

# Boundary Conditions and Calculation of Surface Values for the General Two-Dimensional Electromagnetic Induction Problem

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## *Summary*

General boundary conditions for the problem of electromagnetic induction in a two-dimensional model of a conductor with an arbitrary sub-surface conductivity structure are considered. Program subroutines for both *E*-polarization and *H*-polarization cases are given. These boundary condition subroutines can be used to replace the previously presented subroutines and allow the solution of any conductivity configuration within the conducting region by use of the same numerical technique. An example of a particular model with a sub-surface step structure is illustrated. Also, an improved method of calculating the surface values of the tangential component of the *H*-field (*E*-case) and the tangential component of the *E*-field (*H*-case) at the surface of the conducting region is given for the numerical solution. This new method uses a derivative approximated from the true functional form of the fields instead of a linear approximation and may be applied when a layered or subsurface anomaly is modelled. Some general discussion of the numerical method is given.

## 1. Introduction

At present there is considerable interest in the solution of the problem of electromagnetic induction in the Earth and the local perturbations of the fields when a lateral inhomogeneity is encountered. Jones & Price (1970) considered a two-dimensional problem with a conducting half-space made up of two quarter-spaces of different conductivity, and Jones & Price (1971) considered a surface or buried region of rectangular cross-section of one conductivity surrounded by a region of different conductivity. Jones & Pascoe (1971) extended this work to consider a region of arbitrary shape and of several conductivities surrounded by a region of different conductivity and gave computer programs for the numerical solution of this problem for both the *E*-polarization (*E* parallel to the strike of the structure) and the *H*-polarization (*H* parallel to the strike of the structure) cases.

The programs given by Jones & Pascoe (1971) may be used to consider long cylinders composed of several conductivities and of arbitrary cross-section embedded in a region of uniform conductivity, but cannot be used to solve the problem in which the surrounding region is not uniform. It is important to be able to solve the more general case in which the surrounding medium is a layered one and is not necessarily the same at great distances from the conductivity inhomogeneities on both sides. In

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the previous work (Jones & Pascoe 1971) it was necessary to place the sides of the mesh over which the equations were solved far enough away from any inhomogeneity so that a uniform conducting region could be assumed at these boundaries. It is also true in the more general case presented here that the sides of the mesh must be far from any *vertical* discontinuity in conductivity so that a horizontally layered medium may be assumed at these boundaries. The subroutines for the boundary conditions for the *E*-polarization and *H*-polarization cases which we now present can be used in place of the previous boundary condition subroutines (Jones & Pascoe 1971) and allow the solution of any conductivity configuration within the conducting region as long as the above condition is met.

## 2. The boundary conditions for layered media

Jones (1971) investigated the problem of induction in a two-layered Earth model with a general layer contact topography and derived analytic expressions for the boundary. However, the analytic expressions for the fields at the boundaries in terms of the conductivities and the depth to the interface became cumbersome even for this two-layered case. Therefore, to proceed to a situation which involves more than two layers, a different approach is taken.

### (a) *E*-polarization

If we consider the same co-ordinate system as before, namely with the origin on the surface, *x* and *y* co-ordinates horizontal and the *z* co-ordinate vertically downward, (Jones & Pascoe 1971), then for a uniformly layered conducting region,  $(\partial E_x/\partial y) = 0$  everywhere. The equation which must be solved,

$$\nabla^2 E_x = i\eta^2 E_x$$

where  $\eta^2 = 4\pi\sigma\omega$  reduces to

$$\frac{\partial^2 E_x}{\partial z^2} = i\eta^2 E_x.$$

This equation has the solution

$$E_x = D_1 \exp(-\eta z\sqrt{i}) + D_2 \exp(\eta z\sqrt{i}) \text{ for } \eta \neq 0$$

and

$$E_x = D_1 + D_2 z \text{ for } \eta = 0.$$

If we now consider Fig. 1 and assume that  $E_{x|k}$  and  $(\partial E_x/\partial z)_{|k}$  are known, then by using the proper functional form in region  $\eta_j$  and the boundary conditions [the continuity of  $E_x$  and the tangential component of  $H$ ,  $(\partial E_x/\partial z)$ ] at the horizontal interface,  $k$ , then the constants  $D_1$  and  $D_2$  may be evaluated for layer  $\eta_j$ . Once  $D_1$  and  $D_2$  are known,  $E_{x|j}$  and  $(\partial E_x/\partial z)_{|j}$  may be calculated. In this way a knowledge of  $E_x$  and  $(\partial E_x/\partial z)$  on the lowest grid row allows the determination of  $E_x$  and  $(\partial E_x/\partial z)$  for the remaining grid rows.

### (b) *H*-polarization

For the *H*-polarization case a similar form of the solution is encountered:

$$H_x = D_1 \exp(-\eta z\sqrt{i}) + D_2 \exp(\eta z\sqrt{i}) \text{ for } \eta \neq 0$$

and

$$H_x = D_1 + D_2 z \text{ for } \eta = 0.$$

Again, if  $H_{x|k}$  and  $(\partial H_x/\partial z)_{|k}$  are known, the constants  $D_1$  and  $D_2$  may be determined by using the conditions of continuity of  $H$  and continuity of the tangential

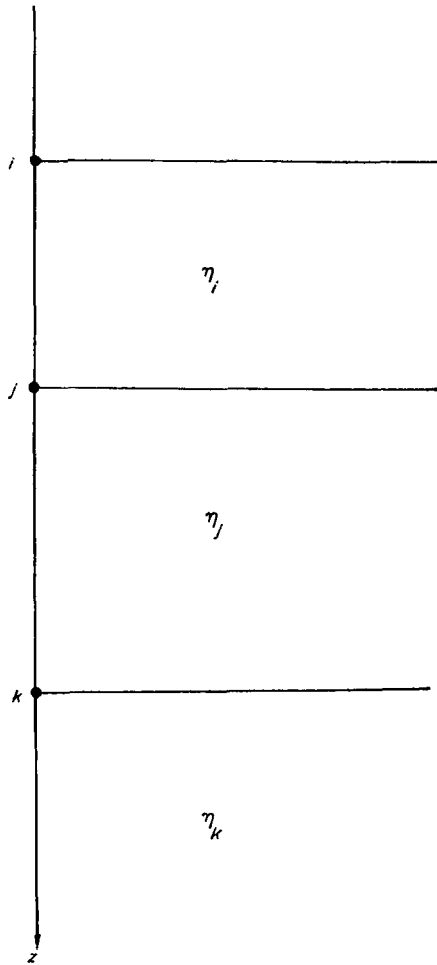


FIG. 1. Notation used to describe boundary of layered medium.

component of  $E$  on line  $k$ . After  $D_1$  and  $D_2$  are thus determined for  $\eta_j$ , we may calculate  $H_{x|j}$  and  $(\partial H_x / \partial z)_{|j}$  and proceed as described in the  $E$  polarization case.

It should be noted that to be certain that  $E_x$  (or  $H_x$ ) remains finite as  $z \rightarrow \infty$ , we have set the value of  $E_x$  (or  $H_x$ ) on the lowermost grid row according to

$$E_x = \exp(-\eta z \sqrt{i})$$

(or  $H_x = \exp(-\eta z \sqrt{i})$ );

that is  $D_1 = 1, D_2 = 0$ .

Also, in the  $H$ -polarization case,  $H_x = H_0$  everywhere in the free-space region (Jones & Price 1970).

In the  $E$ -polarization case, it is necessary to take (Jones & Price 1970)

$$H_{y|left-hand\ surface} = H_{y|right-hand\ surface}$$

and also to place the upper boundary high enough to ensure that any perturbations in  $H$  due to discontinuities in the conductor are negligible there.

Furthermore, since the above applies for a horizontally layered conducting medium, we must ensure that the boundaries are far enough away from any vertical discontinuity so that this assumption holds.

### 3. The boundary value subroutines and example

Figs 2 and 3 give the boundary value subroutine for the  $E$ -polarization case and Figs 4 and 5 for the  $H$ -polarization case. In these subroutines complex variables are used directly, and the real and imaginary parts are separated at the end to accommodate the main program.

Fig. 6 gives the conductive configuration for the example illustrated. The model is that of a layered medium with a step discontinuity. The different conductivities are illustrated by the different letters.

Fig. 7 gives the  $E$ -polarization surface values of the three components, the phase and apparent resistivity. Fig. 8 is the solution for the  $H$ -polarization case. In the model illustrated, a slightly different method of calculating the surface values using a non-linear approximation of the functional form at the surface is employed.

FURTRAN IV G COMPILER	BYCOND	04-13-71	11:56.48	PAGE 0001
0001		SUBROUTINE BYCOND (N)		10
0002		REAL K		20
0003		COMMON F(41,41),G(41,41),H(40),K(40),SCALE,FREQ,REGION(40,40)		30
0004		CCOMPLEX E(41,2),EPRIME(41,2),D1(40,2),D2(40,2),ROOTI,ENZ,EMNZ		40
0005		DIMENSION AETA(40,2),DIST(41)		50
	C			60
	C	SET THE ORIGIN FOR THE Z-AXIS ON GRID ROW 1		70
	C			80
0006		DIST(1)=0.0		90
0007		DO 110 I=2,41		100
0008	110	DIST(I)=DIST(I-1)+K(I-1)		110
0009		ROOTI=CMPLX(1./SQRT(2.),1./SQRT(2.))		120
	C			130
	C	DETERMINE AETA FOR THE LEFT AND RIGHT BOUNDARIES		140
	C			150
0010		DO 120 I=N,40		160
0011		AETA(I,1)=SQRT(REGION(I,1))		170
0012	120	AETA(I,2)=SQRT(REGION(I,40))		180
	C			190
	C	SET THE VALUES OF E AND EPRIME ON THE LAST GRID ROW		200
	C			210
0013		E(41,1)=CEXP(-DIST(41)*AETA(40,1)*ROOTI)		220
0014		E(41,2)=CEXP(-DIST(41)*AETA(40,2)*ROOTI)		230
0015		EPRIME(41,1)=-AETA(40,1)*ROOTI*CEXP(-DIST(41)*AETA(40,1)*ROOTI)		240
0016		EPRIME(41,2)=-AETA(40,2)*ROOTI*CEXP(-DIST(41)*AETA(40,2)*ROOTI)		250
	C			260
	C	SOLVE FOR REMAINING BOUNDARY VALUES		270
	C			280
0017		DO 130 J=1,2		290
0018		DO 130 I=N,4J		300
0019		ENZ=CEXP(DIST(41-I+N)*AETA(40-I+N,J)*ROOTI)		310
0020		EMNZ=CEXP(-DIST(41-I+N)*AETA(40-I+N,J)*ROOTI)		320
0021		D1(40-I+N,J)=((E(41-I+N,J)*AETA(40-I+N,J)*ROOTI*ENZ-EPRIME(41-I+N, J)*EMNZ)/(2.*AETA(40-I+N,J)*ROOTI)		330
0022		D2(40-I+N,J)=((EPRIME(41-I+N,J)*EMNZ+E(41-I+N,J)*AETA(40-I+N,J)*RO OTI*EMNZ)/(2.*AETA(40-I+N,J)*ROOTI)		350
0023		ENZ=CEXP(DIST(40-I+N)*AETA(40-I+N,J)*ROOTI)		370
0024		EMNZ=CEXP(-DIST(40-I+N)*AETA(40-I+N,J)*ROOTI)		380
0025		E(40-I+N,J)=D1(40-I+N,J)*EMNZ+D2(40-I+N,J)*ENZ		390
0026	130	EPRIME(40-I+N,J)=-AETA(40-I+N,J)*ROOTI*D1(40-I+N,J)*EMNZ+AETA(40-I +N,J)*ROOTI*D2(40-I+N,J)*ENZ		400
0027		L=N-1		420
0028		DO 140 J=1,2		430
0029		DO 140 I=1,L		440
0030		D1(N-I,J)=E(N-I+1,J)-EPRIME(N-I+1,J)*DIST(N-I+1)		450
0031		D2(N-I,J)=EPRIME(N-I+1,J)		460
0032		E(N-I,J)=D1(N-I,J)+D2(N-I,J)*DIST(N-I)		470
0033	140	EPRIME(N-I,J)=D2(N-I,J)		480
	C			490
	C	NORMALISE TO (1.0,0.0) ON THE SURFACE ON THE LEFT-HAND-SIDE		500
	C			510
0034		ROOTI=E(N,1)		520
0035		DO 150 I=1,41		530
0036		E(I,1)=E(I,1)/ROOTI		540
0037	150	EPRIME(I,1)=EPRIME(I,1)/ROOTI		550

FIG. 2.  $E$ -polarization boundary condition subroutine.

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      C
      C   ADJUST THE RIGHT-HAND-SIDE TO THE PROPER VALUE
      C
0033      ROOTI=EPRIME(N,2)/EPRIME(N,1)
0039      DO 160 I=1,41
0040      E(I,2)=E(I,2)/ROOTI
0041      160  EPRIME(I,2)=EPRIME(I,2)/ROOTI
      C
      C   SET THE BOUNDARY OF F & G AND INTERPOLATE LINEARLY ACROSS THE GRID
      C
0042      DU 170 I=1,41
0043      F(I,1)=REAL(E(I,1))
0044      F(I,41)=REAL(E(I,2))
0045      G(I,1)=AIMAG(E(I,1))
0046      170  G(I,41)=AIMAG(E(I,2))
0047      DIST(I)=0.0
0048      DU 180 I=2,41
0049      180  DIST(I)=DIST(I-1)+H(I-1)
0050      DO 190 I=1,41
0051      DFDY=(F(I,41)-F(I,1))/DIST(41)
0052      DGDY=(G(I,41)-G(I,1))/DIST(41)
0053      DC 190 J=2,40
0054      F(I,J)=F(I,1)+(DFDY*DIST(J))
0055      190  G(I,J)=G(I,1)+(DGDY*DIST(J))
0056      RETURN
0057      END
    
```

560  
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FIG. 3. *E*-polarization boundary condition subroutine.

**4. Calculation of the surface values**

In the previous work (Jones & Pascoe 1971), it was found that some error was encountered in the calculation of  $E_y$  in the *H*-polarization case and  $H_y$  in the *E*-polarization case. This error is exhibited by a difference between the computed value of the apparent resistivity ( $\rho_A$ ) on the surface over the uniform conducting regions at the extremities of the mesh and the value expected there.

Since for the *H*-polarization case

$$E_y = \frac{1}{4\pi\sigma} \frac{\partial H_x}{\partial z}$$

and for the *E*-polarization case

$$H_y = - \frac{1}{i\omega} \frac{\partial E_x}{\partial z}$$

(Jones & Price 1970), the components  $E_y$  and  $H_y$  were calculated by taking finite differences in the *z* direction. This approximation to the derivative is adequate when the grid spacing is not too large. However, a better approximation which is independent of the grid spacing and which uses the true form of the function can be applied.

For the conducting region, the usual method for approximating  $(\partial E_x/\partial z)$  (or  $(\partial H_x/\partial z)$ ) at the surface is by using a linear approximation to the derivative. For example, from Fig. 9 we would have

$$\frac{\partial F}{\partial z} \Big|_0 \simeq \frac{\Delta F}{\Delta z} \Big|_0 = \frac{F_1 - F_0}{z_1 - z_0},$$

where  $F$  equals  $E_x$  or  $H_x$ . This is a reasonable approximation to the derivative when the grid spacing is small, since the derivative is the value of this gradient in the limit as  $z_1 \rightarrow z_0$ .

FORTRAN IV G COMPILER	BYCOND	04-13-71	12:00.35	PAGE 0001	
0001	SUBROUTINE BYCOND (N)				10
0002	REAL K				20
0003	COMMON F(41,41),G(41,41),H(40),K(40),SCALE,FREQ,REGION(40,40)				30
0004	COMPLEX E(41,2),EPRIME(41,2),D1(40,2),D2(40,2),ROOTI,ENZ,EMNZ				40
0005	DIMENSION AETA(41,2),DIST(41)				50
	C				60
	C SET THE ORIGIN FOR THE Z-AXIS ON GRID ROW 1				70
	C				80
0006	DIST(1)=0.0				90
0007	DO 110 I=2,41				100
0008	110 DIST(I)=DIST(I-1)+K(I-1)				110
0009	ROOTI=CMPLX(1./SQRT(2.),1./SQRT(2.))				120
	C				130
	C DETERMINE AETA FOR THE LEFT AND RIGHT BOUNDARIES				140
	C				150
0010	DO 120 I=N,40				160
0011	AETA(I,1)=SQRT(REGION(I,1))				170
0012	120 AETA(I,2)=SQRT(REGION(I,40))				180
0013	AETA(41,1)=AETA(40,1)				184
0014	AETA(41,2)=AETA(40,2)				188
	C				190
	C SET THE VALUES OF E AND EPRIME ON THE LAST GRID ROW				200
	C				210
0015	E(41,1)=CEXP(-DIST(41)*AETA(40,1)*ROOTI)				220
0016	E(41,2)=CEXP(-DIST(41)*AETA(40,2)*ROOTI)				230
0017	EPRIME(41,1)=-AETA(40,1)*ROOTI*CEXP(-DIST(41)*AETA(40,1)*ROOTI)				240
0018	EPRIME(41,2)=-AETA(40,2)*ROOTI*CEXP(-DIST(41)*AETA(40,2)*ROOTI)				250
	C				260
	C SOLVE FOR REMAINING BOUNDARY VALUES				270
	C				280
0019	DO 130 J=1,2				290
0020	DO 130 I=N,40				300
0021	ENZ=CEXP(DIST(41-I+N)*AETA(40-I+N,J)*ROOTI)				310
0022	EMNZ=CEXP(-DIST(41-I+N)*AETA(40-I+N,J)*ROOTI)				320
0023	D1(40-I+N,J)=((E(41-I+N,J)*AETA(40-I+N,J)*ROOTI*ENZ-((AETA(40-I+N,1J)*AETA(41-I+N,J)**2)*EPRIME(41-I+N,J)*ENZ)/(2.*AETA(40-I+N,J)*RO1CGTI))				350
0024	D2(40-I+N,J)=(((AETA(40-I+N,J)/AETA(41-I+N,J)**2)*EPRIME(41-I+N,1J)*EMNZ+E(41-I+N,J)*AETA(40-I+N,J)*ROOTI*EMNZ)/(2.*AETA(40-I+N,J)*1RCGTI))				360
0025	ENZ=CEXP(DIST(40-I+N)*AETA(40-I+N,J)*ROOTI)				370
0026	EMNZ=CEXP(-DIST(40-I+N)*AETA(40-I+N,J)*ROOTI)				380
0027	E(40-I+N,J)=D1(40-I+N,J)*EMNZ+D2(40-I+N,J)*ENZ				390
0028	130 EPRIME(40-I+N,J)=-AETA(40-I+N,J)*ROOTI*D1(40-I+N,J)*EMNZ+AETA(40-I+N,J)*ROOTI*D2(40-I+N,J)*ENZ				400
0029	L=N-1				410
0030	DO 140 J=1,2				420
0031	DO 140 I=1,L				440
0032	EPRIME(N-I,J)=0.0				460
0033	140 E(N-I,J)=E(N-I+1,J)				470
	C				490
	C NORMALISE TO (1.0,0.0) ON THE SURFACE ON THE LEFT-HAND-SIDE				500
	C				510
0034	ROOTI=E(N,1)				520
0035	DO 150 I=1,41				530

FIG. 4. *H*-Polarization boundary condition subroutine.

FOKTRAN IV G COMPILER	BYCOND	04-13-71	12:00.35	PAGE 0002
0036		E(I,1)=E(I,1)/ROOTI		540
0037	150	EPRIME(I,1)=EPRIME(I,1)/ROOTI		550
	C			560
	C	ADJUST THE RIGHT-HAND-SIDE TO THE PROPER VALUE		570
	C			580
0038		ROOTI=E(N,2)/E(N,1)		590
0039		DU 160 I=1,41		600
0040		E(I,2)=E(I,2)/ROOTI		610
0041	160	EPRIME(I,2)=EPRIME(I,2)/ROOTI		620
	C			630
	C	SET THE BOUNDARY OF F & G AND INTERPOLATE LINEARLY ACROSS THE GRID		640
	C			650
0042		DU 170 I=1,41		660
0043		F(I,1)=REAL(E(I,1))		670
0044		F(I,41)=REAL(E(I,2))		680
0045		G(I,1)=AIMAG(E(I,1))		690
0046	170	G(I,41)=AIMAG(E(I,2))		700
0047		DIST(I)=0.0		710
0048		DO 180 I=2,41		720
0049	180	DIST(I)=DIST(I-1)*H(I-1)		730
0050		DU 190 I=1,41		740
0051		DFDY=(F(I,41)-F(I,1))/DIST(41)		750
0052		DGDY=(G(I,41)-G(I,1))/DIST(41)		760
0053		DC 190 J=2,40		770
0054		F(I,J)=F(I,1)+(DFDY*DIST(J))		780
0055	190	G(I,J)=G(I,1)+(DGDY*DIST(J))		790
0056		RETURN		800
0057		END		810*

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FIG. 5. *H*-polarization boundary condition subroutine.

However, in most instances, the grid spacing is such that the above is only a first approximation to the derivative. If we consider the true form of the function we may obtain a better numerical value for  $(\partial F/\partial z)_{10}$ . For Fig. 1 we have the usual functional form for the conducting region ( $\eta_1$ ):

$$F(z) = D_1 \exp(-\eta_1 z\sqrt{(i)}) + D_2 \exp(\eta_1 z\sqrt{(i)})$$

where  $D_1$  and  $D_2$  are constants.

If we know  $F_{10}$  and  $F_{11}$ ,  $D_1$  and  $D_2$  may be calculated numerically. The value of  $(\partial F/\partial z)_{10}$  may then be determined from

$$\frac{\partial F}{\partial z} \Big|_{10} = -\eta_1 \sqrt{(i)} D_1 + \eta_1 \sqrt{(i)} D_2,$$

where the origin of the *z*-axis has been taken at the surface. This will give a more accurate value for  $(\partial F/\partial z)_{10}$  over a uniformly stratified conducting region, and is likely to be at least as accurate as the linear approximation above regions where lateral discontinuities in conductivity occur. It must be applied with care near regions with discontinuities at the surface, since the above functional form may not necessarily apply near such regions.

### 5. The new surface value subroutine and comparison with the linear approximation

Figs 10 and 11 give the new surface value subroutine for the *H*-polarization case which may be used to replace the previous subroutine (Jones & Pascoe 1971) if the new approximation for the surface values is desired. The altered or inserted statements are those numbered 12–23. The *E*-polarization subroutine would require similar changes. Also, in the new *H*-polarization subroutine a statement (No. 35) is incorporated to indicate where a discontinuity exists at the surface.

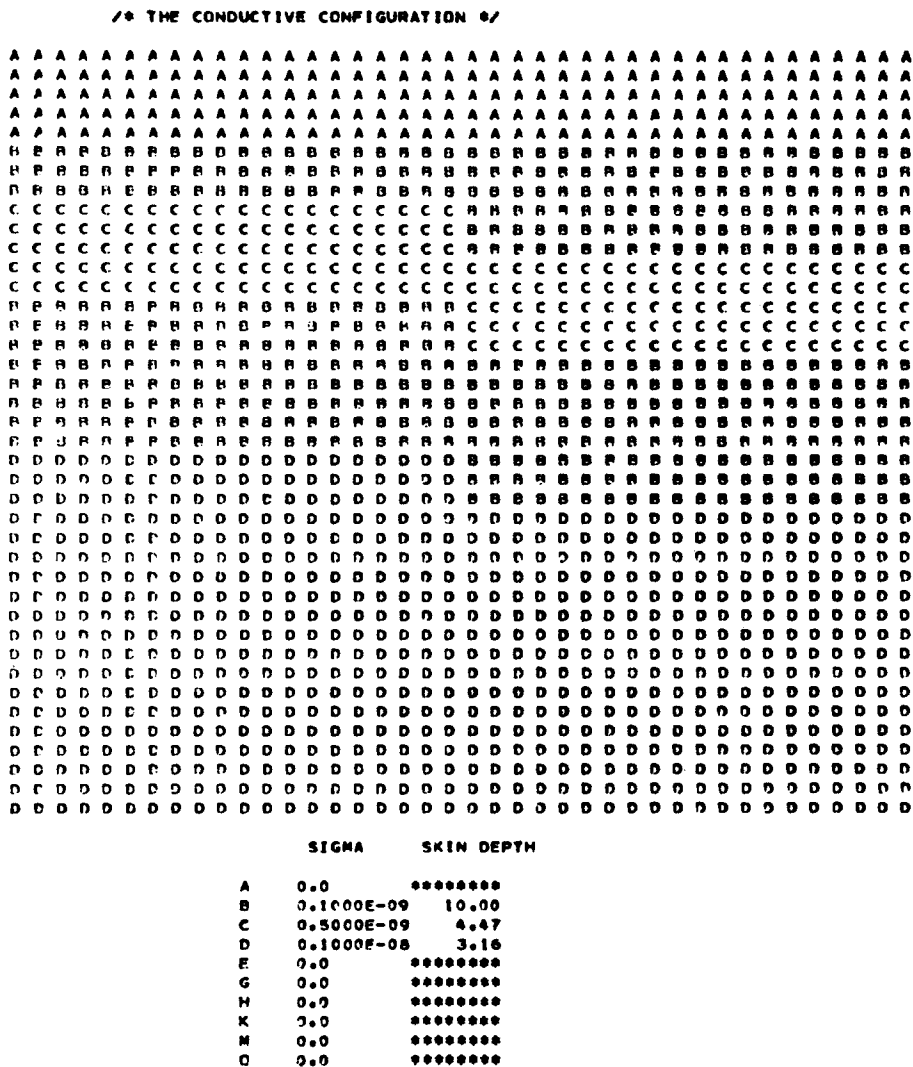


FIG. 6. Conductive configuration for example illustrated. Horizontal (H) and vertical (K) grid dimensions and skin depth are in multiples of scale (cm). Frequency used is 0.000253 Hz.



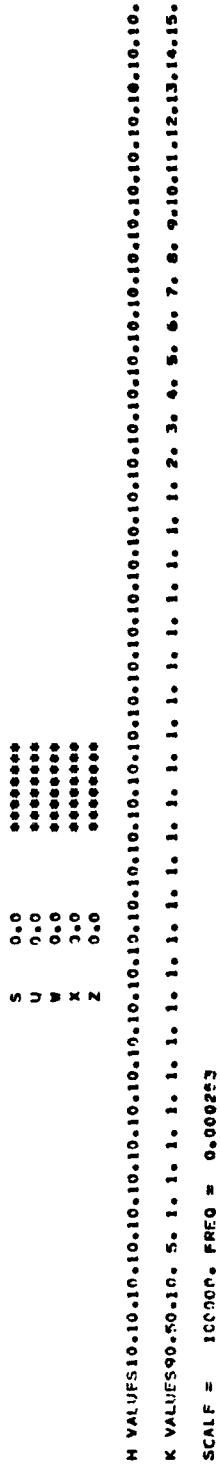


FIG. 6. Continued.

/\* E-POLARISATION \*/

/\* FPS = 0.000100 MAXIMUM NO. OF ITERATIONS = 5000/\*

/\* STOPPED ON ITERATION 238 \*/

/\* SURFACE VALUES \*/

	AME	AMHY	AMHZ	DPHASE	DPHANY	DPHAHZ	APPRES
2	0.999	1.000	0.002	0.006	0.000	-2.743	0.466E 10
3	0.999	1.000	0.000	0.005	-0.001	-2.957	0.466E 10
4	0.999	1.000	0.000	0.004	-0.001	-2.431	0.466E 10
5	0.999	1.000	0.000	0.004	-0.001	-1.645	0.466E 10
6	0.999	1.000	0.000	0.004	-0.001	-1.393	0.466E 10
7	0.999	1.000	0.000	0.004	-0.002	-1.302	0.466E 10
8	1.000	1.000	0.000	0.004	-0.002	-1.273	0.466E 10
9	1.000	1.001	0.000	0.004	-0.002	-1.273	0.466E 10
10	1.000	1.001	0.000	0.004	-0.002	-1.291	0.466E 10
11	1.001	1.001	0.000	0.004	-0.002	-1.325	0.466E 10
12	1.001	1.001	0.000	0.004	-0.002	-1.173	0.466E 10
13	1.002	1.002	0.000	0.003	-0.003	-1.435	0.466E 10
14	1.002	1.002	0.000	0.003	-0.003	-1.407	0.466E 10
15	1.003	1.003	0.001	0.003	-0.004	-1.579	0.466E 10
16	1.004	1.004	0.001	0.002	-0.004	-1.641	0.466E 10
17	1.006	1.006	0.001	0.001	-0.005	-1.692	0.466E 10
18	1.006	1.009	0.003	-0.002	-0.007	-1.700	0.466E 10
19	1.015	1.015	0.004	-0.007	-0.011	-1.514	0.466E 10
20	1.038	1.030	0.030	-0.017	-0.012	-1.072	0.474E 10
21	1.137	1.003	0.052	-0.011	0.003	-0.849	0.594E 10
22	1.251	0.969	0.037	0.016	0.007	-0.836	0.779E 10
23	1.287	0.979	0.012	0.011	0.009	-1.227	0.806E 10
24	1.300	0.987	0.005	0.007	0.007	-1.473	0.809E 10
25	1.305	0.990	0.002	0.005	0.005	-1.566	0.809E 10
26	1.307	0.992	0.001	0.004	0.004	-1.599	0.809E 10
27	1.309	0.994	0.001	0.003	0.003	-1.620	0.809E 10
28	1.310	0.995	0.001	0.002	0.002	-1.637	0.809E 10
29	1.311	0.995	0.000	0.002	0.002	-1.648	0.809E 10
30	1.312	0.996	0.000	0.001	0.001	-1.653	0.809E 10
31	1.312	0.996	0.000	0.001	0.001	-1.654	0.809E 10
32	1.313	0.996	0.000	0.001	0.001	-1.651	0.809E 10
33	1.313	0.997	0.000	0.000	0.001	-1.644	0.809E 10
34	1.313	0.997	0.000	0.000	0.000	-1.627	0.809E 10
35	1.314	0.997	0.000	0.000	0.000	-1.596	0.809E 10
36	1.314	0.993	0.000	-0.000	-0.000	-1.537	0.809E 10
37	1.315	0.994	0.000	-0.000	-0.000	-1.420	0.809E 10
38	1.315	0.999	0.000	-0.000	-0.000	-1.181	0.809E 10
39	1.316	0.999	0.000	-0.000	-0.000	-0.715	0.809E 10
40	1.317	0.999	0.001	0.0	0.0	0.0	0.810E 10

FIG. 7. E-polarization surface values. Amplitudes of components normalized, phase differences in radians, apparent resistivity in emu

/\* H-POLARISATION \*/

/\* FPR = 0.000100 MAXIMUM NO. OF ITERATIONS = 5000 \*/

/\* STOPPED ON ITERATION 167 \*/

/\* SURFACE VALUES \*/

	AMH	AMFY	AMEZ	DPHSH	DPHAFY	DPHAEZ	APRES
2	1.0000	1.0000	0.0	0.0	0.011	0.0	0.470E 10
3	1.0000	1.0000	0.0	0.0	0.011	0.0	0.470E 10
4	1.0000	1.0000	0.0	0.0	0.012	0.0	0.470E 10
5	1.0000	1.0000	0.0	0.0	0.012	0.0	0.470E 10
6	1.0000	1.0000	0.0	0.0	0.012	0.0	0.470E 10
7	1.0000	1.0000	0.0	0.0	0.012	0.0	0.470E 10
8	1.0000	1.0000	0.0	0.0	0.012	0.0	0.470E 10
9	1.0000	0.9999	0.0	0.0	0.012	0.0	0.470E 10
10	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
11	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
12	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
13	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
14	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
15	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
16	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
17	1.0000	0.9999	0.0	0.0	0.012	0.0	0.469E 10
18	1.0000	0.9999	0.0	0.0	0.011	0.0	0.469E 10
19	1.0000	1.0000	0.0	0.0	0.011	0.0	0.470E 10
20	1.0000	1.0000	0.0	0.0	0.015	0.0	0.478E 10
21	1.0000	1.166	0.0	0.0	0.005	0.0	0.534E 10
22	1.0000	1.312	0.0	0.0	-0.011	0.0	0.809E 10
23	1.0000	1.317	0.0	0.0	-0.001	0.0	0.816E 10
24	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
25	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
26	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
27	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
28	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
29	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
30	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
31	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
32	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
33	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
34	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
35	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
36	1.0000	1.318	0.0	0.0	-0.000	0.0	0.817E 10
37	1.0000	1.313	0.0	0.0	-0.000	0.0	0.816E 10
38	1.0000	1.318	0.0	0.0	-0.000	0.0	0.816E 10
39	1.0000	1.318	0.0	0.0	-0.000	0.0	0.816E 10
40	1.0000	1.317	0.0	0.0	0.0	0.0	0.816E 10

FIG. 8. H-polarization surface values. Amplitudes of components normalized, phase differences in radians, apparent resistivity in emu.

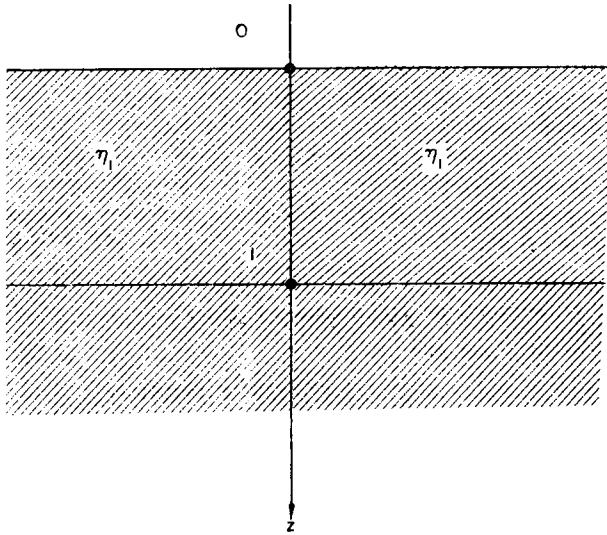


FIG. 9. Notation used to describe surface value calculations.

In Fig. 12 the surface values of apparent resistivity computed by the improved method (solid curves) are compared with the surface values obtained before (dashed curves), (Jones & Pascoe 1971). With the new calculation, the value of apparent resistivity near the boundaries of the mesh approaches the expected value for the uniform subsurface. In the previous calculations, the value of apparent resistivity differed by approximately 6 per cent in the *E*-polarization case and by approximately 11 per cent in the *H*-polarization case from the expected values. The profiles also show some change across the conductor surface when the new and old calculations are compared.

## 6. Dimensional considerations and the iterative procedure

The use of this new method of calculating surface values and the above considerations about accuracy, made it desirable to obtain a better 'feel' for the mode of solution and the relationship between the convergence criteria applied and the various parameters used. To do this, a uniform subsurface case was chosen and the model was run for various convergence conditions and grid-size to skin depth ratios for a uniform square grid. Since in the initialization procedure the boundary values are carried horizontally throughout the mesh, the number of iterations required to satisfy the convergence condition is a measure of the accuracy of the result. Fig. 13 shows a model constructed to illustrate the relationship between the number of iterations (proportional to the height of the rod; terminated at a maximum of 150 iterations), the residual (EPS) and the ratio of grid size to skin depth (GS/SD). It is seen that the iteration should be carried out when the grid size to skin depth ratio is small.

In the solution of the problem the equations are solved by the numerical iterative method over several conductive regions. In each region there is a basic functional form which sets limits on the grid spacing we can use there in order to accommodate the numerical difference equations. For example, in the free-space region the function is approximately linear in both directions and consequently a relatively large grid

FORTRAN IV G COMPILER	SURFVL	04-29-71	09:20:03	PAGE 0001	
0001	SUBROUTINE SURFVL (L)				10
0002	REAL K				20
0003	DIMENSION AMH(41), AMEY(41), AMEZ(41), DPHASH(41), DPHAEY(41), DPHAEZ(41), APPRES(41)				30
0004	COMMON F(41,41),G(41,41),H(40),K(40),SCALE,FREQ,REGION(40,40)				40
0005	COMPLEX ROOTI,H0,H1,DERIVZ				50
	C				55
	C				60
0006	PI=4.0*ATAN(1.0)				60
0007	OMEGA=2.0*PI*FREQ				90
0008	WRITE (6,140)				100
0009	WRITE (6,150)				110
0010	I=L				120
	C				130
0011	DO 110 J=2,40				140
0012	AETAL=SQRT(REGION(I,J-1))				150
0013	AETAR=SQRT(REGION(I,J))				160
0014	AETA=(AETAL+AETAR)/2.0				170
0015	ROOTI=CMPLX(1./SQRT(2.),1./SQRT(2.))				180
0016	H0=CMPLX(F(I,J),G(I,J))				190
0017	H1=CMPLX(F(I+1,J),G(I+1,J))				200
0018	DERIVZ=(AETA*ROOTI*(2.0*H1-H0*(CEXP(AETA*K(I)*ROOTI)+CFXP(-ALTA*K(I)*ROOTI)))/CFXP(ALTA*K(I)*ROOTI)-CEXP(-ALTA*K(I)*ROOTI))				210
0019	DPHASH(J)=ATAN2(G(I,J),F(I,J))				230
0020	DPHAEY(J)=ATAN2(AIMAG(DERIVZ),REAL(DERIVZ))				240
0021	DPHAEZ(J)=0.0				250
0022	AMH(J)=SQRT(F(I,J)**2+G(I,J)**2)				260
0023	AMEY(J)=(OMEGA/(AETA**2))*CABS(DERIVZ)				270
0024	AMEZ(J)={(2.*OMEGA)/(REGION(I,J)+REGION(I,J-1))*SQRT(((F(I,J+1)-F(I,J-1))/(H(J)+H(J-1)))**2+((G(I,J+1)-G(I,J-1))/(H(J)+H(J-1)))**2)}				290
0025	APPRES(J)=(2.0/FREQ)*((AMEY(J)/AMH(J))**2)	110			310
	C				320
	C	THE COMPONENTS AMEY AND AMEZ ARE NORMALIZED WITH RESPECT TO			330
	C	THE FIELD AT INFINITY (POINT 2) AND PHASE DIFFERENCES AMF			340
	C	CALCULATED RELATIVE TO POINT 40			350
0026	AMF=SQRT(AMEY(2)**2+AMEZ(2)**2)				360
0027	DO 120 J=2,40				370
0028	AMEY(J)=AMEY(J)/AMF				380
0029	AMEZ(J)=AMEZ(J)/AMF				390
0030	DPHASH(J)=DPHASH(J)-DPHASH(40)				400
0031	DPHAEY(J)=DPHAEY(J)-DPHAEY(40)				410
0032	DPHAEZ(J)=DPHAEZ(J)-DPHAEZ(40)	120			420
	C				430
0033	DO 130 J=2,40				440
0034	WRITE (6,160) J,AMH(J),AMEY(J),AMEZ(J),DPHASH(J),DPHAEY(J),DPHAEZ(J),APPRES(J)				450
0035	IF (REGION(L,J-1).NE.REGION(L,J)) WRITE (6,170)				460
0036	CONTINUE	130			468
0037	REWIND 3				470
0038	WRITE (3) F,G,H,SCALE,AMH,AMEY,AMEZ,DPHASH,DPHAEY,DPHAEZ,APPRES				480
0039	RETURN				490
	C				500
	C				510
	C				520
0040	140 FORMAT (1H0,40X,20H/* SURFACE VALUES **//)				530

FIG. 10. H-polarization surface value subroutine.

```

0041_   150   FORMAT (1H0,T9,'AMH',T21,'AMEY',T33,'AMEZ',T45,'DPHASH',T57,'DPHAE
          1Y',T69,'DPHAEZ',T81,'APPRES'//)
0042   160   FORMAT (1H ,12,6(2X,F10.3),/12.3)
0043   170   FORMAT (1H+,T95,'DISCONTINUITY')
0044           END

```

540  
550  
560  
565  
570\*

TOTAL MEMORY REQUIRMENTS 000E88 BYTES

FIG. 11. *H*-polarization surface value subroutine.

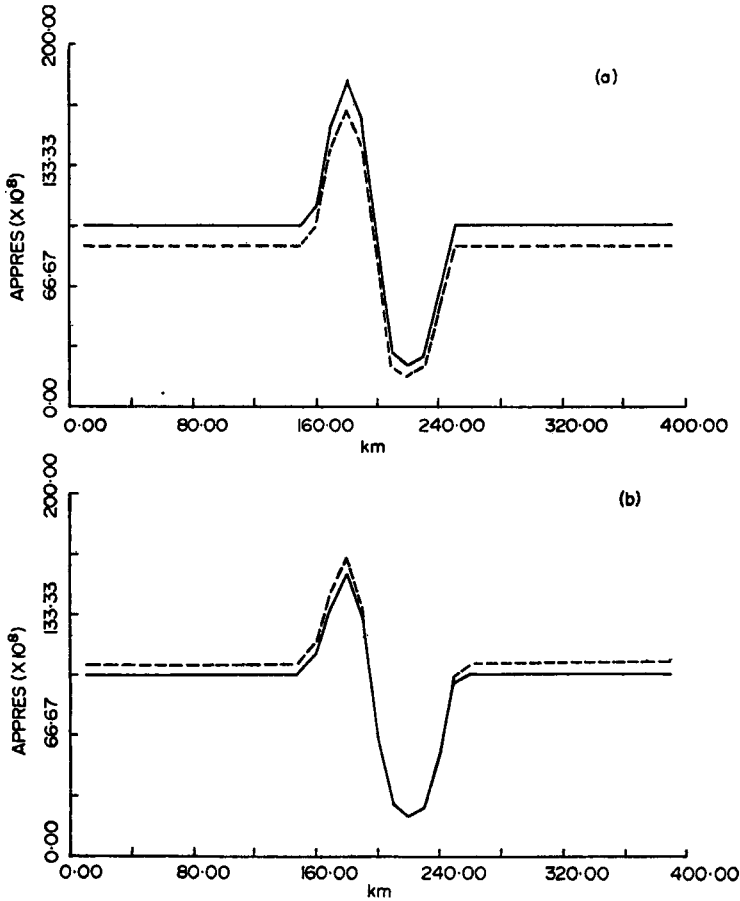


FIG. 12. Apparent resistivity profiles for the previous (dashed lines) calculation and the improved (solid lines) calculation (emu). (a) *H*-polarization. (b) *E*-polarization.

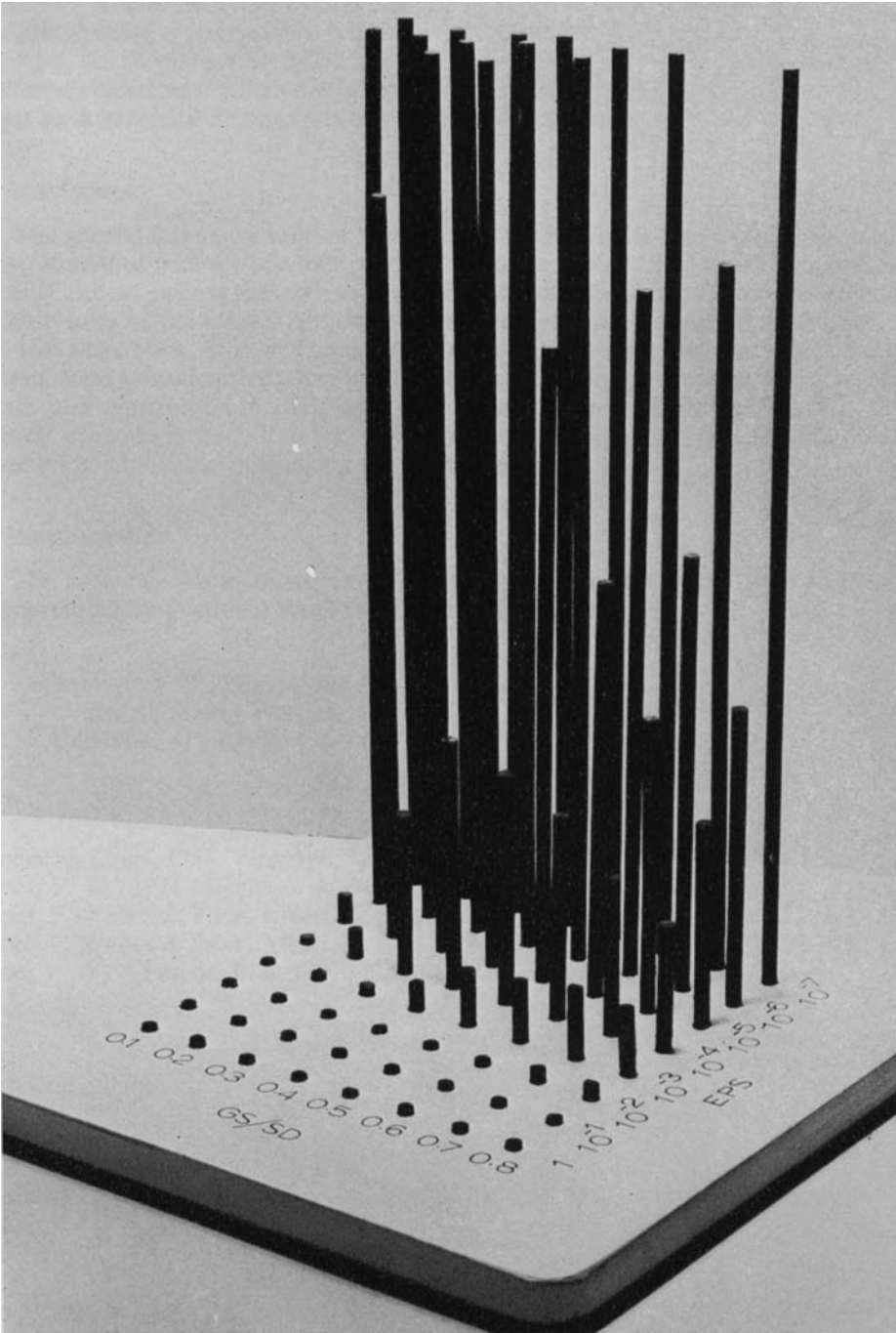


FIG. 13. Model constructed to investigate iteration procedure and relationship of various parameters.

[Facing p. 192

spacing can be used. In the conducting regions we have taken the phenomenon as approximately linear in the horizontal direction, but exponential in the vertical direction. Therefore, in the conducting region, in order to fit the form of the function the grid spacing is restricted in the vertical direction.

Also, the solution is based on the assumption that as  $y \rightarrow \pm \infty$ , the subsurface is uniformly stratified. To accommodate this restriction in the model the boundaries must be several skin depths from any vertical discontinuity.

## 7. Conclusions

The general boundary subroutines as presented provide a two-dimensional model when combined with the previous program (Jones & Pascoe 1971) which is completely general, subject only to the restriction that a uniformly stratified subsurface is required at both sides of the mesh. Although the subsurface must be uniformly stratified at the two extremities, it is not necessary that both sides have identical stratification. In use, some consideration must be given to the relationship between grid size, skin depth and dimensions of the mesh when particular models are considered. This general program is flexible to use and provides a modelling technique which should be of considerable use in studying observed phenomena.

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