

NAG Toolbox

nag_fit_glin_l1sol (e02ga)

1 Purpose

nag_fit_glin_l1sol (e02ga) calculates an l_1 solution to an over-determined system of linear equations.

2 Syntax

```
[a, b, x, resid, irank, iter, ifail] = nag_fit_glin_l1sol(a, b, 'm', m, 'nplus2',
nplus2, 'toler', toler)
```

```
[a, b, x, resid, irank, iter, ifail] = e02ga(a, b, 'm', m, 'nplus2', nplus2,
'toler', toler)
```

3 Description

Given a matrix A with m rows and n columns ($m \geq n$) and a vector b with m elements, the function calculates an l_1 solution to the over-determined system of equations

$$Ax = b.$$

That is to say, it calculates a vector x , with n elements, which minimizes the l_1 norm (the sum of the absolute values) of the residuals

$$r(x) = \sum_{i=1}^m |r_i|,$$

where the residuals r_i are given by

$$r_i = b_i - \sum_{j=1}^n a_{ij}x_j, \quad i = 1, 2, \dots, m.$$

Here a_{ij} is the element in row i and column j of A , b_i is the i th element of b and x_j the j th element of x . The matrix A need not be of full rank.

Typically in applications to data fitting, data consisting of m points with coordinates (t_i, y_i) are to be approximated in the l_1 norm by a linear combination of known functions $\phi_j(t)$,

$$\alpha_1\phi_1(t) + \alpha_2\phi_2(t) + \dots + \alpha_n\phi_n(t).$$

This is equivalent to fitting an l_1 solution to the over-determined system of equations

$$\sum_{j=1}^n \phi_j(t_i)\alpha_j = y_i, \quad i = 1, 2, \dots, m.$$

Thus if, for each value of i and j , the element a_{ij} of the matrix A in the previous paragraph is set equal to the value of $\phi_j(t_i)$ and b_i is set equal to y_i , the solution vector x will contain the required values of the α_j . Note that the independent variable t above can, instead, be a vector of several independent variables (this includes the case where each ϕ_i is a function of a different variable, or set of variables).

The algorithm is a modification of the simplex method of linear programming applied to the primal formulation of the l_1 problem (see Barrodale and Roberts (1973) and Barrodale and Roberts (1974)). The modification allows several neighbouring simplex vertices to be passed through in a single iteration, providing a substantial improvement in efficiency.

4 References

Barrodale I and Roberts F D K (1973) An improved algorithm for discrete l_1 linear approximation *SIAM J. Numer. Anal.* **10** 839–848

Barrodale I and Roberts F D K (1974) Solution of an overdetermined system of equations in the l_1 -norm *Comm. ACM* **17(6)** 319–320

5 Parameters

5.1 Compulsory Input Parameters

1: **a**(*lda*, **nplus2**) – REAL (KIND=nag_wp) array

lda, the first dimension of the array, must satisfy the constraint $lda \geq \mathbf{m} + 2$.

a(*i*, *j*) must contain a_{ij} , the element in the *i*th row and *j*th column of the matrix *A*, for $i = 1, 2, \dots, m$ and $j = 1, 2, \dots, n$. The remaining elements need not be set.

2: **b**(**m**) – REAL (KIND=nag_wp) array

b(*i*) must contain b_i , the *i*th element of the vector *b*, for $i = 1, 2, \dots, m$.

5.2 Optional Input Parameters

1: **m** – INTEGER

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

The number of equations, *m* (the number of rows of the matrix *A*).

Constraint: $\mathbf{m} \geq n \geq 1$.

2: **nplus2** – INTEGER

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

$n + 2$, where *n* is the number of unknowns (the number of columns of the matrix *A*).

Constraint: $3 \leq \mathbf{nplus2} \leq \mathbf{m} + 2$.

3: **toler** – REAL (KIND=nag_wp)

Default: 0.0.

A non-negative value. In general **toler** specifies a threshold below which numbers are regarded as zero. The recommended threshold value is $\epsilon^{2/3}$ where ϵ is the *machine precision*. The recommended value can be computed within the function by setting **toler** to zero. If premature termination occurs a larger value for **toler** may result in a valid solution.

5.3 Output Parameters

1: **a**(*lda*, **nplus2**) – REAL (KIND=nag_wp) array

$lda = \mathbf{m} + 2$.

Contains the last simplex tableau generated by the simplex method.

2: **b**(**m**) – REAL (KIND=nag_wp) array

The *i*th residual r_i corresponding to the solution vector *x*, for $i = 1, 2, \dots, m$.

3: **x**(**nplus2**) – REAL (KIND=nag_wp) array

x(*j*) contains the *j*th element of the solution vector *x*, for $j = 1, 2, \dots, n$. The elements **x**($n + 1$) and **x**($n + 2$) are unused.

- 4: **resid** – REAL (KIND=nag_wp)
The sum of the absolute values of the residuals for the solution vector x .
- 5: **irank** – INTEGER
The computed rank of the matrix A .
- 6: **iter** – INTEGER
The number of iterations taken by the simplex method.
- 7: **ifail** – INTEGER
ifail = 0 unless the function detects an error (see Section 5).

6 Error Indicators and Warnings

Errors or warnings detected by the function:

ifail = 1 (*warning*)

An optimal solution has been obtained but this may not be unique.

ifail = 2

The calculations have terminated prematurely due to rounding errors. Experiment with larger values of **toler** or try scaling the columns of the matrix (see Section 9).

ifail = 3

On entry, **nplus2** < 3,
or **nplus2** > **m** + 2,
or **lda** < **m** + 2.

ifail = -99

An unexpected error has been triggered by this routine. Please contact NAG.

ifail = -399

Your licence key may have expired or may not have been installed correctly.

ifail = -999

Dynamic memory allocation failed.

7 Accuracy

Experience suggests that the computational accuracy of the solution x is comparable with the accuracy that could be obtained by applying Gaussian elimination with partial pivoting to the n equations satisfied by this algorithm (i.e., those equations with zero residuals). The accuracy therefore varies with the conditioning of the problem, but has been found generally very satisfactory in practice.

8 Further Comments

The effects of m and n on the time and on the number of iterations in the Simplex Method vary from problem to problem, but typically the number of iterations is a small multiple of n and the total time taken is approximately proportional to mn^2 .

It is recommended that, before the function is entered, the columns of the matrix A are scaled so that the largest element in each column is of the order of unity. This should improve the conditioning of the matrix, and also enable the argument **toler** to perform its correct function. The solution x obtained will

then, of course, relate to the scaled form of the matrix. Thus if the scaling is such that, for each $j = 1, 2, \dots, n$, the elements of the j th column are multiplied by the constant k_j , the element x_j of the solution vector x must be multiplied by k_j if it is desired to recover the solution corresponding to the original matrix A .

9 Example

Suppose we wish to approximate a set of data by a curve of the form

$$y = Ke^t + Le^{-t} + M$$

where K , L and M are unknown. Given values y_i at 5 points t_i we may form the over-determined set of equations for K , L and M

$$e^{x_i}K + e^{-x_i}L + M = y_i, \quad i = 1, 2, \dots, 5.$$

nag_fit_glin_llsol (e02ga) is used to solve these in the l_1 sense.

9.1 Program Text

```
function e02ga_example

fprintf('e02ga example results\n\n');

a = zeros(7, 5);
for i = 1:5
    a(i, 1) = exp((i-1)/5);
    a(i, 2) = exp(-(i-1)/5);
    a(i, 3) = 1;
end
b = [4.501; 4.36; 4.333; 4.418; 4.625];

[a, b, x, resid, irank, iter, ifail] = ...
    e02ga(a, b);

fprintf('Resid = %8.4f Rank = %5d Iterations = %5d\n\n', resid, irank, iter);
disp('Solution:');
disp(x(1:irank)');
```

9.2 Program Results

```
e02ga example results

Resid = 0.0028 Rank = 3 Iterations = 5

Solution:
1.0014 2.0035 1.4960
```
