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## THE DIRECT SOLUTION OF THE DISCRETE POISSON EQUATION ON IRREGULAR REGIONS

BY

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STAN-CS-71-195 DECEMBER, 1970

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NATIONAL TECHNICAL NFORMATION SERVICE Springfield, Va. 22151

Unclassified						
Security Classification						
DOCUMENT CONTROL DATA - R & D						
Security classification of title, body of abstract and indexing i	annotation must be e					
Computer Science Department			SSIFIED			
•		25. GROUP				
Stanford, Calif. 94305		20. 04000				
3 REPORT TITLE						
THE DIRECT SOLUTION OF THE DISCRETE POISSO	ON EQUATION C	N IRREGUL	AR REGIONS			
A DESCRIPTIVE NOTES (Type of report and inclusive dates) Technical Report, December 1970						
5. AUTHOR(S) (First name, middle initial, last name)						
B.L. Buzbee, F.W. Dorr, J.A. George, G.H.	Golub	2				
6. REPORT DATE	78. TOTAL NO. OF	PAGES	76. NO. OF REFS			
December 1970	30		24			
B. CONTRACT OR GRANT NO.	Se. ORIGINATOR'S	REPORT NUM	6ER(\$)			
D. PROJECT NO. NR 044-211	NO0014-67-A-00112-0029 B. PROJECT NO. NR 044-211 Technical Report No. STAN-CS-71-195					
c. also A.E.C. AT (04-3) 326, PA 30	(04-3) 326, PA 30 9b. OTHER REPORT NO(5) (Any other numbers that may be assigned this report)					
d.	Los Alamo	s Laborato	ory Report LA 4553			
10. DISTRIBUTION STATEMENT						
Releasable without limitations on dissemin						
11. SUPPLEMENTARY NOTES	SUPPLEMENTARY NOTES 12. SPONSORING MILITARY ACTIVITY Mathematics Program					
*	Office of Naval Research					
	Arlington, Virginia 22217					
13. ABSTRACT		<u>, , , , , , , , , , , , , , , , , , , </u>				
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KEY WORDS	ROLE	WT	ROLE	WT	ROLE	
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Poisson equation						
Discrete			·			
Partial differential equation						
Elliptic difference equation						
Computer solution					-	
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## THE DIRECT SOLUTION OF THE DISCRETE

POISSON EQUATION ON IRREGULAR REGIONS

by

B. L. Buzbee, F. W. Dorr, J. A. George, G. H. Golub

To be issued simultaneously as Los Alamos Scientific

Laboratory Report LA 4553.

Computer Science Department, Stanford University, Stanford, California 94305. Research supported by the Office of Naval Research under grant No. NOOL-67-A-00112-0029, and by the Atomic Energy Commission under grant No. AT(04-3)326, PA30.

Los Alamos Scientific Laboratory of the University of California, Los Alamos, New Mexico 87544. Research supported by the U. S. Atomic Energy Commission under Contract No. W-7405-ENG-36.

## Abstract

There are several very fast direct methods which can be used to solve the discrete Poisson equation on rectangular domains. We show that these methods can also be used to treat problems on irregular regions. 1. <u>Introduction</u>. Within the past few years, several very fast and accurate direct methods have been developed for solving finite difference approximations to the Poisson equation,

$$\begin{array}{c} \Delta u = f & \text{in } \mathbb{R}, \\ u = g & \text{on } \partial \mathbb{R}. \end{array}$$

These methods can usually be applied only on rectangular regions, although the differential operator and boundary conditions can be more general than those in the Poisson equation. In this paper, we will show how these algorithms for rectangular domains can also be used effectively on irregular regions. The approach used is similar to that employed by Hockney [16, 17], Buneman [7], and George [14]. We also mention the work of Angel [1-4], Angel and Kalaba [5], Collins and Angel [9], Kalaba [20], and Roache [22] on the use of direct methods for problems in irregular regions.

We will not discuss the details of any specific direct method. A survey of these procedures is given in [11], and in particular we cite the recent work of Buneman [6], Buzbee, Golub, and Nielson [8], and Hockney [16].

We will also not consider the derivation of the finite difference equations that approximate the partial differential equation. This subject is treated in detail by Forsythe and Wasow [13], and we assume that the problem has been reduced to finding the solution of a matrix equation Ax = y. The matrix A is frequently very large and sparse, but its structure does not permit the application of the most efficient direct methods. For our computational procedure, we alter certain rows of A to obtain a matrix B, and we will show how to define a modified right-hand side z so that the solution x also satisfies the equation Ex = z. The matrix B is chosen so that these equations can be solved by the direct methods.

This method is computationally advantageous when we are solving a sequence of equations  $Ax_i = y_i$ . This situation frequently arises in time-dependent partial differential equations, in nonlinear problems, and in linear problems where the right-hand side is varied but the region and differential operator remain the same. After some initial computation, each solution  $x_i$  can be obtained in approximately twice the time required for the solution of an equation Bx = z.

In Sections 2 and 3 we derive this algorithm in a general form. We describe a number of applications of the method in Sections 4 and 5, and in Section 6 we present some computational results.

2. <u>Method of Solution if  $det B \neq 0$ </u>. Suppose that we are given an n by n matrix A and an integer p with  $1 \le p \le n$ . We wish to modify p rows of A to obtain another matrix B. Without loss of generality we assume that the first p rows of A are to be changed, since we can achieve this situation by multiplying A by a suitable permutation matrix. However, we emphasize that this multiplication should not be done explicitly in the computational procedure. Rather, the rearrangement of rows should be done implicitly by indexing. The direct methods mentioned later in the paper require that B has a particular structure, which could be altered by the permutation transformation.

Partition A in the form

$$\mathbf{A} = \begin{pmatrix} \mathbf{A}_1 \\ \mathbf{A}_2 \end{pmatrix} ,$$

where  $A_1$  is a p by n matrix and  $A_2$  is an (n-p) by n matrix. We then write

$$B = \begin{pmatrix} B_1 \\ A_2 \end{pmatrix} ,$$

where  $B_1$  is a p by n matrix. For the remainder of this section, we assume that det  $B \neq 0$ .

Suppose we are given a linear equation A = y. We partition y in the same way as A, and write

$$\underline{y} = \begin{pmatrix} \underline{y}_1 \\ \underline{y}_2 \end{pmatrix}$$
.

Let  $\overline{y}$  be any vector of the form

$$\overline{\mathbf{y}} = \begin{pmatrix} \overline{\mathbf{y}}_1 \\ \mathbf{y}_2 \end{pmatrix}$$

If W is an arbitrary nonsingular p by p matrix, we define an n by p matrix  $\overline{W}$  by

$$\overline{\mathbf{W}} = \left(\begin{array}{c} \mathbf{W} \\ \mathbf{O} \end{array}\right) \quad .$$

Define the p by p matrix C by

$$C = A_1 B^{-1} \overline{W}$$

Following Hockney [16], we call C the <u>capacitance matrix</u>.<sup>\*/</sup> Assume that there exists a p by 1 vector  $\beta$  that is a solution to the equation

$$C \stackrel{\beta}{\sim} = \underbrace{y_1}_{1} - A_1 B^{-1} \underbrace{\overline{y}}_{1} . \qquad (1)$$

Since A and B differ only in the first p rows, it is easy to verify that a solution x to the equation Ax = y is given by

$$\underset{\sim}{\mathbf{x}} = \mathbf{B}^{-1} \left( \underbrace{\overline{\mathbf{y}}}_{\mathbf{x}} + \overline{\mathbf{w}} \underset{\sim}{\mathbf{\beta}} \right)$$

We first show that this method of obtaining the solution x will be valid whenever the original system Ax = y is consistent.

Theorem. If  $jet B \neq 0$ , then

$$\det C = \frac{(\det A)(\det W)}{\det B}$$

\*/ Hockney actually refers to C<sup>-1</sup> as the capacitance matrix. Since C may be singular in our development, we have adopted the present notation.

If the system Ax = y is consistent, then Eq. (1) is also consistent. <u>Proof.</u> Partition B<sup>-1</sup> in the form

$$B^{-1} = \begin{pmatrix} D_1 & D_2 \end{pmatrix},$$

where  $D_1$  is n by p and  $D_2$  is n by (n-p). It then follows that

$$BB^{-1} = \begin{pmatrix} B_1 D_1 & B_1 D_2 \\ & & \\ A_2 D_1 & A_2 D_2 \end{pmatrix} = \begin{pmatrix} I & 0 \\ & & \\ 0 & I \end{pmatrix},$$

and

$$AB^{-1} = \begin{pmatrix} A_1 D_1 & A_1 D_2 \\ A_2 D_1 & A_2 D_2 \end{pmatrix} = \begin{pmatrix} A_1 D_1 & A_1 D_2 \\ 0 & I \end{pmatrix}.$$

Thus we have

$$C = A_1 B^{-1} \overline{W} = A_1 D_1 W ,$$

and so

$$\det C = \det (A_1 D_1) \det W$$
$$= \det (A B^{-1}) \det W$$
$$= \frac{(\det A)(\det W)}{\det B}$$

To prove the consistency statement, suppose  $\mathbf{C}^{T} \stackrel{\mathbf{\gamma}}{\sim} = \stackrel{\mathbf{0}}{\sim}$ . Write

$$\mathbf{A}\mathbf{B}^{-1} = \begin{pmatrix} \mathbf{C}\mathbf{W}^{-1} & \mathbf{E}_{\mathbf{1}} \\ & \mathbf{I} \\ \mathbf{O} & \mathbf{I} \end{pmatrix}$$

,

and define an n by 1 vector  $\overline{r}$  by

$$\overline{\mathfrak{L}} = \begin{pmatrix} \mathfrak{L} \\ - \mathbf{E}_{1}^{\mathrm{T}} \mathfrak{L} \end{pmatrix}$$

We then have

$$\left(\mathbf{A} \mathbf{B}^{-1}\right)^{\mathrm{T}} \mathbf{\overline{\chi}} = \begin{pmatrix} \left(\mathbf{W}^{-1}\right)^{\mathrm{T}} \mathbf{c}^{\mathrm{T}} & \mathbf{o} \\ & & \\ \mathbf{E}_{1}^{\mathrm{T}} & \mathbf{I} \end{pmatrix} \begin{pmatrix} \mathbf{\chi} \\ & \\ \mathbf{E}_{1}^{\mathrm{T}} & \mathbf{I} \end{pmatrix} = \mathbf{0}$$

Since the system  $A_x = y$  is assumed to be consistent, we therefore have  $\overline{\chi}^T y = 0$ , which is the same as  $\chi^T (y_1 - E_1 y_2) = 0$ . But then

$$\chi^{T}(\underline{y}_{1} - \underline{A}_{1} \underline{B}^{-1} \overline{\underline{y}}) = \chi^{T}(\underline{y}_{1} - \underline{C} \underline{W}^{-1} \overline{\underline{y}}_{1} - \underline{E}_{1} \underline{y}_{2})$$
$$= -(\underline{C}^{T} \underline{\chi})^{T} \underline{W}^{-1} \overline{\underline{y}}_{1}$$
$$= 0 ,$$

which is the consistency condition for Eq. (1).

The Woodbury formula [18, pp. 123 - 124] for the inverse of a matrix (B + FG) is

$$(B + FG)^{-1} = B^{-1} (I - F(I + GB^{-1}F)^{-1}GB^{-1})$$

This equation has been used in direct methods for solving the Poisson equation by George [14], and for the biharmonic equation by Golub [15]. If A is nonsingular we write

$$\mathbf{A} = \mathbf{B} + \mathbf{F}\mathbf{G} ,$$

where  $F = \overline{W}$ , and G is the p by n matrix given by

$$G = W^{-1} (A_1 - B_1)$$

For the case in which A is nonsingular, the algorithm we have derived is equivalent to using the Woodbury formula for  $A^{-1}$ .

Suppose that we have a very efficient method for solving equations of the form Bz = w. The solution of the equation Ax = y then proceeds in the following steps:

(1) Compute 
$$C = A_1 B^{-1} \overline{W}$$

(2) Compute 
$$\bar{\mathbf{x}} = \mathbf{B}^{-1} \bar{\mathbf{y}}$$
,

(3) Solve the equation  $C\beta = y_1 - A_1 \overline{x}$ .

The solution x can then be obtained from the formula

$$\sum_{n=1}^{\infty} = B^{-1} \left( \overline{y} + \overline{W} \frac{\beta}{n} \right) \quad . \tag{2}$$

If it is possible to store the vector  $\overline{x}$  and the matrix  $\overline{B} = B^{-1}\overline{W}$ , then x = a can also be computed from

$$\mathbf{x} = \mathbf{\bar{x}} + \mathbf{\bar{B}}\mathbf{\beta} \quad . \tag{3}$$

The decision whether to use Eq. (2) or Eq. (3) would be made on consideration of storage requirements, and on the relative speed of solving the system in Eq. (2) versus multiplying by the matrix in Eq. (3). For problems arising from elliptic difference equations, it is frequently better to use Eq. (2) because B has a band structure, but the matrix  $\overline{B}$  may be full.

The type of application we have in mind for this method is one in which we have to solve a number of equations  $A_{X_i} = y_i$ . In this case, we compute the capacitance matrix and factor it as part of a preprocessing stage. The solution of each equation  $A_{X_i} = y_i$  is then approximately as fast as the time it takes to solve two equations Bz = w.

To be specific, let  $\theta(n)$  denote the number of arithmetic operations necessary to solve a system Bz = w. Then to compute C and form its LU decomposition in a preprocessing stage requires approximately

$$p \theta(n) + k_1 p^2 n + k_2 p^3$$

operations (cf. [19, Sec. 2.1]). In many cases the matrix  $A_1$  is sparse, and this estimate is

$$p \theta(n) + k_3 p^3$$
 (4)

operations. To compute the solution to a particular equation  $A_{X} = y$  using Eq. (2) takes an additional

$$2\theta(n) + k_{\mu}pn + k_{5}p^{2}$$

operations. If  $A_1$  is sparse and we let W = I, this estimate can be replaced by

$$2\theta(n) + k_{6}p^{2}$$
<sup>(5)</sup>

operations. To compute a particular solution using Eq. (3) requires

$$\theta(n) + k_7 p n + k_8 p^2$$
 (6)

operations. In general this estimate cannot be reduced, because the matrix  $\overline{B}$  may be full.

3. <u>Method of Solution if rank(B) = n-l</u>. The method derived in Section 2 gives a procedure for finding a p by 1 vector  $\delta$  such that a solution x to Ax = y also satisfies the equation

$$Bx = y + \begin{pmatrix} \delta \\ \tilde{} \\ 0 \\ \tilde{} \end{pmatrix}$$

If B is singular, it may not be possible to find such a vector  $\delta$ . To show this, suppose  $B^T v = 0$  but  $v \neq 0$ . In order for  $\delta$  to exist, we must satisfy the consistency condition

$$\underbrace{\mathbf{v}}_{\mathbf{u}}^{\mathrm{T}}(\mathbf{y} + \sum_{i=1}^{p} \delta_{i} \underbrace{\mathbf{e}}_{i}) = 0 \quad .$$
(7)

If  $v^{T}e_{i} = 0$  for  $1 \le i \le p$  and  $v^{T}y \ne 0$ , it is not possible to satisfy Eq. (7). However, if A is nonsingular this difficulty does not arise, because then the only vector v satisfying  $B^{T}v = 0$  and  $v^{T}e_{i} = 0$  for  $1 \le i \le p$  is v = 0.

We will now describe an algorithm we have used when rank(B) = n-1 and A is nonsingular. There are two advantages in treating this particular case. First, the construction is quite simple, and it is easy to see how the method could be extended to a more general matrix B. Second, the case rank(B) = n-1has a special significance in the solution of partial differential equations, because this condition is satisfied by the matrix corresponding to the Neumann problem. For simplicity, we assume that the matrix W of Section 2 is the identity matrix.

<u>Theorem 2</u>. Assume that A is nonsingular and rank(B) = n-1, and let u and v be two non-zero vectors satisfying Bu =  $B^{T}v = 0$ . Then there exists an integer k with  $1 \le k \le p$  such that  $v^{T}e_{k} \ne 0$ . Define a constant

$$\alpha = (\mathbf{v}^{\mathrm{T}} \mathbf{e}_{k})^{-1}$$
,

and let  $\bar{x}$  be a solution to

$$B\bar{x} = \bar{y} - (\alpha v^T \bar{y}) e_k$$
.

For  $1 \leq i \leq p$  and  $i \neq k$  let  $\underline{\eta}_i$  be a solution to

$$B_{-i} = e_i - (\alpha v^T e_i) e_k$$
,

and let  $\underline{\eta}_k = \underline{u}$ . Let C be the p by p matrix whose i-th column is the vector  $A_1 \underline{\eta}_i$ . Then C is nonsingular, and, if  $\beta$  is the solution to

$$C_{\tilde{\mu}}^{\beta} = y_{1} - A_{1} \bar{x} ,$$

<u>the solution</u>  $\underset{\sim}{x}$  <u>to</u>  $Ax = \underset{\sim}{y}$  <u>is given</u> by

$$\sum_{n=1}^{\infty} = \sum_{n=1}^{\infty} + \sum_{i=1}^{p} \beta_{i} \eta_{i}$$

<u>Proof</u>. If we partition v in the same way as y, we have  $A^{T}v = A_{1}^{T}v_{1} + A_{2}^{T}v_{2}$  and  $B^{T}v = B_{1}^{T}v_{1} + A_{2}^{T}v_{2}$ . Thus if  $B^{T}v = 0$  and  $v_{1} = 0$  we would have  $A^{T}v = 0$ . Since A is nonsingular and  $v \neq 0$  this cannot happen, and hence  $v_{1} \neq 0$ .

To prove that C is nonsingular, we show that  $C\beta = 0$  implies  $\beta = 0$ . Suppose  $\beta$  is an arbitrary vector such that  $C\beta = 0$ . Then  $x = \sum_{i=1}^{p} \beta_i \eta_i$  satisfies Ax = 0, and hence x = 0. This implies that Bx = 0, or

$$\sum_{i=1}^{p} \beta_{i} e_{i} = \alpha \left( \sum_{i=1}^{p} \beta_{i} v^{T} e_{i} \right) e_{k}$$

Thus  $\beta_i = 0$  for  $1 \le i \le p$  and  $i \ne k$ , and the condition x = 0 then implies that  $\beta_k = 0$ . Thus  $\beta = 0$ , and so C is nonsingular.

Remark. As we discussed in Section 2, the computation proceeds in the following steps:

- (1) Compute (and factor) C,
- (2) Compute  $\bar{x}$ ,
- (3) Solve for  $\beta$  .

The solution x can then be obtained from the formula

$$\underset{\sim}{x} = \underset{\sim}{\bar{x}} + \underset{i=1}{\overset{p}{\sum}} \beta_{i} \underset{\sim}{\eta}_{i} .$$

However, if the problem arises from a partial differential equation, it is more efficient computationally to obtain x in the form

$$\mathbf{x} = \hat{\mathbf{x}} + \hat{\boldsymbol{\beta}} \mathbf{u} ,$$

where  $\hat{x}$  is a solution to

$$B\hat{x} = \bar{y} - (\alpha v^{T} \bar{y}) e_{k} + \sum_{\substack{i=1\\i \neq k}}^{p} \beta_{i}(e_{i} - (\alpha v^{T} e_{i})e_{k})$$

and

$$\hat{\boldsymbol{\beta}} = \left[ \underbrace{\boldsymbol{u}}^{\mathrm{T}}_{-} \cdot \underbrace{\boldsymbol{x}}_{+} + \sum_{i=1}^{p} \boldsymbol{\beta}_{i} \underbrace{\boldsymbol{u}}^{\mathrm{T}}_{-} \cdot \underbrace{\boldsymbol{\eta}}_{+} - \underbrace{\boldsymbol{u}}^{\mathrm{T}}_{-} \cdot \underbrace{\boldsymbol{x}}_{-} \right] \left[ \underbrace{\boldsymbol{u}}^{\mathrm{T}}_{-} \cdot \underbrace{\boldsymbol{u}}_{-} \right]^{-1} .$$

4. Applications to Partial Differential Equations by Imbedding. Suppose we are given a two-dimensional bounded region R in the x - y plane, and we wish to find a solution u to the Poisson equation,

$$\begin{array}{c} \Delta u = f & \text{in } R \\ u = g & \text{on } \partial R \end{array} \right\}$$

We assume that this differential equation is approximated by a finite difference equation (cf. Forsythe and Wasow [13]). Thus we have a finite set of unknowns  $\{U_i \mid 1 \le i \le n_0\}$  which approximate the solution u at the grid points. If we denote by  $\Delta_h$  a finite difference approximation to the Laplacian operator  $\Delta$ , by  $R_h$  the discrete interior of the grid, and by  $\partial R_h$  the discrete boundary of the grid, then the discrete Poisson equation can be written in the form

$$\begin{array}{c} \Delta_{\mathbf{h}} \mathbf{U} = \mathbf{f} & \operatorname{in} \mathbf{R}_{\mathbf{h}} \\ \mathbf{U} = \mathbf{g} & \operatorname{on} \partial \mathbf{R}_{\mathbf{h}} \end{array} \right\}$$
(8)

Let  $R'_h$  be a discrete rectangular region such that  $R_h \subset R'_h$  and  $\partial R_h \subset R'_h \cup \partial R'_h$ , and let  $S_h = \partial R_h \cap R'_h$ . Extend the functions f and g to the regions  $R'_h$  and  $\partial R_h \cup \partial R'_h$  respectively, and consider the equation

$$\Delta_{\mathbf{h}} \mathbf{U} = \mathbf{f} \quad \text{in } \mathbf{R}_{\mathbf{h}}' - \mathbf{S}_{\mathbf{h}} , \\ \mathbf{U} = \mathbf{g} \quad \text{on } \mathbf{S}_{\mathbf{h}} \cup \partial \mathbf{R}_{\mathbf{h}}' . \end{cases}$$

$$(9)$$

We will solve Eq. (9), and the solution U will then also satisfy Eq. (8).

Equation (9) is a linear equation in the unknowns  $\{U_i \mid 1 \le i \le n\}$ . Observe that we may have increased the number of unknowns by the imbedding process, so that  $n_0 \le n$ . We write Eq. (9) as a matrix equation AU = V, and the matrix A can frequently be chosen to be block tr.diagonal with tridiagonal matrices as the non-zero blocks (cf. [15]).

Let p be the number of grid points in  $S_h$ . We modify the p rows of A and V corresponding to the equations

$$U = g$$
 on  $S_h$ ,

and replace them with the equations

$$\Delta_h U = f$$
 on  $S_h$ .

This defines a new matrix B and a new right-hand side  $\overline{\nabla}$ . An equation BU =  $\overline{\nabla}$  corresponds to the difference equation

$$\begin{array}{c|c} \Delta_{\mathbf{h}} U = \mathbf{f} & \operatorname{in} \mathbf{R}_{\mathbf{h}}' \\ U = g & \operatorname{on} \partial \mathbf{R}_{\mathbf{h}}' \end{array} \right\} .$$
(10)

Since  $R'_h$  is a rectangular region, we have very fast methods for solving Eq. (10). We can now apply the method of Section 2 to solve the equation AU = V by using the modified matrix B.

To illustrate this construction, let R be a rectangular region with an interior rectangle removed, such as that shown in Figure 1. For simplicity, we assume that the discrete boundary  $\partial R_h$  is a subset of  $\partial R$ . The imbedding rectangle is  $R'_h = R_h \cup S_h \cup T_h$ . The only function extension required for this example is that f be defined (arbitrarily) in  $S_h \cup T_h$ . To define this extension, we can set  $f \equiv 0$  in  $S_h \cup T_h$ , or we can define f so that it is continuous in all of  $R'_h$ . The advantage of using a continuous f is that the solution to Eq. (10) is then smooth. However, the direct methods used to solve Eq. (10) are so accurate that the smoothness of the solution does not appear to influence the computational results. Therefore,

1,

in the examples we have considered, we extend f by setting  $f \equiv 0$  in  $S_h \cup T_h$ .

If we let W = I in the method of Section 2, this algorithm is closely connected with the discrete Green's function for the region  $R'_h$  (cf. [13, pp. 314-318]). In fact, the method is then equivalent to adding suitable multiples of the discrete Green's function for the points on  $S_h$  so that the boundary conditions on  $S_h$  will be satisfied. Since we have Dirichlet boundary conditions on  $S_h$ , by a proper ordering of the unknowns we can write

$$\mathbf{A}_{\mathbf{1}} = (\mathbf{I} \quad \mathbf{0}) \quad .$$

Since B is positive definite and

$$\mathbf{C} = (\mathbf{I} \quad \mathbf{0}) \mathbf{B}^{-1} \begin{pmatrix} \mathbf{I} \\ \mathbf{0} \end{pmatrix},$$

we see that C is also positive definite in this case. This is advantageous because Cholesky decomposition can then be used to compute an  $LL^T$  decomposition of C (cf. [12, Chap. 23]).

If the grid on  $R'_h$  has N points on a side, we have  $n = N^2$ . In that case, we can solve the system  $BU = \overline{V}$  in approximately

$$\theta(N) = 5 N^2 \log_2 N$$

operations (cf. [11, p. 260]). The preprocessing then takes

$$5 \text{ p N}^2 \log_2 \text{N} + \text{k}_3 \text{ p}^3$$

operations (cf. Eq. (4)). To solve Eq. (8) for a particular choice of f and g by using Eq. (2) with W = I takes an additional

$$10 N^2 \log_2 N + k_6 p^2$$

operations (cf. Eq. (5)). If we use Eq. (3) to compute the solution, it takes an additional

$$5 N^2 \log_2 N + k_7 p N^2 + k_8 p^2$$

operations (cf. Eq. (6)). Thus if  $p \gg \log_2 N$  it is faster to use Eq. (2) to compute the solution. We also observe that for this problem the matrix  $B^{-1}$  is full [23, p. 85], so to store  $\overline{B}$  in using Eq. (3) would require  $p N^2$  locations. Thus for large values of p and N it is both faster and more economical in terms of storage to use Eq. (2) to compute the solution to a particular equation.

It should be clear that the imbedding procedure can be applied to other elliptic difference operators with other types of boundary conditions. To be a practical procedure, we simply require that we have a fast method for solving the imbedded problem in the rectangular region.

As another example, consider the region shown in Figure 2. This problem arises in the time-dependent study of a rotating fluid [10], and the fluid surface is moving slowly. We are given Dirichlet boundary data on  $S_h$ , and Neumann boundary data on  $\partial R_h - S_h$ . The imbedding rectangle is  $R'_h = R_h \cup S_h \cup T_h$ , and we use Neumann boundary conditions on  $\partial R'_h$ . Thus B corresponds to the Neumann problem on  $R'_h$ , and the rank of B is n-1. The method of Section 3 can then be applied, and direct methods for solving the rectangular Neumann problem are given in [8].

For an example with the Poisson operator in another geometry, consider the region in the z-r plane shown in Figure 3. This problem arises in the

time-dependent study of a plasma [21], and a Poisson equation must be solved at each time step. The boundary conditions are Dirichlet on  $S_h$  and Neumann on  $\partial R_h^{(1)}$ . We use Neumann boundary conditions on  $\partial R_h^{(1)}$  and  $\partial R_h^{\prime(3)}$  and Dirichlet boundary conditions on  $\partial R_h^{\prime(2)}$  and  $\partial R_h^{\prime(4)}$  for the imbedding region  $R'_h = R_h \cup S_h \cup T_h$ . The elliptic difference equation in  $R'_h$  is solved by the method of matrix decomposition [8]. 5. <u>Applications to Partial Differential Equations by Splitting</u>. There are many problems for which the imbedding approach is not an economical algorithm. For example, imbedding the region in a rectangle may introduce an excessively large number of additional unknowns that are not necessary to the solution of the original problem. Another instance is one in which the differential operator or the mesh size changes in different parts of the region. In this section, we give two such examples. In each case, the method of Section 2 can be used to split the problem into two rectangular problems, which can be solved by the usual direct methods.

Consider the elongated L-shaped region in Figure 4, and the equation

 $\Delta_h U = f \quad \text{in } R_h ,$   $U = g \quad \text{on } \partial R_h .$ 

We assume that points on the line marked  $T_h$  are all grid points. To define the matrix B, we replace the equations

$$\Delta_h U = f$$
 on  $T_h$ 

by the equations

$$U = g$$
 on  $T_h$ ,

where g has been (arbitrarily) extended to  $T_h$ . The solution of an equation  $B \bigcup = \overline{V}$  now consists of solving the two rectangular problems

$$\Delta_{h} U_{i} = f \quad in \ R_{h}^{(i)},$$
$$U_{i} = g \quad on \ \partial \ R_{h}^{(i)},$$

for i = 1, 2. We can then apply the method of Section 2 to solve the original

problem. This algorithm is similar to one developed in [8, Sec. 9] for non-rectangular regions.

As another example, consider the multiple-material problem shown in Figure 5. The differential equation is

$$\frac{\partial}{\partial x} \left( \sigma(x) \frac{\partial u}{\partial x} \right) + \frac{\partial}{\partial y} \left( \tau(y) \frac{\partial u}{\partial y} \right) = f(x, y)$$

and

The functions  $\sigma_1(x)$ ,  $\sigma_2(x)$ , and  $\tau(y)$  are assumed to be smooth. Dirichlet data is given on  $\partial R$ , and we require that  $\sigma \frac{\partial u}{\partial x}$  be continuous across the boundary between  $R^{(1)}$  and  $R^{(2)}$ . The computational procedure is essentially the same as that for the L-shaped region. The only difference is that in forming the matrix B we replace the equations for the continuity of  $\sigma \frac{\partial u}{\partial x}$ across the line  $T_h$  by the equations

$$U = g$$
 on  $T_h$ .

As before, the equation  $B U = \overline{V}$  corresponds to the two rectangular problems

$$\frac{\partial}{\partial x} \left( \sigma_{i}(x) \frac{\partial u_{i}}{\partial x} \right) + \frac{\partial}{\partial y} \left( \tau(y) \frac{\partial u_{i}}{\partial y} \right) = f(x, y) \quad \text{in } \mathbb{R}^{(i)},$$
$$u_{i}(x, y) = g(x, y) \quad \text{on } \partial \mathbb{R}^{(i)},$$

for i = 1, 2. These problems can be solved directly by the method of matrix decomposition [8, Sec. 8]. A similar method can be used for the case in which  $\tau(y)$  is only piecewise smooth.

It is clear that this splitting method can be applied to the Poisson equation in regions such as that in Figure 5 when different mesh sizes are used in  $R^{(1)}$  and  $R^{(2)}$ . The method developed in [8, Sec. 8] can also be adapted to include rectangular problems with irregular meshes.

6. <u>Computational Results</u>. In Table 1 we have tabulated some computational results for two regions of the form of Figure 1. In each case, a square with sides of length 1 has a symmetrically located square removed from its center. For region 1 the inner square has sides of length  $\frac{1}{8}$ , and for region 2 the inner sides are of length  $\frac{1}{4}$ . We solve the Poisson equation with Dirichlet boundary conditions for the function  $u(x, y) = x^2 + y^2$ . This function was selected because there is no truncation error, and all of the measured error is due to inaccuracies in the solution of the difference equations. All of the computations were performed on a CDC 6600 computer.

The iterative methods used are:

- SOR : point successive overrelaxation [23, p. 58],
- SLOR: successive line overrelaxation [23, p. 80],
- ADI : Peaceman-Rachford alternating direction implicit iteration [24, Chap. 6].

The iteration parameters used are those for the imbedding rectangle  $R'_h$ , and for ADI the parameters for cycles of length four are calculated by the Wachspress algorithm [24, Chap. 6]. The initial guess is identically zero, and the iterations are terminated when the maximum difference between iterates is less than 10<sup>-5</sup>.

The direct method used is variant one of the Buneman algorithm [8, Sec. 11]. Preprocessing times are given in Table 2. Computational results for a similar problem are given in [14].

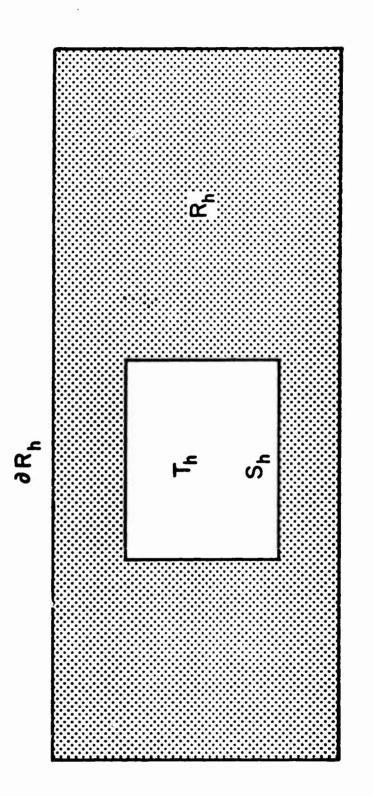
The problem described in Section 4 for the region in Figure 2 has been treated by Daly and Nichols [10]. The mesh used has  $23 \times 40 = 920$  points. Using the direct method of matrix decomposition, a particular solution requires about 30 - 50% of the time required for a point Gauss-Seidel iterative procedure.

The problem discussed in Section 4 for the region in Figure 3 has been treated by Morse and Rudsinski [21]. The mesh used has  $52 \times 98 = 5096$ points, and the preprocessing time is approximately 150 seconds. The region and differential operator are very seldom changed, so the factored capacitance matrix is stored on magnetic tape. Thus there is essentially no preprocessing time for the execution of the program. To solve for a particular solution requires about 2 seconds, which is approximately 40% of the time required for a successive line overrelaxation iterative procedure. Table 1. Computational results for solving the discrete Poisson equation.

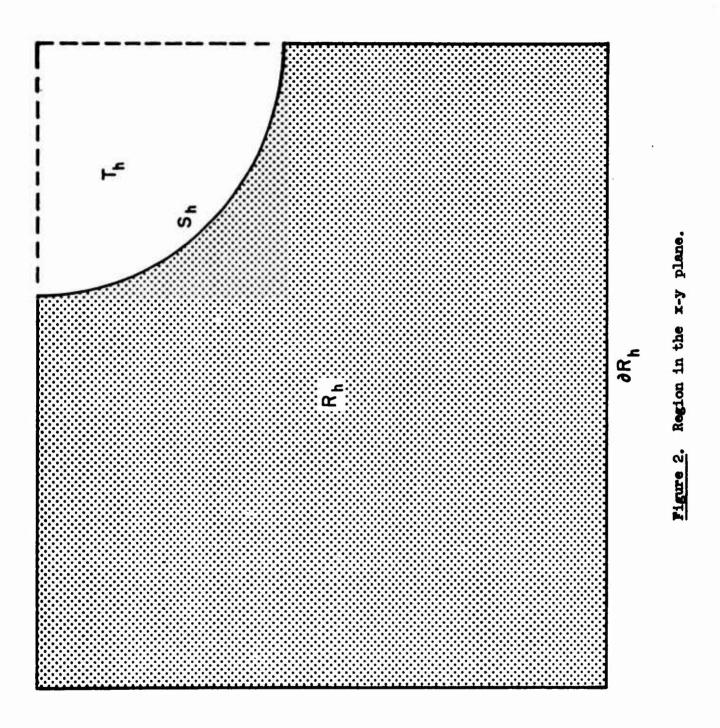
Region	h	р	Method.	Maximum Error	Computation Time (Sec.)	Scaled Computation Time
			SOR	5.02 (- 6)	3.586	21,866
	$\frac{1}{32}$	16	SLOR	7.63 (-6)	2.654	16.183
			ADI	2.36 (-6)	1.128	6.878
1			Direct	4.44 (-13)	0.164	1.000
			SOR	8.12 (- 6)	29.388	43.994
	1 64	32	SLOR	7.95 (-6)	21.424	32.072
			ADI	3.41 (-6)	5.642	8.446
			Direct	1.90 (-12)	0.668	1.000
			SOR	2.35 (-6)	3.570	21.250
	1 32	32	SLOR	6.48 (-6)	2.558	15.226
			ADI	2.11 (-6)	0.870	5.179
2			Direct	3.77 (-13)	0.168	1.000
			SOR	2.02 (-6)	29.624	43.565
	1 64	64	SLOR	9.96 (-6)	20.510	30.162
			ADI	3.57 (-6)	5.332	7.841
			Direct	1.54 (-12)	0.680	1.000

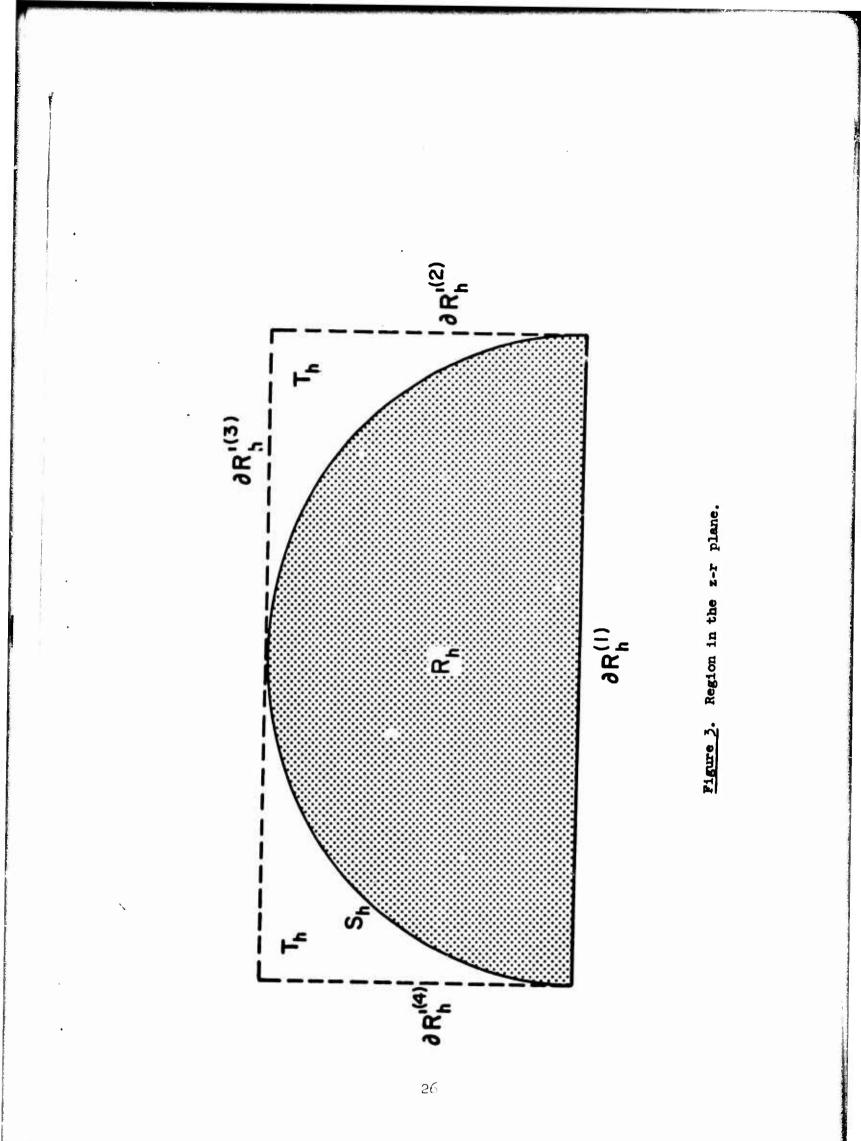
Table 2. Preprocessing time for the direct method results in Table 1.

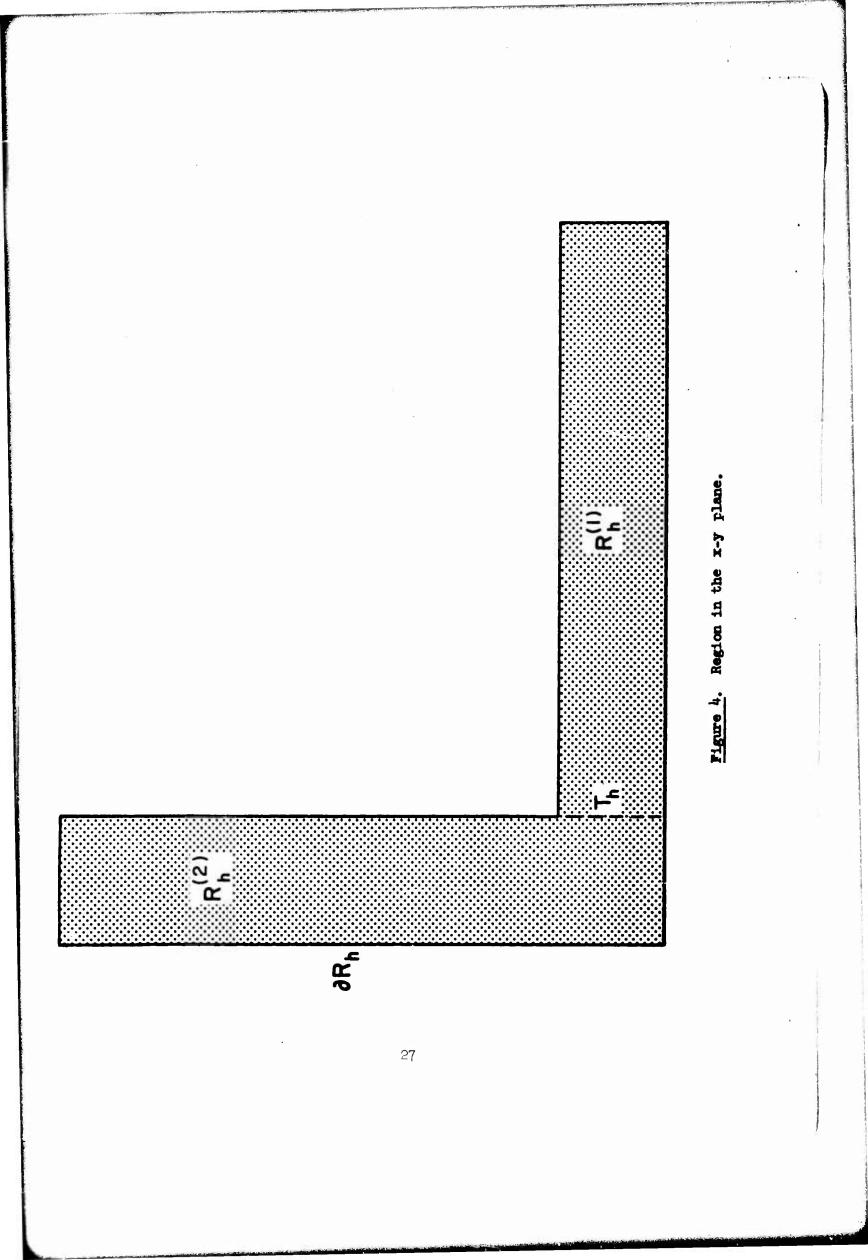
Region	h	Preprocessing Time (Sec.)
1	$\frac{1}{32}$	1.062
	1 64	8.670
2	$\frac{1}{32}$	2.188
	1 64	17.698

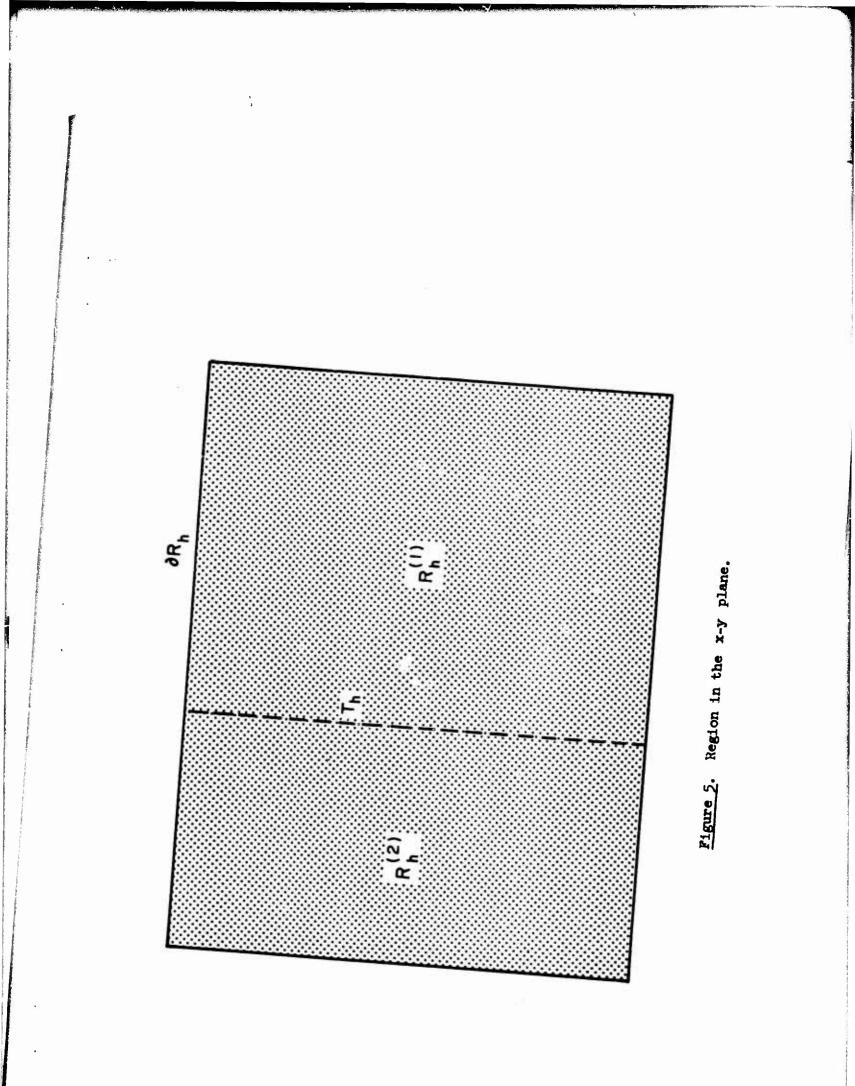












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