THE FINITE ELEMENT METHOD FOR SOLVING THE POISSON PARTIAL DIFFERENTIAL EQUATION

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INTRODUCTION

Further, we will present the Poisson equations with mixed conditions on frontier solution, using a method of finite element.

1. THE PROBLEM

Let be $D \subset \mathbf{R}^2$ a bordered domain, having the regular frontier \mathbf{S} and an operator A of Laplace type with next expression:

$$A = -\nabla K \nabla, \qquad (1)$$

where $K = (k_{ij} (x, y))$, I, j=1, 2 is a symmetrical matrix of continuous functions on **D**, express of physical anisotropy of the domain **D** on which we study the described phenomena by Poisson equations (3).

We will noted with D_A the whole set of the functions definite on \overline{D} that satisfy the following conditions:

$$u, \frac{\partial u}{\partial x}, \frac{\partial u}{\partial y} \quad continuous \ functions$$
$$\frac{\partial^2 u}{\partial x^2}, \frac{\partial^2 u}{\partial x^{\partial y}}, \frac{\partial^2 u}{\partial y^2} \quad continuous \ functions \ on \ portions$$

They satisfy a mixed condition on the frontier S with this expression:

$$K\nabla u * n + p(s)u - h(s) = 0$$
(2)

where:

n is the normal unit vector toward the exterior frontier S;

n p(s), h(s) are continuous functions on S. Further, we will determine a function $u_0 \in D_A$, solution for the equation:

Au = g, (3) where the function g named in engineering problems as "source function" is at least integrable on **D**.

It is demonstrated that the operator A definite through (3) is symmetrical and positive definite and the solution of the equation (3) with the frontier condition (2) is equivalence on D_A with the functional minimisation

$$F(u) = \langle -Au, u \rangle -2 \langle g, u \rangle =$$

$$\iint_{D} [(\nabla u)K(\nabla u) - 2gu] dx dy$$

$$- \int_{S} (p(s)u^{2} - 2h(s)u) ds$$
(4)

For the function $u_o \in D_A$ determination that accomplish minimum of (4) further we will use the known Riesz method.

2. THE DISCRETE PROBLEM

To the formulate problem is associated a discrete problem. We will build a minifying sequence of functions w_1, w_2, \ldots with imposed form, minimise sequence for the functional (4), that is:

$$\lim_{M\to\infty}\left\|W_M\right\|_E=\left\|u_0\right\|_E,$$

where $\|\cdot\|_{F}$ is the energetic norm.

The idea of finite element method consists in the domain decomposing into disjoint subsets reunion T_i (named elements) satisfying the conditions:

$$\begin{array}{l}
\stackrel{\scriptscriptstyle M}{U}\overline{T}_{i} = D; \\
\overline{T}_{i} \cap \overline{T}_{j} = \begin{cases} a \text{ po int;} \\
a \text{ curve;} \\
\Phi \end{cases} \qquad i,j = \overline{1,M} \quad (5)
\end{array}$$

Corresponding to this divisions the functional F(u) will be:

$$F(u) = \sum_{i=1}^{M} F_{i}(u),$$
 (6)

where: $F_i(u)$ express the energy on the element T_i . To every element T_m of the division $T_1, T_2,...,T_M$ will be associated a number of characteristically points P_m named nodal points, obtaining N nodes on the whole domain. We will associate to every interior node "i" of (x_i, y_i) co-ordinates, a function $f_i(x,y)$ with the following proprieties:

a. f_i continuous on D, i=1,2,...,N,
b. f_i(x_i, y_i) =
$$\delta_{ij}$$
, (7)

c. $\langle \mathbf{f}_{i}, \mathbf{f}_{j} \rangle = 0$, for $i \neq j$,

d. $\forall i \in 1, M$, the function f_i satisfies the condition (2) on **S**.

Let be E_N the subspace of D_A , generated by the function $f_1, f_2, ..., f_N$.

A certain function $w_N \in E_N$ will have the next form:

$$W_{N} = \sum_{i=1}^{N} C_i f_i$$
(8)

The function $u(x, y) \in D_A$ will be approximated on every T_i element with a continuous on D function w(x,y) with continuous on portions partial derivatives of first order.

For a certain element T_m of the proposed division, the approximation function can be writhe as:

$$w_m(x, y) = \sum_{i=1}^{P_m} f_{m_i}(x, y) w_{m_i},$$
(9)

Where f_m is the function f₁, f₂, ... f_N

restriction to T_m element which contains the node "i" and w_{m_i} is the w(x, y) function value in the node "i" and have part of variations parameter . We will insert the matric notations:

$$F_{m} = \begin{bmatrix} f_{m_{1}}, f_{m_{2}}, \dots f_{p_{m}} \end{bmatrix}$$
(10)

$$W_m^T = [w_{m_1}, w_{m_2}, ..., w_{p_m}],$$
 (11)

the relation (9) became:

$$w_m(x, y) = f_m w_m \tag{12}$$

At every element T_m it is associated an matrix L_m with P_m lines and N columns, which put into biunique correspondence the T_m element nodes and obtained nodes on D through performing the divisions in the elements $T_1, ..., T_M$.

Making the notations:

$$W^{T} = [W_{1}, W_{2}, ..., W_{N}]$$
 (13)

and using the matrix L_m we will obtain the link between the relations (12) and (13):

$$W_m = L_m W, \tag{14}$$

and (7) became:

$$\mathbf{w}_{\mathrm{m}}(\mathbf{x},\mathbf{y}) = W^{T} L_{m}^{T} F_{m}^{T}$$
(15)

By means of formula (15) we obtained a interpolation on T_m for the function from D_A . If in (6) the function u is replaced with w_m it is obtained:

$$F(W) = \sum_{m=1}^{M} \left(\iint_{T_m} \left(\nabla W_m \, K \nabla W_m - 2W_m \, g_m \right) dx dy + \int_{S_m} \left(2h_m W_m - p_m W_m^2 \right) ds \right), \tag{16}$$

where:

 g_m , h_m , p_m are the functions g, h, p restrictions at the elements T_m on the frontier S_m .

Taking into consideration (15) the expression of $\nabla W_m(x, y)$ is:

$$\nabla W_m(x, y) = B_m L_m W_m$$
, where:

$$B_{m} = \begin{pmatrix} \frac{\partial f_{m1}}{\partial x} \dots \frac{\partial f_{mP_{m}}}{\partial x} \\ \frac{\partial f_{m1}}{\partial y} \dots \frac{\partial f_{mP_{m}}}{\partial y} \end{pmatrix}$$

The similar expressions with (12) can be obtain also for the functions g, h, p:

 $g(x,y) = F_m g_m,$ $h(x,y) = F_m h_m,$ $p(x,y) = F_m p_m,$ where :

 $g_{m,}h_{m,}p_{m}$ are vectors that contain functions values g, h, p in the element T_{m} nodes.

The expression of the functional (16) became:

$$F(W) = W^{T} \left[\sum_{m=1}^{M} \left(L_{m}^{T} G_{m} L_{m} \right) W - 2 \sum_{m=1}^{M} L_{m}^{T} H_{gm} + 2 \sum_{m=1}^{M} L_{m}^{T} H_{hm} - \sum_{m=1}^{M} L_{m}^{T} H_{p_{m}} W \right]$$
(17)

where we made the notations:

$$H_{g_m} = \left(\iint_{T_m} F_m^T F_m dx dy \right) g_m,$$

$$H_{h_m} = \left(\iint_{S_m} F_m^T F_m ds \right) h_m,$$

$$H_{p_m} = \left(\iint_{S_m} F_m^T F_m ds \right) p_m,$$

The function (17) minimisation relative with W_1 , W_2 ,..., W_N leads at the linear system solution:

(19)

$$\begin{cases} \frac{\partial F}{\partial w_1} = 0\\ \dots \\ \frac{\partial F}{\partial w_n} = 0 \end{cases}$$
(18)

which can be also write as:

 $GW=H_g - H_h + H_p$, where we noted with:

$$egin{aligned} G &= \sum_{m=1}^{M} L_m^T G_m L_m, \ H_g &= \sum_{m=1}^{M} L_m^T H_{g_m}, \ H_h &= \sum_{m=1}^{M} L_m^T H_{h_m}, \ H_p &= \sum_{m=1}^{M} L_m^T H_{p_m}. \end{aligned}$$

The system (19) unknowns are: functions values u_0 that minimises the energy functional and so, the problem (3) solution in net nodes formed through the domain D division into finite elements.

This system solutions is quite comfortable because in case of the definite and symmetrical positive operators, the system (19) matrix is symmetrical having moreover an band structure that means an substantial advantage in numerical solution of system (19) by computer way.

3. THE COMPUTER PROGRAM

The computer program in C++ language solves an undetermined compatible system of n linear equations with n unknowns. For round-off errors reduction which are made by computer, before system and subsystems processing it is replace the equation that must to contain the one equation pivot from subsystem so that the obtained pivot to have the maximum absolute value.

The terms of problem are the system dimension, the coefficients and the free terms. #include <stdio.h> #include <conio.h> #include <conio.h> #include <math.h> int n, i, j, k, p, t; double a [10] [11], x[10], max, s; void main (void) {cout << "Dimensiunea sistemului n="; cin >> n; cout << "Introduceți coeficienții : " <<endl ; for (i=1 ; i<=n ; i++) $\{ \text{cout } << "b ("<< i <<") = ";$ $cin >> a[i] [n+1]; \}$ k=1; t=0; do { max = fabs (a [k] [k]); p = k; for (i = k+1; i < = n; i + +)if $(\max < \text{fabs}(a [i] [k]))$ $\{ \max = \text{fabs } (a [i] [k]); \}$ $p = i; \}$ if (max = = 0)t = 1;else { if (p|=k)for (j = k; j < n+1; j + +) $\{ s = a [i] [k]; \}$ a[i][k] = a[p][j]; $a[p][j] = s; \}$ for $(i = k+1; i \le n; i + +)$ for $(j = k+1; j \le n+1; j++)$ a[i][j] - = a[i][k] * a[k][j] / a[k][k]; $k + +; \}$ while ((k < n) & & (t = = 0)) cout << "Sistemul nu este compatibil determinat ! " << endl ; else { for (i = n; i > 0; i - -) $\{ s = 0;$ for $(j = i+1; j \le n; j + +)$ s + = a [i] [j] * x[j]; $x[i] = (a [i] [n+1] - s / a[i] [i]; \}$ cout << "Soluția sistemului:" << endl; for $(i = 1; i \le n; i + +)$ cout << x(" << i << ") = " << x[i] << endl; $getch(); \}$

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