# **NAG Toolbox for MATLAB**

# f08wv

### 1 Purpose

f08wv balances a pair of complex square matrices (A,B) of order n. Balancing usually improves the accuracy of computed generalized eigenvalues and eigenvectors.

# 2 Syntax

# 3 Description

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

$$Ax = \lambda Bx$$

f08wv 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:

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

$$PBP^{\mathrm{T}} = G = \left( egin{array}{ccc} G_{11} & G_{12} & G_{13} \ & G_{22} & G_{23} \ & & G_{33} \ \end{array} 
ight)$$

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:

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

$$DGD^{-1} = \begin{pmatrix} I & 0 & 0 \\ 0 & D_{22} & 0 \\ 0 & 0 & I \end{pmatrix} \begin{pmatrix} G_{11} & G_{12} & G_{13} \\ & G_{22} & G_{23} \\ & & G_{33} \end{pmatrix} \begin{pmatrix} I & 0 & 0 \\ 0 & \hat{D}_{22} & 0 \\ 0 & 0 & I \end{pmatrix}$$

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

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#### 4 References

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

#### 5 Parameters

### 5.1 Compulsory Input Parameters

#### 1: **job** – **string**

Specifies the operations to be performed on matrices A and B.

job = 'N'

No balancing is done. Initialize ilo = 1, ihi = n, lscale(i) = 1.0 and rscale(i) = 1.0, for i = 1, ..., n.

job = 'P'

Only permutations are used in balancing.

job = 'S'

Only scalings are are used in balancing.

job = 'B'

Both permutations and scalings are used in balancing.

Constraint: job = 'N', 'P', 'S' or 'B'.

#### 2: a(lda,\*) - complex array

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

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

The n by n matrix A.

#### 3: b(ldb,\*) – complex array

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

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

The n by n matrix B.

### 5.2 Optional Input Parameters

#### 1: n - int32 scalar

Default: The second dimension of the array a The second dimension of the array b.

n, the order of the matrices A and B.

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

# 5.3 Input Parameters Omitted from the MATLAB Interface

lda, ldb, work

### 5.4 Output Parameters

# 1: a(lda,\*) - complex array

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

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

a contains the balanced matrix. If job = 'N', a is not referenced.

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#### 2: b(ldb,\*) - complex array

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

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

**b** contains the balanced matrix. If job = 'N', **b** is not referenced.

- 3: ilo int32 scalar
- 4: ihi int32 scalar

 $i_{lo}$  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 
$$\mathbf{job} = 'N'$$
 or 'S',  $i_{lo} = 1$  and  $i_{hi} = n$ .

### 5: lscale(\*) - double array

**Note**: the dimension of the array **Iscale** must be at least  $max(1, \mathbf{n})$ .

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**(
$$i$$
) =  $P_i$ , for  $i = 1, ..., i_{lo} - 1$ ;

**Iscale**(
$$i$$
) =  $d_i$ , for  $i = i_0, \dots, i_{hi}$ ;

**Iscale**(
$$i$$
) =  $P_i$ , for  $i = i_{hi} + 1, \dots, n$ .

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

# 6: rscale(\*) - double array

**Note**: the dimension of the array **rscale** must be at least  $max(1, \mathbf{n})$ .

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**(
$$j$$
) =  $P_i$ , for  $j = 1, ..., i_{lo} - 1$ ;

$$\mathbf{rscale}(j) = \hat{d}_i$$
, for  $j = i_{10}, \dots, i_{hi}$ ;

$$\mathbf{rscale}(j) = P_i$$
, for  $j = i_{hi} + 1, \dots, n$ .

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

# 7: info – int32 scalar

info = 0 unless the function detects an error (see Section 6).

### 6 Error Indicators and Warnings

Errors or warnings detected by the function:

$$info = -i$$

If info = -i, parameter i had an illegal value on entry. The parameters are numbered as follows:

It is possible that **info** refers to a parameter that is omitted from the MATLAB interface. This usually indicates that an error in one of the other input parameters has caused an incorrect value to be inferred.

### 7 Accuracy

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

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#### **8** Further Comments

f08wv is usually the first step in computing the complex generalized eigenvalue problem but it is an optional step. The matrix B is reduced to the triangular form using the QR factorization function f08as and the unitary transformation Q is applied to the matrix A by calling f08au. This is followed by f08ws 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, f08ww must be called.

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

The real analogue of this function is f08wh.

# 9 Example

```
job = 'B';
a = [complex(1, +3), complex(1, +4), complex(1, +5), complex(1, +6);
    complex(2, +2), complex(4, +3), complex(8, +4), complex(16, +5);
    complex(3, +1), complex(9, +2), complex(27, +3), complex(81, +4);
      complex(4, +0), complex(16, +1), complex(64, +2), complex(256, +3)];
b = [complex(1, +0), complex(2, +1), complex(3, +2), complex(4, +3);
      complex(1, +1), complex(4, +2), complex(9, +3), complex(16, +4);
complex(1, +2), complex(8, +3), complex(27, +4), complex(64, +5);
complex(1, +3), complex(16, +4), complex(81, +5), complex(256, +6)];
[aOut, bOut, ilo, ihi, lscale, rscale, info] = f08wv(job, a, b)
aOut =
     1.0000 + 3.0000i
                               1.0000 + 4.0000i
                                                         0.1000 + 0.5000i
                                                                                    0.1000 +
0.6000i
     2.0000 + 2.0000i
                               4.0000 + 3.0000i
                                                         0.8000 + 0.4000i
                                                                                    1.6000 +
0.5000i
     0.3000 + 0.1000i
                               0.9000 + 0.2000i
                                                         0.2700 + 0.0300i
                                                                                    0.8100 +
0.0400i
                                1.6000 + 0.1000i
                                                          0.6400 + 0.0200i
    0.4000
                                                                                    2.5600 +
0.0300i
bOut =
    1.0000
                                2.0000 + 1.0000i
                                                          0.3000 + 0.2000i
                                                                                    0.4000 +
0.3000i
                               4.0000 + 2.0000i
     1.0000 + 1.0000i
                                                         0.9000 + 0.3000i
                                                                                    1.6000 +
0.4000i
     0.1000 + 0.2000i
                               0.8000 + 0.3000i
                                                         0.2700 + 0.0400i
                                                                                    0.6400 +
0.0500i
     0.1000 + 0.3000i
                               1.6000 + 0.4000i
                                                         0.8100 + 0.0500i
                                                                                    2.5600 +
0.0600i
ilo =
ihi =
lscale =
     1.0000
     1.0000
     0.1000
     0.1000
rscale =
     1.0000
     1.0000
     0.1000
     0.1000
info =
              0
```

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