

NAG Library Function Document

nag_dpqr (f08jgc)

1 Purpose

nag_dpqr (f08jgc) computes all the eigenvalues and, optionally, all the eigenvectors of a real symmetric positive definite tridiagonal matrix, or of a real symmetric positive definite matrix which has been reduced to tridiagonal form.

2 Specification

```
#include <nag.h>
#include <nagf08.h>

void nag_dpqr (Nag_OrderType order, Nag_ComputeZType compz, Integer n,
              double d[], double e[], double z[], Integer pdz, NagError *fail)
```

3 Description

nag_dpqr (f08jgc) computes all the eigenvalues and, optionally, all the eigenvectors of a real symmetric positive definite tridiagonal matrix T . In other words, it can compute the spectral factorization of T as

$$T = Z\Lambda Z^T,$$

where Λ is a diagonal matrix whose diagonal elements are the eigenvalues λ_i , and Z is the orthogonal matrix whose columns are the eigenvectors z_i . Thus

$$Tz_i = \lambda_i z_i, \quad i = 1, 2, \dots, n.$$

The function may also be used to compute all the eigenvalues and eigenvectors of a real symmetric positive definite matrix A which has been reduced to tridiagonal form T :

$$\begin{aligned} A &= QTQ^T, \text{ where } Q \text{ is orthogonal} \\ &= (QZ)\Lambda(QZ)^T. \end{aligned}$$

In this case, the matrix Q must be formed explicitly and passed to nag_dpqr (f08jgc), which must be called with **compz** = Nag_UpdateZ. The functions which must be called to perform the reduction to tridiagonal form and form Q are:

full matrix	nag_dsytrd (f08fec) and nag_dorgtr (f08ffc)
full matrix, packed storage	nag_dsprtd (f08gec) and nag_dopgtr (f08gfc)
band matrix	nag_dsbrtd (f08hec) with vect = Nag_FormQ.

nag_dpqr (f08jgc) first factorizes T as LDL^T where L is unit lower bidiagonal and D is diagonal. It forms the bidiagonal matrix $B = LD^{\frac{1}{2}}$, and then calls nag_dbdsqr (f08mec) to compute the singular values of B which are the same as the eigenvalues of T . The method used by the function allows high relative accuracy to be achieved in the small eigenvalues of T . The eigenvectors are normalized so that $\|z_i\|_2 = 1$, but are determined only to within a factor ± 1 .

4 References

Barlow J and Demmel J W (1990) Computing accurate eigensystems of scaled diagonally dominant matrices *SIAM J. Numer. Anal.* **27** 762–791

5 Arguments

- 1: **order** – Nag_OrderType *Input*
On entry: the **order** argument specifies the two-dimensional storage scheme being used, i.e., row-major ordering or column-major ordering. C language defined storage is specified by **order** = Nag_RowMajor. See Section 3.2.1.3 in the Essential Introduction for a more detailed explanation of the use of this argument.
Constraint: **order** = Nag_RowMajor or Nag_ColMajor.
- 2: **compz** – Nag_ComputeZType *Input*
On entry: indicates whether the eigenvectors are to be computed.
compz = Nag_NotZ
 Only the eigenvalues are computed (and the array **z** is not referenced).
compz = Nag_InitZ
 The eigenvalues and eigenvectors of T are computed (and the array **z** is initialized by the function).
compz = Nag_UpdateZ
 The eigenvalues and eigenvectors of A are computed (and the array **z** must contain the matrix Q on entry).
Constraint: **compz** = Nag_NotZ, Nag_UpdateZ or Nag_InitZ.
- 3: **n** – Integer *Input*
On entry: n , the order of the matrix T .
Constraint: $n \geq 0$.
- 4: **d**[*dim*] – double *Input/Output*
Note: the dimension, *dim*, of the array **d** must be at least $\max(1, n)$.
On entry: the diagonal elements of the tridiagonal matrix T .
On exit: the n eigenvalues in descending order, unless **fail.code** = NE_CONVERGENCE or NE_POS_DEF, in which case **d** is overwritten.
- 5: **e**[*dim*] – double *Input/Output*
Note: the dimension, *dim*, of the array **e** must be at least $\max(1, n - 1)$.
On entry: the off-diagonal elements of the tridiagonal matrix T .
On exit: **e** is overwritten.
- 6: **z**[*dim*] – double *Input/Output*
Note: the dimension, *dim*, of the array **z** must be at least
 $\max(1, \mathbf{pdz} \times n)$ when **compz** = Nag_UpdateZ or Nag_InitZ and **order** = Nag_ColMajor;
 $\max(1, \times \mathbf{pdz})$ when **compz** = Nag_UpdateZ or Nag_InitZ and **order** = Nag_RowMajor;
 1 when **compz** = Nag_NotZ.
 The (i, j)th element of the matrix Z is stored in
 $\mathbf{z}[(j - 1) \times \mathbf{pdz} + i - 1]$ when **order** = Nag_ColMajor;
 $\mathbf{z}[(i - 1) \times \mathbf{pdz} + j - 1]$ when **order** = Nag_RowMajor.
On entry: if **compz** = Nag_UpdateZ, **z** must contain the orthogonal matrix Q from the reduction to tridiagonal form.
 If **compz** = Nag_InitZ, **z** need not be set.

On exit: if **compz** = Nag_InitZ or Nag_UpdateZ, the n required orthonormal eigenvectors stored as columns of Z ; the i th column corresponds to the i th eigenvalue, where $i = 1, 2, \dots, n$, unless **fail.code** = NE_CONVERGENCE or NE_POS_DEF.

If **compz** = Nag_NotZ, **z** is not referenced.

7: **pdz** – Integer *Input*

On entry: the stride separating row or column elements (depending on the value of **order**) in the array **z**.

Constraints:

if **order** = Nag_ColMajor,
 if **compz** = Nag_InitZ or Nag_UpdateZ, **pdz** \geq max(1, **n**);
 if **compz** = Nag_NotZ, **pdz** \geq 1.;
 if **order** = Nag_RowMajor,
 if **compz** = Nag_UpdateZ or Nag_InitZ, **pdz** \geq max(1, **n**);
 if **compz** = Nag_NotZ, **pdz** \geq 1..

8: **fail** – NagError * *Input/Output*

The NAG error argument (see Section 3.6 in the Essential Introduction).

6 Error Indicators and Warnings

NE_ALLOC_FAIL

Dynamic memory allocation failed.

NE_BAD_PARAM

On entry, argument $\langle value \rangle$ had an illegal value.

NE_CONVERGENCE

The algorithm to compute the singular values of the Cholesky factor B failed to converge; $\langle value \rangle$ off-diagonal elements did not converge to zero.

NE_ENUM_INT_2

On entry, **compz** = $\langle value \rangle$, **pdz** = $\langle value \rangle$ and **n** = $\langle value \rangle$.
 Constraint: if **compz** = Nag_InitZ or Nag_UpdateZ, **pdz** \geq max(1, **n**);
 if **compz** = Nag_NotZ, **pdz** \geq 1.

On entry, **compz** = $\langle value \rangle$, **pdz** = $\langle value \rangle$, **n** = $\langle value \rangle$.
 Constraint: if **compz** = Nag_UpdateZ or Nag_InitZ, **pdz** \geq max(1, **n**);
 if **compz** = Nag_NotZ, **pdz** \geq 1.

NE_INT

On entry, **n** = $\langle value \rangle$.
 Constraint: **n** \geq 0.

On entry, **pdz** = $\langle value \rangle$.
 Constraint: **pdz** $>$ 0.

NE_INTERNAL_ERROR

An internal error has occurred in this function. Check the function call and any array sizes. If the call is correct then please contact NAG for assistance.

NE_POS_DEF

The leading minor of order $\langle value \rangle$ is not positive definite and the Cholesky factorization of T could not be completed. Hence T itself is not positive definite.

7 Accuracy

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

To be more precise, let H be the tridiagonal matrix defined by $H = DTD$, where D is diagonal with $d_{ii} = t_{ii}^{-\frac{1}{2}}$, and $h_{ii} = 1$ for all i . If λ_i is an exact eigenvalue of T and $\tilde{\lambda}_i$ is the corresponding computed value, then

$$|\tilde{\lambda}_i - \lambda_i| \leq c(n)\epsilon\kappa_2(H)\lambda_i$$

where $c(n)$ is a modestly increasing function of n , ϵ is the *machine precision*, and $\kappa_2(H)$ is the condition number of H with respect to inversion defined by: $\kappa_2(H) = \|H\| \cdot \|H^{-1}\|$.

If z_i is the corresponding exact eigenvector of T , and \tilde{z}_i is the corresponding computed eigenvector, then the angle $\theta(\tilde{z}_i, z_i)$ between them is bounded as follows:

$$\theta(\tilde{z}_i, z_i) \leq \frac{c(n)\epsilon\kappa_2(H)}{relgap_i}$$

where $relgap_i$ is the relative gap between λ_i and the other eigenvalues, defined by

$$relgap_i = \min_{i \neq j} \frac{|\lambda_i - \lambda_j|}{(\lambda_i + \lambda_j)}$$

8 Parallelism and Performance

`nag_dpteqr` (f08jgc) is threaded by NAG for parallel execution in multithreaded implementations of the NAG Library.

`nag_dpteqr` (f08jgc) makes calls to BLAS and/or LAPACK routines, which may be threaded within the vendor library used by this implementation. Consult the documentation for the vendor library for further information.

Please consult the Users' Note for your implementation for any additional implementation-specific information.

9 Further Comments

The total number of floating-point operations is typically about $30n^2$ if `compz` = Nag_NotZ and about $6n^3$ if `compz` = Nag_UpdateZ or Nag_InitZ, but depends on how rapidly the algorithm converges. When `compz` = Nag_NotZ, the operations are all performed in scalar mode; the additional operations to compute the eigenvectors when `compz` = Nag_UpdateZ or Nag_InitZ can be vectorized and on some machines may be performed much faster.

The complex analogue of this function is `nag_zpteqr` (f08juc).

10 Example

This example computes all the eigenvalues and eigenvectors of the symmetric positive definite tridiagonal matrix T , where

$$T = \begin{pmatrix} 4.16 & 3.17 & 0.00 & 0.00 \\ 3.17 & 5.25 & -0.97 & 0.00 \\ 0.00 & -0.97 & 1.09 & 0.55 \\ 0.00 & 0.00 & 0.55 & 0.62 \end{pmatrix}.$$

10.1 Program Text

```

/* nag_dpqr (f08jgc) Example Program.
 *
 * Copyright 2001 Numerical Algorithms Group.
 *
 * Mark 7, 2001.
 */

#include <stdio.h>
#include <nag.h>
#include <nag_stdlib.h>
#include <nagf08.h>
#include <nagx04.h>

int main(void)
{
    /* Scalars */
    Integer      i, j, n, pdz, d_len, e_len;
    Integer      exit_status = 0;
    NagError     fail;
    Nag_OrderType order;
    /* Arrays */
    double       *z = 0, *d = 0, *e = 0;

#ifdef NAG_COLUMN_MAJOR
#define Z(I, J) z[(J - 1) * pdz + I - 1]
    order = Nag_ColMajor;
#else
#define Z(I, J) z[(I - 1) * pdz + J - 1]
    order = Nag_RowMajor;
#endif

    INIT_FAIL(fail);

    printf("nag_dpqr (f08jgc) Example Program Results\n\n");

    /* Skip heading in data file */
    scanf("%*[\n] ");
    scanf("%ld%*[\n] ", &n);
    pdz = n;
    d_len = n;
    e_len = n - 1;

    /* Allocate memory */
    if (!(z = NAG_ALLOC(n * n, double)) ||
        !(d = NAG_ALLOC(d_len, double)) ||
        !(e = NAG_ALLOC(e_len, double)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }

    /* Read T from data file */
    for (i = 0; i < d_len; ++i)
        scanf("%lf", &d[i]);
    for (i = 0; i < e_len; ++i)
        scanf("%lf", &e[i]);
    /* Calculate all the eigenvalues and eigenvectors of T */

```

```

/* nag_dpqr (f08jgc).
 * All eigenvalues and eigenvectors of real symmetric
 * positive-definite tridiagonal matrix, reduced from real
 * symmetric positive-definite matrix
 */
nag_dpqr(order, Nag_InitZ, n, d, e, z, pdz, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_dpqr (f08jgc).\n%s\n", fail.message);
    exit_status = 1;
    goto END;
}
/* Normalize the eigenvectors */
for(j=1; j<=n; j++)
{
    for(i=n; i>=1; i--)
    {
        z(i, j) = z(i, j) / z(1,j);
    }
}
/* Print eigenvalues and eigenvectors */
printf(" Eigenvalues\n");
for (i = 0; i < n; ++i)
    printf(" %7.4lf", d[i]);
printf("\n\n");
/* nag_gen_real_mat_print (x04cac).
 * Print real general matrix (easy-to-use)
 */
fflush(stdout);
nag_gen_real_mat_print(order, Nag_GeneralMatrix, Nag_NonUnitDiag, n, n,
                        z, pdz, "Eigenvectors", 0, &fail);
if (fail.code != NE_NOERROR)
{
    printf("Error from nag_gen_real_mat_print (x04cac).\n%s\n",
           fail.message);
    exit_status = 1;
    goto END;
}
END:
NAG_FREE(d);
NAG_FREE(e);
NAG_FREE(z);
return exit_status;
}

```

10.2 Program Data

```

nag_dpqr (f08jgc) Example Program Data
4                               :Value of N
4.16  5.25  1.09  0.62
3.17  -0.97  0.55                :End of matrix T

```

10.3 Program Results

```

nag_dpqr (f08jgc) Example Program Results

```

```

Eigenvalues
8.0023  1.9926  1.0014  0.1237

```

```

Eigenvectors
      1      2      3      4
1      1.0000      1.0000      1.0000      1.0000
2      1.2121     -0.6837     -0.9964     -1.2733
3     -0.1711      0.9721     -1.0962     -3.4611
4     -0.0127      0.3895     -1.5807      3.8354

```
