

nag_2d_spline_fit_scatter (e02ddc)

1. Purpose

nag_2d_spline_fit_scatter (e02ddc) computes a bicubic spline approximation to a set of scattered data. The knots of the spline are located automatically, but a single parameter must be specified to control the trade-off between closeness of fit and smoothness of fit.

2. Specification

```
#include <nag.h>
#include <nage02.h>

void nag_2d_spline_fit_scatter(Nag_Start start, Integer m, double x[],
    double y[], double f[], double weights[], double s,
    Integer nxest, Integer nyest, double *fp, Integer *rank,
    double *warmstartinf, Nag_2dSpline *spline,
    NagError *fail)
```

3. Description

This function determines a smooth bicubic spline approximation $s(x, y)$ to the set of data points (x_r, y_r, f_r) with weights w_r , for $r = 1, 2, \dots, m$. The approximation domain is considered to be the rectangle $[x_{\min}, x_{\max}] \times [y_{\min}, y_{\max}]$, where x_{\min} (y_{\min}) and x_{\max} (y_{\max}) denote the lowest and highest data values of $x(y)$.

The spline is given in the B-spline representation

$$s(x, y) = \sum_{i=1}^{n_x-4} \sum_{j=1}^{n_y-4} c_{ij} M_i(x) N_j(y), \quad (1)$$

where $M_i(x)$ and $N_j(y)$ denote normalised cubic B-splines, the former defined on the knots λ_i to λ_{i+4} and the latter on the knots μ_j to μ_{j+4} . For further details, see Hayes and Halliday (1974) for bicubic splines and de Boor (1972) for normalised B-splines.

The total numbers n_x and n_y of these knots and their values $\lambda_1, \dots, \lambda_{n_x}$ and μ_1, \dots, μ_{n_y} are chosen automatically by the function. The knots $\lambda_5, \dots, \lambda_{n_x-4}$ and $\mu_5, \dots, \mu_{n_y-4}$ are the interior knots; they divide the approximation domain $[x_{\min}, x_{\max}] \times [y_{\min}, y_{\max}]$ into $(n_x - 7) \times (n_y - 7)$ subpanels $[\lambda_i, \lambda_{i+1}] \times [\mu_j, \mu_{j+1}]$, for $i = 4, 5, \dots, n_x - 4$, $j = 4, 5, \dots, n_y - 4$. Then, much as in the curve case (see nag_1d_spline_fit (e02bec)); the coefficients c_{ij} are determined as the solution of the following constrained minimization problem:

minimize

$$\eta, \quad (2)$$

subject to the constraint

$$\theta = \sum_{r=1}^m \varepsilon_r^2 \leq S, \quad (3)$$

where

η is a measure of the (lack of) smoothness of $s(x, y)$. Its value depends on the discontinuity jumps in $s(x, y)$ across the boundaries of the subpanels. It is zero only when there are no discontinuities and is positive otherwise, increasing with the size of the jumps (see Dierckx (1981a) for details). ε_r denotes the weighted residual $w_r(f_r - s(x_r, y_r))$, and S is a non-negative number to be specified by the user.

By means of the parameter S , ‘the smoothing factor’, the user will then control the balance between smoothness and closeness of fit, as measured by the sum of squares of residuals in (3). If S is too large, the spline will be too smooth and signal will be lost (underfit); if S is too small, the spline will pick up too much noise (overfit). In the extreme cases the method would return an interpolating

spline ($\theta = 0$) if S were set to zero, and the least-squares bicubic polynomial ($\eta = 0$) if S is set very large. Experimenting with S values between these two extremes should result in a good compromise. (See Section 6.3 for advice on choice of S .) Note however, that this function, unlike `nag_1d_spline_fit` (e02bec) and `nag_1d_spline_deriv` (e02bcc), does not allow S to be set exactly to zero.

The method employed is outlined in Section 6.6 and fully described in Dierckx (1981a) and Dierckx (1981b). It involves an adaptive strategy for locating the knots of the bicubic spline (depending on the function underlying the data and on the value of S), and an iterative method for solving the constrained minimization problem once the knots have been determined.

Values of the computed spline can subsequently be computed by calling `nag_2d_spline_eval` (e02dec) or `nag_2d_spline_eval_rect` (e02dfc) as described in Section 6.7.

4. Parameters

start

Input: **start** must be set to **Nag_Cold** or **Nag_Warm**

If **start** = **Nag_Cold** (cold start), the function will build up the knot set starting with no interior knots. No values need be assigned to **spline.nx** and **spline.ny** and memory will be internally allocated to **spline.lamda**, **spline.mu** and **spline.c**

If **start** = **Nag_Warm** (warm start), the function will restart the knot-placing strategy using the knots found in a previous call of the routine. In this case, all parameters except **s** must be unchanged from that previous call. This warm start can save much time in searching for a satisfactory value of S .

Constraint: **start** = **Nag_Cold** or **Nag_Warm**.

m

Input: m , the number of data points.

The number of data points with non-zero weight (see **weights** below) must be at least 16.

x[m]

y[m]

f[m]

Input: **x**[$r - 1$], **y**[$r - 1$], **f**[$r - 1$] must be set to the co-ordinates of (x_r, y_r, f_r) , the r th data point, for $r = 1, 2, \dots, m$. The order of the data points is immaterial.

weights[m]

Input: **weights**[$r - 1$] must be set to w_r , the r th value in the set of weights, for $r = 1, 2, \dots, m$. Zero weights are permitted and the corresponding points are ignored, except when determining x_{\min} , x_{\max} , y_{\min} and y_{\max} (see Section 6.5). For advice on the choice of weights, see the Chapter Introduction.

Constraint: the number of data points with non-zero weight must be at least 16.

s

Input: the smoothing factor, S . For advice on the choice of S , see Section 3 and Section 6.3.

Constraint: $s > 0.0$.

nxest

nyest

Input: an upper bound for the number of knots n_x and n_y required in the x and y directions respectively. In most practical situations, **nxest** = **nyest** = $5 + \sqrt{m}$ is sufficient. See also Section 6.4.

Constraints: **nxest** ≥ 8 and **nyest** ≥ 8 .

fp

Output: the weighted sum of squared residuals, θ , of the computed spline approximation. **fp** should equal S within a relative tolerance of 0.001 unless **spline.nx** = **spline.ny** = 8, when the spline has no interior knots and so is simply a bicubic polynomial. For knots to be inserted, S must be set to a value below the value of **fp** produced in this case.

rank

Output: **rank** gives the rank of the system of equations used to compute the final spline (as determined by a suitable machine-dependent threshold). When **rank** = (**spline.nx**−4) × (**spline.ny**−4), the solution is unique; otherwise the system is rank-deficient and the minimum-norm solution is computed. The latter case may be caused by too small a value of *S*.

warmstartinf

Output: if the warm start option is used, its value must be left unchanged from the previous call.

spline

Input/Output: Pointer to structure of type Nag_2dSpline with the following members:

nx - Integer

Input: if the warm start option is used, the value of **nx** must be left unchanged from the previous call.

Output: the total number of knots, n_x , of the computed spline with respect to the *x* variable.

lamda - double *

Input: a pointer to which if **start** = **Nag_Cold**, memory of size **nxest** is internally allocated. If the warm start option is used, the values **spline.lamda**[0], **spline.lamda**[1], ..., **spline.lamda**[**spline.nx**−1] must be left unchanged from the previous call.

Output: **spline.lamda** contains the complete set of knots λ_i associated with the *x* variable, i.e. the interior knots **spline.lamda**[4], **spline.lamda**[5], ..., **spline.lamda**[**spline.nx**−5] as well as the additional knots **spline.lamda**[0] = **spline.lamda**[1] = **spline.lamda**[2] = **spline.lamda**[3] = x_{\min} and **spline.lamda**[**spline.nx**−4] = **spline.lamda**[**spline.nx**−3] = **spline.lamda**[**spline.nx**−2] = **spline.lamda**[**spline.nx**−1] = x_{\max} needed for the B-spline representation (where x_{\min} and x_{\max} are as described in Section 3).

ny - Integer

Input: if the warm start option is used, the value of **ny** must be left unchanged from the previous call.

Output: the total number of knots, n_y , of the computed spline with respect to the *y* variable.

mu - double *

Input: a pointer to which if **start** = **Nag_Cold**, memory of size **nyest** is internally allocated. If the warm start option is used, the values **spline.mu**[0], **spline.mu**[1], ..., **spline.mu**[**spline.ny**−1] must be left unchanged from the previous call.

Output: **spline.mu** contains the complete set of knots μ_i associated with the *y* variable, i.e. the interior knots **spline.mu**[4], **spline.mu**[5], ..., **spline.mu**[**spline.ny**−5] as well as the additional knots **spline.mu**[0] = **spline.mu**[1] = **spline.mu**[2] = **spline.mu**[3] = y_{\min} and **spline.mu**[**spline.ny**−4] = **spline.mu**[**spline.ny**−3] = **spline.mu**[**spline.ny**−2] = **spline.mu**[**spline.ny**−1] = y_{\max} needed for the B-spline representation (where y_{\min} and y_{\max} are as described in Section 3).

c - double *

Output: a pointer to which, if **start** = **Nag_Cold**, memory of size (**nxest**−4) × (**nyest**−4) is internally allocated. **spline.c**[($n_y - 4$) × ($i - 1$) + $j - 1$] is the coefficient c_{ij} defined in Section 3.

fail

The NAG error parameter, see the Essential Introduction to the NAG C Library.

5. Error Indications and Warnings

NE_BAD_PARAM

On entry, parameter **start** had an illegal value.

NE_INT_ARG_LT

On entry, **nxest** must not be less than 8: **nxest** = *<value>*.

On entry, **nyest** must not be less than 8: **nyest** = *<value>*.

NE_REAL_ARG_LE

On entry, **s** must not be less than or equal to 0.0: **s** = $\langle \text{value} \rangle$.

NE_NON_ZERO_WEIGHTS

On entry, the number of data points with non-zero weights = $\langle \text{value} \rangle$. Constraint: the number of non-zero weights ≥ 16 .

NE_ALLOC_FAIL

Memory allocation failed.

NE_ENUMTYPE_WARM

start has been set to **Nag_Warm** at the first call of this function. It must be set to **Nag_Cold** at the first call.

NE_ALL_ELEMENTS_EQUAL

On entry, all the values in the array **x** must not be equal.

On entry, all the values in the array **y** must not be equal.

NE_NUM_KNOTS_2D_GT_SCAT

The number of knots required is greater than allowed by **nxest** or **nyest**, **nxest** = $\langle \text{value} \rangle$, **nyest** = $\langle \text{value} \rangle$. Possibly **s** is too small, especially if **nxest**, **nyest** $> 5 + \sqrt{\mathbf{m}}$. **s** = $\langle \text{value} \rangle$, **m** = $\langle \text{value} \rangle$.

NE_NUM_COEFF_GT

No more knots can be added because the number of B-spline coefficients already exceeds **m**. Either **m** or **s** is probably too small: **m** = $\langle \text{value} \rangle$, **s** = $\langle \text{value} \rangle$.

NE_NO_ADDITIONAL_KNOTS

No more knots added; the additional knot would coincide with an old one. Possibly an inaccurate data point has too large a weight, or **s** is too small. **s** = $\langle \text{value} \rangle$.

NE_SPLINE_COEFF_CONV

The iterative process has failed to converge. Possibly **s** is too small: **s** = $\langle \text{value} \rangle$.

If the function fails with an error exit of **NE_NUM_KNOTS_2D_GT_SCAT**, **NE_NUM_COEFF_GT**, **NE_NO_ADDITIONAL_KNOTS** or **NE_SPLINE_COEFF_CONV**, then a spline approximation is returned, but it fails to satisfy the fitting criterion (see (2) and (3) in Section 3) – perhaps by only a small amount, however.

6. Further Comments

6.1. Accuracy

On successful exit, the approximation returned is such that its weighted sum of squared residuals **fp** is equal to the smoothing factor S , up to a specified relative tolerance of 0.001 – except that if $n_x = 8$ and $n_y = 8$, **fp** may be significantly less than S : in this case the computed spline is simply the least-squares bicubic polynomial approximation of degree 3, i.e. a spline with no interior knots.

6.2. Timing

The time taken for a call of `nag_2d_spline_fit_scatter` depends on the complexity of the shape of the data, the value of the smoothing factor S , and the number of data points. If `nag_2d_spline_fit_scatter` is to be called for different values of S , much time can be saved by setting **start** = **Nag_Warm** after the first call.

It should be noted that choosing S very small considerably increases computation time.

6.3. Choice of S

If the weights have been correctly chosen (see the Chapter Introduction), the standard deviation of $w_r f_r$ would be the same for all r , equal to σ , say. In this case, choosing the smoothing factor S in the range $\sigma^2(m \pm \sqrt{2m})$, as suggested by Reinsch (1967), is likely to give a good start in the search for a satisfactory value. Otherwise, experimenting with different values of S will be required from the start.

In that case, in view of computation time and memory requirements, it is recommended to start with a very large value for S and so determine the least-squares bicubic polynomial; the value

returned for **fp**, call it **fp**₀, gives an upper bound for S . Then progressively decrease the value of S to obtain closer fits – say by a factor of 10 in the beginning, i.e. $S = \mathbf{fp}_0/10$, $S = \mathbf{fp}_0/100$, and so on, and more carefully as the approximation shows more details.

To choose S very small is strongly discouraged. This considerably increases computation time and memory requirements. It may also cause rank-deficiency (as indicated by the parameter **rank**) and endanger numerical stability.

The number of knots of the spline returned, and their location, generally depend on the value of S and on the behaviour of the function underlying the data. However, if `nag_2d_spline_fit_sc` is called with **start** = **Nag_Warm**, the knots returned may also depend on the smoothing factors of the previous calls. Therefore if, after a number of trials with different values of S and **start** = **Nag_Warm**, a fit can finally be accepted as satisfactory, it may be worthwhile to call `nag_2d_spline_fit_sc` once more with the selected value for S but now using **start** = **Nag_Cold**. Often, `nag_2d_spline_fit_sc` then returns an approximation with the same quality of fit but with fewer knots, which is therefore better if data reduction is also important.

6.4. Choice of **nxest** and **nyest**

The number of knots may also depend on the upper bounds **nxest** and **nyest**. Indeed, if at a certain stage in `nag_2d_spline_fit_sc` the number of knots in one direction (say n_x) has reached the value of its upper bound (**nxest**), then from that moment on all subsequent knots are added in the other (y) direction. This may indicate that the value of **nxest** is too small. On the other hand, it gives the user the option of limiting the number of knots the function locates in any direction. For example, by setting **nxest** = 8 (the lowest allowable value for **nxest**), the user can indicate that he wants an approximation which is a simple cubic polynomial in the variable x .

6.5. Restriction of the Approximation Domain

The fit obtained is not defined outside the rectangle $[\lambda_4, \lambda_{n_x-3}] \times [\mu_4, \mu_{n_y-3}]$. The reason for taking the extreme data values of x and y for these four knots is that, as is usual in data fitting, the fit cannot be expected to give satisfactory values outside the data region. If, nevertheless, the user requires values over a larger rectangle, this can be achieved by augmenting the data with two artificial data points $(a, c, 0)$ and $(b, d, 0)$ with zero weight, where $[a, b] \times [c, d]$ denotes the enlarged rectangle.

6.6. Outline of Method Used

First suitable knot sets are built up in stages (starting with no interior knots in the case of a cold start but with the knot set found in a previous call if a warm start is chosen). At each stage, a bicubic spline is fitted to the data by least-squares and θ , the sum of squares of residuals, is computed. If $\theta > S$, a new knot is added to one knot set or the other so as to reduce θ at the next stage. The new knot is located in an interval where the fit is particularly poor. Sooner or later, we find that $\theta \leq S$ and at that point the knot sets are accepted. The function then goes on to compute a spline which has these knot sets and which satisfies the full fitting criterion specified by (2) and (3). The theoretical solution has $\theta = S$. The function computes the spline by an iterative scheme which is ended when $\theta = S$ within a relative tolerance of 0.001. The main part of each iteration consists of a linear least-squares computation of special form. The minimal least-squares solution is computed wherever the linear system is found to be rank-deficient.

An exception occurs when the function finds at the start that, even with no interior knots ($n_x = n_y = 8$), the least-squares spline already has its sum of squares of residuals $\leq S$. In this case, since this spline (which is simply a bicubic polynomial) also has an optimal value for the smoothness measure η , namely zero, it is returned at once as the (trivial) solution. It will usually mean that S has been chosen too large.

For further details of the algorithm and its use see Dierckx (1981a).

6.7. Evaluation of Computed Spline

The values of the computed spline at the points $(\mathbf{tx}[r-1], \mathbf{ty}[r-1])$, for $r = 1, 2, \dots, \mathbf{n}$, may be obtained in the array **ff**, of length at least **n**, by the following code:

```
e02dec(n, tx, ty, ff, &spline, &fail)
```

where **spline** is a structure of type **Nag_2dSpline** which is an output parameter of `nag_2d_spline_fit_sc`.

To evaluate the computed spline on a **kx** by **ky** rectangular grid of points in the x - y plane, which is defined by the x co-ordinates stored in **tx**[$q - 1$], for $q = 1, 2, \dots, \mathbf{kx}$, and the y co-ordinates stored in **ty**[$r - 1$], for $r = 1, 2, \dots, \mathbf{ky}$, returning the results in the array **fg** which is of length at least $\mathbf{kx} \times \mathbf{ky}$, the following call may be used:

```
e02dfc(kx, ky, tx, ty, fg, &spline, &fail)
```

where **spline** is a structure of type **Nag_2dSpline** which is an output parameter of nag_2d_spline_fit_scatter. The result of the spline evaluated at grid point (q, r) is returned in element $[\mathbf{ky} \times (q - 1) + r - 1]$ of the array **fg**.

6.8. References

- De Boor C (1972) On Calculating with B-splines *J. Approx. Theory* **6** 50–62.
 Dierckx P (1981a) An Algorithm for Surface Fitting with Spline Functions *IMA J. Num. Anal.* **1** 267–283.
 Dierckx P (1981b) *An Improved Algorithm for Curve Fitting with Spline Functions* Dept. Computer Science, K.U.Leuven, Report TW54.
 Hayes J G and Halliday J (1974) The Least-squares Fitting of Cubic Spline Surfaces to General Data Sets *J. Inst. Math. Appl.* **14** 89–103.
 Peters G and Wilkinson J H (1970) The Least-squares Problem and Pseudo-inverses *Comput. J.* **13** 309–316.
 Reinsch C H (1967) Smoothing by Spline Functions *Num. Math.* **10** 177–183.

7. See Also

```
nag_2d_spline_interpolant (e01dac)
nag_1d_spline_fit (e02bec)
nag_2d_spline_fit_grid (e02dcc)
nag_2d_spline_eval (e02dec)
nag_2d_spline_eval_rect (e02dfc)
```

8. Example

This example program reads in a value of **m**, followed by a set of **m** data points (x_r, y_r, f_r) and their weights w_r . It then calls nag_2d_spline_fit_scatter to compute a bicubic spline approximation for one specified value of S, and prints the values of the computed knots and B-spline coefficients. Finally it evaluates the spline at a small sample of points on a rectangular grid.

8.1. Program Text

```
/* nag_2d_spline_fit_scatter(e02ddc) Example Program
 *
 * Copyright 1991 Numerical Algorithms Group.
 *
 * Mark 2, 1991.
 */

#include <math.h>;
#include <stdio.h>;
#include <stdlib.h>;
#include "e02.h";

#define MMAX 30
#define NMAX 7
main()
{
    double f[MMAX], weights[MMAX], x[MMAX], y[MMAX], fg[NMAX*NMAX],
    px[NMAX], py[NMAX];
    Integer i, j, m, rank, nx, ny, npx, npy, nxest, nyest;
    double s, delta, fp, xhi, yhi, xlo, ylo;
    Nag_Start start;
    Nag_2dSpline spline;
    double warmstartinf;

    nxest = 14;
```

```

nyest = 14;
Vprintf("e02ddc Example Program Results\n");
Vscanf("%*[\n]"); /* Skip heading in data file */
/* Input the number of data-points m. */
Vscanf("%ld",&m);
if (m>0 && m<=MMAX)
{
    /* Input the data-points and the weights. */
    for (i=0; i<m; i++)
        Vscanf("%lf%lf%lf%lf",&x[i],&y[i],&f[i],&weights[i]);
    start = Nag_Cold;
    if (scanf("%lf",&s) != EOF)
    {
        /* Determine the spline approximation. */

        e02ddc(start, m, x, y, f, weights, s, nxest, nyest, &fp,
            &rank, &warmstartinf, &spline, NAGERR_DEFAULT);
        nx = spline.nx;
        ny = spline.ny;
        Vprintf("\nCalling with smoothing factor s = %11.4e, nx = %ld,\n",
            ny = %ld\n",s,nx,ny);

        Vprintf("rank deficiency = %ld\n\n",(nx-4)*(ny-4)-rank);

        /* Print the knot sets, lamda and mu. */
        Vprintf("Distinct knots in x direction located at\n");
        for (j=3; j<spline.nx-3; j++)
            Vprintf("%12.4f%s",spline.lamda[j],((j-3)%5==4 || j==spline.nx-4)
                ? "\n" : " ");
        Vprintf("\nDistinct knots in y direction located at\n");
        for (j=3; j<spline.ny-3; j++)
            Vprintf("%12.4f%s",spline.mu[j],((j-3)%5==4 || j==spline.ny-4)
                ? "\n" : " ");
        Vprintf("\nThe B-spline coefficients:\n\n");
        for (i=0; i<ny-4; i++)
        {
            for (j=0; j<nx-4; j++)
                Vprintf("%9.2f",spline.c[i+j*(ny-4)]);
            Vprintf("\n");
        }

        Vprintf("\n Sum of squared residuals fp = %11.4e\n",fp);
        if (nx==8 && ny==8)
            Vprintf("The spline is the least-squares bi-cubic polynomial\n");

        /* Evaluate the spline on a rectangular grid at npy*nx points
         * over the domain (xlo to xhi) x (ylo to yhi).
         */

        Vscanf("%ld%lf%lf",&npx,&xlo,&xhi);
        Vscanf("%ld%lf%lf",&npy,&ylo,&yhi);

        if (npx<=NMAX && npy<=NMAX)
        {
            delta = (xhi-xlo)/(npx-1);
            for (i=0; i<npx; i++)
                px[i] = MIN(xhi,xlo+i*delta);
            for (i=0; i<npy; i++)
                py[i] = MIN(yhi,ylo+i*delta);

            e02dfc(npx, npy, px, py, fg, &spline, NAGERR_DEFAULT);

            Vprintf("\nValues of computed spline:\n\n");
            Vprintf("      x");
            for (i=0; i<npx; Vprintf("%8.2f ",px[i]), i++);
            Vprintf("\n      y\n");
            for (i=npy-1; i>=0; i--)
            {
                Vprintf("%8.2f ",py[i]);
                for (j=0; j<npx; Vprintf("%8.2f ",fg[npy*j+i]), j++);
            }
        }
    }
}

```

```

        }
        }
    }
    else
    {
        Vfprintf(stderr,"npx or npy is out of range: npx = %4ld,\
npy = %4ld\n",npx,npy);
        exit(EXIT_FAILURE);
    }
    exit(EXIT_SUCCESS);
}
else
{
    Vfprintf(stderr,"m is out of range: m = %4ld,\n",m);
    exit(EXIT_FAILURE);
}
}

```

8.2. Program Data

e02ddc Example Program Data

```

30
11.16  1.24 22.15 1.00
12.85  3.06 22.11 1.00
19.85 10.72 7.97 1.00
19.72  1.39 16.83 1.00
15.91  7.74 15.30 1.00
 0.00 20.00 34.60 1.00
20.87 20.00  5.74 1.00
 3.45 12.78 41.24 1.00
14.26 17.87 10.74 1.00
17.43  3.46 18.60 1.00
22.80 12.39  5.47 1.00
 7.58  1.98 29.87 1.00
25.00 11.87  4.40 1.00
 0.00  0.00 58.20 1.00
 9.66 20.00  4.73 1.00
 5.22 14.66 40.36 1.00
17.25 19.57  6.43 1.00
25.00  3.87  8.74 1.00
12.13 10.79 13.71 1.00
22.23  6.21 10.25 1.00
11.52  8.53 15.74 1.00
15.20  0.00 21.60 1.00
 7.54 10.69 19.31 1.00
17.32 13.78 12.11 1.00
 2.14 15.03 53.10 1.00
 0.51  8.37 49.43 1.00
22.69 19.63  3.25 1.00
 5.47 17.13 28.63 1.00
21.67 14.36  5.52 1.00
 3.31  0.33 44.08 1.00
10.0
 7  3.0 21.0
 6  2.0 17.0

```

8.3. Program Results

e02ddc Example Program Results

Calling with smoothing factor s = 1.0000e+01, nx = 10, ny = 9
rank deficiency = 0

Distinct knots in x direction located at
0.0000 9.7575 18.2582 25.0000

Distinct knots in y direction located at
0.0000 9.0008 20.0000

The B-spline coefficients:

58.16	46.31	6.01	32.00	5.86	-23.78
63.78	46.74	33.37	18.30	14.36	15.95
40.84	-33.79	5.17	13.10	-4.13	19.37
75.44	111.92	6.94	17.33	7.09	-13.24
34.61	-42.61	25.20	-1.96	10.37	-9.09

Sum of squared residuals fp = 1.0002e+01

Values of computed spline:

	x	3.00	6.00	9.00	12.00	15.00	18.00	21.00
y								
17.00		40.74	28.62	19.84	14.29	11.21	9.46	7.09
14.00		48.34	33.97	21.56	14.71	12.32	10.82	7.15
11.00		37.26	24.46	17.21	14.14	13.02	11.23	7.29
8.00		30.25	19.66	16.90	16.28	15.21	12.71	8.99
5.00		36.64	26.75	23.07	21.13	18.97	15.90	11.98
2.00		45.04	33.70	26.25	22.88	21.62	19.39	13.40
