

NAG Toolbox

nag_lapack_zgesv (f07an)

1 Purpose

nag_lapack_zgesv (f07an) computes the solution to a complex system of linear equations

$$AX = B,$$

where A is an n by n matrix and X and B are n by r matrices.

2 Syntax

```
[a, ipiv, b, info] = nag_lapack_zgesv(a, b, 'n', n, 'nrhs_p', nrhs_p)
[a, ipiv, b, info] = f07an(a, b, 'n', n, 'nrhs_p', nrhs_p)
```

3 Description

nag_lapack_zgesv (f07an) uses the LU decomposition with partial pivoting and row interchanges to factor A as

$$A = PLU,$$

where P is a permutation matrix, L is unit lower triangular, and U is upper triangular. The factored form of A is then used to solve the system of equations $AX = B$.

4 References

Anderson E, Bai Z, Bischof C, Blackford S, Demmel J, Dongarra J J, Du Croz J J, Greenbaum A, Hammarling S, McKenney A and Sorensen D (1999) *LAPACK Users' Guide* (3rd Edition) SIAM, Philadelphia <http://www.netlib.org/lapack/lug>

Golub G H and Van Loan C F (1996) *Matrix Computations* (3rd Edition) Johns Hopkins University Press, Baltimore

5 Parameters

5.1 Compulsory Input Parameters

1: **a**(lda,:) – COMPLEX (KIND=nag_wp) array

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

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

The n by n coefficient matrix A .

2: **b**(ldb,:) – COMPLEX (KIND=nag_wp) array

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

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

The n by r right-hand side matrix B .

5.2 Optional Input Parameters

1: **n** – INTEGER

Default: the first dimension of the array **b**.

n , the number of linear equations, i.e., the order of the matrix A .

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

2: **nrhs_p** – INTEGER

Default: the second dimension of the array **b**.

r , the number of right-hand sides, i.e., the number of columns of the matrix B .

Constraint: **nrhs_p** ≥ 0 .

5.3 Output Parameters

1: **a**(*lda*,:) – COMPLEX (KIND=nag_wp) array

The first dimension of the array **a** will be $\max(1, \mathbf{n})$.

The second dimension of the array **a** will be $\max(1, \mathbf{n})$.

The factors L and U from the factorization $A = PLU$; the unit diagonal elements of L are not stored.

2: **ipiv**(**n**) – INTEGER array

If no constraints are violated, the pivot indices that define the permutation matrix P ; at the i th step row i of the matrix was interchanged with row **ipiv**(i). **ipiv**(i) = i indicates a row interchange was not required.

3: **b**(*ldb*,:) – COMPLEX (KIND=nag_wp) array

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

The second dimension of the array **b** will be $\max(1, \mathbf{nrhs_p})$.

If **info** = 0, the n by r solution matrix X .

4: **info** – INTEGER

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

6 Error Indicators and Warnings

info < 0

If **info** = $-i$, argument i had an illegal value. An explanatory message is output, and execution of the program is terminated.

info > 0 (*warning*)

Element $\langle value \rangle$ of the diagonal is exactly zero. The factorization has been completed, but the factor U is exactly singular, so the solution could not be computed.

7 Accuracy

The computed solution for a single right-hand side, \hat{x} , satisfies the equation of the form

$$(A + E)\hat{x} = b,$$

where

$$\|E\|_1 = O(\epsilon)\|A\|_1$$

and ϵ is the *machine precision*. An approximate error bound for the computed solution is given by

$$\frac{\|\hat{x} - x\|_1}{\|x\|_1} \leq \kappa(A) \frac{\|E\|_1}{\|A\|_1}$$

where $\kappa(A) = \|A^{-1}\|_1 \|A\|_1$, the condition number of A with respect to the solution of the linear equations. See Section 4.4 of Anderson *et al.* (1999) for further details.

Following the use of `nag_lapack_zgesv` (f07an), `nag_lapack_zgecon` (f07au) can be used to estimate the condition number of A and `nag_lapack_zgerfs` (f07av) can be used to obtain approximate error bounds. Alternatives to `nag_lapack_zgesv` (f07an), which return condition and error estimates directly are `nag_linsys_complex_square_solve` (f04ca) and `nag_lapack_zgesvx` (f07ap).

8 Further Comments

The total number of floating-point operations is approximately $\frac{8}{3}n^3 + 8n^2r$, where r is the number of right-hand sides.

The real analogue of this function is `nag_lapack_dgesv` (f07aa).

9 Example

This example solves the equations

$$Ax = b,$$

where A is the general matrix

$$A = \begin{pmatrix} -1.34 + 2.55i & 0.28 + 3.17i & -6.39 - 2.20i & 0.72 - 0.92i \\ -0.17 - 1.41i & 3.31 - 0.15i & -0.15 + 1.34i & 1.29 + 1.38i \\ -3.29 - 2.39i & -1.91 + 4.42i & -0.14 - 1.35i & 1.72 + 1.35i \\ 2.41 + 0.39i & -0.56 + 1.47i & -0.83 - 0.69i & -1.96 + 0.67i \end{pmatrix} \quad \text{and} \quad b = \begin{pmatrix} 26.26 + 51.78i \\ 6.43 - 8.68i \\ -5.75 + 25.31i \\ 1.16 + 2.57i \end{pmatrix}.$$

Details of the LU factorization of A are also output.

9.1 Program Text

```
function f07an_example

fprintf('f07an example results\n\n');

% Linear system Ax = B
a = [ -1.34 + 2.55i, 0.28 + 3.17i, -6.39 - 2.2i, 0.72 - 0.92i;
      -0.17 - 1.41i, 3.31 - 0.15i, -0.15 + 1.34i, 1.29 + 1.38i;
      -3.29 - 2.39i, -1.91 + 4.42i, -0.14 - 1.35i, 1.72 + 1.35i;
      2.41 + 0.39i, -0.56 + 1.47i, -0.83 - 0.69i, -1.96 + 0.67i];
b = [ 26.26 + 51.78i;
      6.43 - 8.68i;
      -5.75 + 25.31i;
      1.16 + 2.57i];

[LU, ipiv, x, info] = f07an(a, b);

disp('Solution');
disp(x');
disp('Details of factorization');
disp(LU);
disp('Pivot indices');
disp(double(ipiv'));
```

9.2 Program Results

f07an example results

Solution

1.0000 - 1.0000i 2.0000 + 3.0000i -4.0000 + 5.0000i 0.0000 - 6.0000i

Details of factorization

-3.2900 - 2.3900i -1.9100 + 4.4200i -0.1400 - 1.3500i 1.7200 + 1.3500i
0.2376 + 0.2560i 4.8952 - 0.7114i -0.4623 + 1.6966i 1.2269 + 0.6190i
-0.1020 - 0.7010i -0.6691 + 0.3689i -5.1414 - 1.1300i 0.9983 + 0.3850i
-0.5359 + 0.2707i -0.2040 + 0.8601i 0.0082 + 0.1211i 0.1482 - 0.1252i

Pivot indices

3 2 3 4
