a block diagonal matrix and IBLOCK( $i$ ) contains the block number of the eigenvalue stored in W( $i$ ), for  $i = 1, 2, \dots, m$ . Note that IBLOCK( $i$ ) < 0 for some  $i$  whenever INFO = 1 (see Section 5).
- 16:** ISPLIT(\*) — INTEGER array *Global 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 the matrix  $T$  splits up into submatrices (blocks) as follows. The first submatrix consists of rows/columns 1 to ISPLIT(1), the second submatrix consists of rows/columns ISPLIT(1) + 1 to ISPLIT(2), and the last submatrix consists of rows/columns ISPLIT(NSPLIT-1) + 1 to ISPLIT(NSPLIT) (=  $n$ ).
- 17:** WORK(\*) — DOUBLE PRECISION array *Local Workspace*  
**Note:** the dimension of WORK must be at least  $\max(1, LWORK)$ . The minimum value of LWORK required to successfully call this routine can be obtained by setting LWORK = -1. The required size is returned in array element WORK(1).  
*On exit:* WORK(1) contains the minimum dimension of the array WORK required to successfully complete the task.
- 18:** LWORK — INTEGER *Local Input*  
*On entry:* either -1 (see WORK) or the dimension of the array WORK required to successfully complete the task. If LWORK is set to -1 on entry this routine simply performs some initial error checking and then, if these checks are successful, calculates the minimum size of LWORK required.  
*Constraints:*  
 either LWORK = -1  
 or LWORK  $\geq \max(7, 5 \times N)$ .
- 19:** IWORK(\*) — INTEGER array *Local Workspace*  
**Note:** the minimum value of LIWORK required to successfully call this routine can be obtained by setting LIWORK = -1. The required size is returned in array element IWORK(1).  
*On exit:* IWORK(1) contains the minimum dimension of array IWORK required to successfully complete the task.

**20: LIWORK — INTEGER***Local Input*

*On entry:* either  $-1$  (see IWORK) or the dimension of the array WORK required to successfully complete the task. If LIWORK is set to  $-1$  on entry this routine simply performs some initial error checking and then, if these checks are successful, calculates the minimum size of LIWORK required.

*Constraints:*

either  $\text{LIWORK} = -1$   
or  $\text{LIWORK} \geq \max(14, 4 \times N)$ .

**21: INFO — INTEGER***Global Output*

The NAG Parallel Library provides a mechanism, via the routine Z02EAFP, to reduce the amount of parameter validation performed by this routine. For a full description refer to the Z02 Chapter Introduction.

*On exit:*  $\text{INFO} = 0$  (or  $-9999$  if reduced error checking is enabled) unless the routine detects an error (see Section 5).

## 5 Errors and Warnings

If  $\text{INFO} \neq 0$  an explanatory message is output and control returned to the calling program.

$\text{INFO} = -i$

On entry, the  $i$ th argument was invalid. This error occurred either because a global argument did not have the same value on all logical processors, or because its value on one or more processors was incorrect. An explanatory message distinguishes between these two cases.

$\text{INFO} = 1$

The algorithm failed to compute some (or all) of the required eigenvalues to the desired accuracy. More precisely,  $\text{IBLOCK}(i) < 0$  indicates that the  $i$ th eigenvalue (stored in  $W(i)$ ) failed to converge. The effect is that the eigenvalues may not be as accurate as specified by the tolerance.

$\text{INFO} = 2$

There is a mismatch between the number of eigenvalues computed and the desired number.

$\text{INFO} = 3$

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

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

## 6 Further Comments

### 6.1 Algorithmic Detail

The diagonal matrix  $D(\tau)$  of the  $LDL^T$  of the shifted tridiagonal matrix  $T - \tau I$  are computed where  $\tau$  is a shift. The inertia of the matrix  $T - \tau I$  is then given by  $D(\tau)$ . The number of negative values of the diagonal entries of  $D(\tau)$  gives the number of eigenvalues of  $T$  in the interval  $(-\infty, \tau]$  which is the crucial information required to locate eigenvalues. See Kahan [1].

### 6.2 Parallelism Detail

The algorithm is wholly parallel and inter-process communication is minimal.

### 6.3 Accuracy

See Blackford *et al.* [2].

## 7 References

- [1] Kahan W (1966) Accurate eigenvalues of a symmetric tridiagonal matrix *Report CS41* Stanford University
- [2] Blackford L S, Choi J, Cleary A, D’Azevedo E, Demmel J, Dhillon I, Dongarra J, Hammarling S, Henry G, Petitet A, Stanley K, Walker D and Whaley R C (1997) ScaLAPACK Users’ Guide *SIAM* 3600 University City Science Center, Philadelphia, PA 19104-2688, USA. URL: [http://www.netlib.org/scalapack/slug/scalapack\\_slug.html](http://www.netlib.org/scalapack/slug/scalapack_slug.html)

## 8 Example

The example program illustrates the computation of eigenvalues of a 10 by 10 dense symmetric matrix  $A_s$ . The  $(i, j)$ th element of this matrix  $A_s$  is taken to be  $\max(i, j)$  and the matrix is generated using the routine F01ZQFP. The value UPLO = 'U' is used so that only the upper triangular part of  $A_s$  is referenced.

The matrix  $A_s$  is first tridiagonalised using the routine F08FEFP (PDSYTRD).

The diagonal vector  $d$  and the off-diagonal vector  $e$  are gathered to each logical processor by calling the routine F01ZFPF twice.

Finally, the vector  $d$  (denoted locally by the array DL) and the the vector  $e$  (denoted locally by the array EL) are used by the routine F08JJFP (PDSTEBZ) to compute the eigenvalues of the tridiagonal  $T$ .

### 8.1 Example Text

```
*      F08JJFP Example Program Text
*      NAG Parallel Library Release 3 Revised. NAG Copyright 1999.
*      .. Parameters ..
INTEGER          NOUT
PARAMETER       (NOUT=6)
INTEGER          N
PARAMETER       (N=12)
INTEGER          NA
PARAMETER       (NA=20)
INTEGER          MP, NP
PARAMETER       (MP=2, NP=2)
INTEGER          NB
PARAMETER       (NB=2)
INTEGER          LDA, TDA
PARAMETER       (LDA=NA/MP+NB, TDA=NA/NP+NB)
INTEGER          LWORK, LIWORK
PARAMETER       (LWORK=80, LIWORK=80)
*      .. Local Scalars ..
INTEGER          I, IA, ICNTXT, IFAIL, INFO, JA, M, NSPLIT,
+               MPROC, NPROC
LOGICAL          ROOT
CHARACTER        UPLO
*      .. Local Arrays ..
DOUBLE PRECISION A(LDA, TDA), D(TDA), DL(N), E(TDA), EL(N),
+               TAU(TDA), W(N), WORK(LWORK)
INTEGER          IBLOCK(N), IDESCA(9), ISPLIT(N), IWORK(LIWORK)
*      .. External Functions ..
LOGICAL          Z01ACFP
EXTERNAL         Z01ACFP
*      .. External Subroutines ..
EXTERNAL         F01ZFPF, F01ZQFP, F08FEFP, F08JJFP, GMATA,
+               Z01AAFP, Z01ABFP
*      .. Executable Statements ..
```

```

ROOT = Z01ACFP()
IF (ROOT) THEN
  WRITE (NOUT,*) 'F08JJFP Example Program Results'
  WRITE (NOUT,*)
END IF
*
MPROC = MP
NPROC = NP
IFAIL = 0
CALL Z01AAFP(ICNTXT,MPROC,NPROC,IFAIL)
*
* Set the starting address and array descriptor for A
*
IA = 1
JA = 1
IDESCA(1) = 1
IDESCA(2) = ICNTXT
IDESCA(3) = NA
IDESCA(4) = NA
IDESCA(5) = NB
IDESCA(6) = NB
IDESCA(7) = 0
IDESCA(8) = 0
IDESCA(9) = LDA
*
* Generate the matrix A
*
IFAIL = 0
CALL F01ZQFP(GMATA,N,N,A,IA,JA,IDESCA,IFAIL)
*
* Reduce A to tridiagonal form
*
UPLO = 'U'
CALL F08FEFP(UPLO,N,A,IA,JA,IDESCA,D,E,TAU,WORK,LWORK,INFO)
*
* Gather the diagonal D to each logical processor
*
IFAIL = 0
CALL F01ZFPF(N,IA,JA,IDESCA,D,DL,WORK,LWORK,IFAIL)
*
* Gather the off-diagonal E to each logical processor
*
IFAIL = 0
CALL F01ZFPF(N,IA,JA,IDESCA,E,EL,WORK,LWORK,IFAIL)
*
* Compute the eigenvalues of T (and of A)
*
IFAIL = 0
IF (UPLO.EQ.'L') THEN
  CALL F08JJFP(ICNTXT,'A','E',N,0.0D0,0.0D0,0,0,-1.0D0,DL,EL,M,
+           NSPLIT,W,IBLOCK,ISPLIT,WORK,LWORK,IWORK,LIWORK,
+           INFO)
ELSE IF (UPLO.EQ.'U') THEN
  CALL F08JJFP(ICNTXT,'A','E',N,0.0D0,0.0D0,0,0,-1.0D0,DL,EL(2),
+           M,NSPLIT,W,IBLOCK,ISPLIT,WORK,LWORK,IWORK,LIWORK,
+           INFO)
END IF
*

```

```

*      Print the computed eigenvalues
*
      IF (ROOT) THEN
        WRITE (NOUT,*) 'Eigenvalues'
        WRITE (NOUT,*)
        DO 20 I = 1, N
          WRITE (NOUT,*) I, W(I)
20      CONTINUE
      END IF
*
      IFAIL = 0
      CALL Z01ABFP(ICNTXT,'N',IFAIL)
*
      STOP
      END
*
      SUBROUTINE GMATA(I1,I2,J1,J2,AL,LDAL)
*
*      GMATA generates the block A(I1: I2, J1: J2) of the upper
*      triangular part of the matrix A such that
*
*          a(i,j) = max(i,j)
*
*      in the array AL.
*
*      .. Scalar Arguments ..
      INTEGER          I1, I2, J1, J2, LDAL
*      .. Array Arguments ..
      DOUBLE PRECISION AL(LDAL,*)
*      .. Local Scalars ..
      INTEGER          I, J, K, L
*      .. Intrinsic Functions ..
      INTRINSIC        MAX
*      .. Executable Statements ..
      L = 1
      DO 40 J = J1, J2
        K = 1
        DO 20 I = I1, I2
          IF (I.LE.J) THEN
            AL(K,L) = MAX(I,J)
          END IF
          K = K + 1
20      CONTINUE
        L = L + 1
40     CONTINUE
*
      RETURN
*
      END

```

## 8.2 Example Data

None.

### 8.3 Example Results

#### F08JJFP Example Program Results

##### Eigenvalues

1	-18.2653445961538
2	-3.91505976820175
3	-1.74107763047893
4	-1.01013587928877
5	-0.678451626799023
6	-0.501676952612396
7	-0.397980864225928
8	-0.333705083138936
9	-0.293061251985502
10	-0.268013496211147
11	-0.254347826770162
12	105.658854975866

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