# **NAG Library Function Document**

# nag\_dggbal (f08whc)

#### 1 Purpose

nag\_dggbal (f08whc) balances a pair of real square matrices (A, B) of order *n*. Balancing usually improves the accuracy of computed generalized eigenvalues and eigenvectors.

#### 2 Specification

```
#include <nag.h>
#include <naqf08.h>
```

```
void nag_dggbal (Nag_OrderType order, Nag_JobType job, Integer n, double a[],
Integer pda, double b[], Integer pdb, Integer *ilo, Integer *ihi,
double lscale[], double rscale[], NagError *fail)
```

#### **3** Description

Balancing may reduce the 1-norms of the matrices and improve the accuracy of the computed eigenvalues and eigenvectors in the real generalized eigenvalue problem

$$Ax = \lambda Bx.$$

nag\_dggbal (f08whc) is usually the first step in the solution of the above generalized eigenvalue problem. Balancing is optional but it is highly recommended.

The term 'balancing' covers two steps, each of which involves similarity transformations on A and B. The function can perform either or both of these steps. Both steps are optional.

1. The function first attempts to permute A and B to block upper triangular form by a similarity transformation:

$$PAP^{\mathrm{T}} = F = \begin{pmatrix} F_{11} & F_{12} & F_{13} \\ & F_{22} & F_{23} \\ & & F_{33} \end{pmatrix}$$

$$PBP^{\mathsf{T}} = G = \begin{pmatrix} G_{11} & G_{12} & G_{13} \\ & G_{22} & G_{23} \\ & & G_{33} \end{pmatrix}$$

where P is a permutation matrix,  $F_{11}$ ,  $F_{33}$ ,  $G_{11}$  and  $G_{33}$  are upper triangular. Then the diagonal elements of the matrix pairs  $(F_{11}, G_{11})$  and  $(F_{33}, G_{33})$  are generalized eigenvalues of (A, B). The rest of the generalized eigenvalues are given by the matrix pair  $(F_{22}, G_{22})$  which are in rows and columns  $i_{lo}$  to  $i_{hi}$ . Subsequent operations to compute the generalized eigenvalues of (A, B) need only be applied to the matrix pair  $(F_{22}, G_{22})$ ; this can save a significant amount of work if  $i_{lo} > 1$  and  $i_{hi} < n$ . If no suitable permutation exists (as is often the case), the function sets  $i_{lo} = 1$  and  $i_{hi} = n$ .

2. The function applies a diagonal similarity transformation to (F, G), to make the rows and columns of  $(F_{22}, G_{22})$  as close in norm as possible:

$$DF\hat{D} = \begin{pmatrix} I & 0 & 0 \\ 0 & D_{22} & 0 \\ 0 & 0 & I \end{pmatrix} \begin{pmatrix} F_{11} & F_{12} & F_{13} \\ F_{22} & F_{23} \\ F_{33} \end{pmatrix} \begin{pmatrix} I & 0 & 0 \\ 0 & \hat{D}_{22} & 0 \\ 0 & 0 & I \end{pmatrix}$$

0 0 *I* 

$$DG\hat{D} = \begin{pmatrix} I & 0 & 0 \\ 0 & D_{22} & 0 \\ 0 & 0 & I \end{pmatrix} \begin{pmatrix} G_{11} & G_{12} & G_{13} \\ & G_{22} & G_{23} \\ & & G_{33} \end{pmatrix} \begin{pmatrix} I & 0 \\ 0 & \hat{D}_{22} \\ 0 & 0 \end{pmatrix}$$

This transformation usually improves the accuracy of computed generalized eigenvalues and eigenvectors.

#### 4 References

f08whc

Ward R C (1981) Balancing the generalized eigenvalue problem SIAM J. Sci. Stat. Comp. 2 141-152

#### 5 Arguments

1: **order** – Nag\_OrderType

*On entry*: the **order** argument specifies the two-dimensional storage scheme being used, i.e., rowmajor 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: **job** – Nag\_JobType

On entry: specifies the operations to be performed on matrices A and B.

- **job** = Nag\_DoNothing No balancing is done. Initialize **ilo** = 1, **ihi** = **n**, **lscale**[i-1] = 1.0 and **rscale**[i-1] = 1.0, for i = 1, 2, ..., n.
- **job** = Nag\_Permute Only permutations are used in balancing.
- **job** = Nag\_Scale Only scalings are are used in balancing.
- job = Nag\_DoBoth

Both permutations and scalings are used in balancing.

Constraint: job = Nag\_DoNothing, Nag\_Permute, Nag\_Scale or Nag\_DoBoth.

3: **n** – Integer

On entry: n, the order of the matrices A and B.

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

4:  $\mathbf{a}[dim] - double$ 

Note: the dimension, *dim*, of the array **a** must be at least  $max(1, pda \times n)$ .

Where A(i, j) appears in this document, it refers to the array element

 $\mathbf{a}[(j-1) \times \mathbf{pda} + i - 1]$  when order = Nag\_ColMajor;

 $\mathbf{a}[(i-1) \times \mathbf{pda} + j - 1]$  when  $\mathbf{order} = \text{Nag}_{\text{RowMajor}}$ .

On entry: the n by n matrix A.

On exit: **a** is overwritten by the balanced matrix. If  $\mathbf{job} = \text{Nag}_D$  on Nothing, **a** is not referenced.

Input/Output

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Input

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

Constraint:  $\mathbf{pda} \ge \max(1, \mathbf{n})$ .

array a.

pda – Integer

5:

6:

 $\mathbf{b}[dim]$  – double Note: the dimension, dim, of the array **b** must be at least max(1, pdb × n).

Where  $\mathbf{B}(i, j)$  appears in this document, it refers to the array element

 $\mathbf{b}[(j-1) \times \mathbf{pdb} + i - 1]$  when order = Nag\_ColMajor;  $\mathbf{b}[(i-1) \times \mathbf{pdb} + j - 1]$  when order = Nag\_RowMajor.

On entry: the n by n matrix B.

On exit: **b** is overwritten by the balanced matrix. If  $\mathbf{job} = \text{Nag}_D$  on Nothing, **b** is not referenced.

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

7: **pdb** – Integer

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

*Constraint*:  $\mathbf{pdb} \ge \max(1, \mathbf{n})$ .

- 8: ilo Integer \*
- 9: ihi Integer \*

On exit:  $i_{lo}$  and  $i_{hi}$  are set such that  $\mathbf{A}(i, j) = 0$  and  $\mathbf{B}(i, j) = 0$  if i > j and  $1 \le j < i_{lo}$  or  $i_{hi} < i \le n$ .

If **job** = Nag\_DoNothing or Nag\_Scale,  $i_{lo} = 1$  and  $i_{hi} = n$ .

10: lscale[n] - double

On exit: details of the permutations and scaling factors applied to the left side of the matrices A and B. If  $P_i$  is the index of the row interchanged with row i and  $d_i$  is the scaling factor applied to row i, then

**Iscale** $[i - 1] = P_i$ , for  $i = 1, 2, ..., i_{lo} - 1$ ; **Iscale** $[i - 1] = d_i$ , for  $i = i_{lo}, ..., i_{hi}$ ; **Iscale** $[i - 1] = P_i$ , for  $i = i_{hi} + 1, ..., n$ .

The order in which the interchanges are made is n to  $i_{\rm hi} + 1$ , then 1 to  $i_{\rm lo} - 1$ .

#### 11: rscale[n] - double

On exit: details of the permutations and scaling factors applied to the right side of the matrices A and B.

If  $P_j$  is the index of the column interchanged with column j and  $\hat{d}_j$  is the scaling factor applied to column j, then

rscale $[j-1] = P_j$ , for  $j = 1, 2, ..., i_{lo} - 1$ ; rscale $[j-1] = \hat{d}_j$ , for  $j = i_{lo}, ..., i_{hi}$ ; rscale $[j-1] = P_j$ , for  $j = i_{hi} + 1, ..., n$ .

The order in which the interchanges are made is n to  $i_{hi} + 1$ , then 1 to  $i_{lo} - 1$ .

# Input/Output

Input/Output

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Output

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Output

Input

# 6 Error Indicators and Warnings

## NE\_ALLOC\_FAIL

Dynamic memory allocation failed. See Section 3.2.1.2 in the Essential Introduction for further information.

## NE\_BAD\_PARAM

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

### NE\_INT

On entry,  $\mathbf{n} = \langle value \rangle$ . Constraint:  $\mathbf{n} \ge 0$ .

On entry,  $\mathbf{pda} = \langle value \rangle$ . Constraint:  $\mathbf{pda} > 0$ .

On entry,  $\mathbf{pdb} = \langle value \rangle$ . Constraint:  $\mathbf{pdb} > 0$ .

## NE\_INT\_2

On entry,  $\mathbf{pda} = \langle value \rangle$  and  $\mathbf{n} = \langle value \rangle$ . Constraint:  $\mathbf{pda} \ge \max(1, \mathbf{n})$ .

On entry,  $\mathbf{pdb} = \langle value \rangle$  and  $\mathbf{n} = \langle value \rangle$ . Constraint:  $\mathbf{pdb} \geq \max(1, \mathbf{n})$ .

## 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.

An unexpected error has been triggered by this function. Please contact NAG. See Section 3.6.6 in the Essential Introduction for further information.

### NE\_NO\_LICENCE

Your licence key may have expired or may not have been installed correctly. See Section 3.6.5 in the Essential Introduction for further information.

# 7 Accuracy

The errors are negligible, compared to those in subsequent computations.

# 8 Parallelism and Performance

nag\_dggbal (f08whc) is not threaded by NAG in any implementation.

nag\_dggbal (f08whc) 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 X06 Chapter Introduction for information on how to control and interrogate the OpenMP environment used within this function. Please also consult the Users' Note for your implementation for any additional implementation-specific information.

# 9 Further Comments

nag\_dggbal (f08whc) is usually the first step in computing the real generalized eigenvalue problem but it is an optional step. The matrix B is reduced to the upper triangular form using the QR factorization function nag\_dgeqrf (f08aec) and this orthogonal transformation Q is applied to the matrix A by calling

nag\_dormqr (f08agc). This is followed by nag\_dgghrd (f08wec) which reduces the matrix pair into the generalized Hessenberg form.

If the matrix pair (A, B) is balanced by this function, then any generalized eigenvectors computed subsequently are eigenvectors of the balanced matrix pair. In that case, to compute the generalized eigenvectors of the original matrix, nag\_dggbak (f08wjc) must be called.

The total number of floating-point operations is approximately proportional to  $n^2$ .

The complex analogue of this function is nag\_zggbal (f08wvc).

# 10 Example

See Section 10 in nag\_dhgeqz (f08xec) and nag\_dtgevc (f08ykc).