

Techniques of Water-Resources Investigations  
of the United States Geological Survey

**Chapter C1**

**FINITE-DIFFERENCE MODEL FOR  
AQUIFER SIMULATION IN  
TWO DIMENSIONS  
WITH RESULTS OF  
NUMERICAL EXPERIMENTS**

**By P. C. Trescott, G. F. Pinder, and S. P. Larson**

**Book 7**

**AUTOMATED DATA PROCESSING AND COMPUTATIONS**

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## PREFACE

The series of manuals on techniques describes procedures for planning and executing specialized work in water-resources investigations. The material is grouped under major headings called books and further subdivided into sections and chapters; section C of Book 7 is on computer programs.

“Finite-difference model for aquifer simulation in two dimensions with results of numerical experiments” supersedes the report published in 1970 entitled, “A digital model for aquifer evaluation” by G. F. Pinder as Chapter C1 of Book 7. The new Chapter C1 represents a significant improvement in the computational capability to solve the flow equations and has greater flexibility in the hydrologic situations that can be simulated.

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# FINITE-DIFFERENCE MODEL FOR AQUIFER SIMULATION IN TWO DIMENSIONS WITH RESULTS OF NUMERICAL EXPERIMENTS

By P. C. Trescott, G. F. Pinder, and S. P. Larson

## Abstract

The model will simulate ground-water flow in an artesian aquifer, a water-table aquifer, or a combined artesian and water-table aquifer. The aquifer may be heterogeneous and anisotropic and have irregular boundaries. The source term in the flow equation may include well discharge, constant recharge, leakage from confining beds in which the effects of storage are considered, and evapotranspiration as a linear function of depth to water.

The theoretical development includes presentation of the appropriate flow equations and derivation of the finite-difference approximations (written for a variable grid). The documentation emphasizes the numerical techniques that can be used for solving the simultaneous equations and describes the results of numerical experiments using these techniques. Of the three numerical techniques available in the model, the strongly implicit procedure, in general, requires less computer time and has fewer numerical difficulties than do the iterative alternating direction implicit procedure and line successive overrelaxation (which includes a two-dimensional correction procedure to accelerate convergence).

The documentation includes a flow chart, program listing, an example simulation, and sections on designing an aquifer model and requirements for data input. It illustrates how model results can be presented on the line printer and pen plotters with a program that utilizes the graphical display software available from the Geological Survey Computer Center Division. In addition the model includes options for reading input data from a disk and writing intermediate results on a disk.

## Introduction

The finite-difference aquifer model documented in this report is designed to simulate in two dimensions the response of an aquifer to an imposed stress. The aquifer may be

artesian, water table, or a combination of artesian and water table; it may be heterogeneous and anisotropic and have irregular boundaries. The model permits leakage from confining beds in which the effects of storage are considered, constant recharge, evapotranspiration as a linear function of depth to water, and well discharge. Although it was not designed for cross-sectional problems, the model has been used with some success for this type of simulation.

The aquifer simulator has evolved from Pinder's (1970) original model and modifications by Pinder (1969) and Trescott (1973). The model documented by Trescott (1973) incorporates several features described by Prickett and Lonquist (1971) and has been applied to a variety of aquifer simulation problems by various users. The model described in this report is basically the same as the 1973 version but includes minor modifications to the logic and data input. In addition, the user may choose an equation solving scheme from among the alternating direction implicit procedure, line successive overrelaxation, and the strongly implicit procedure. The program is arranged so that other techniques for solving simultaneous equations can be coded and substituted for the iterative techniques included with the model.

The documentation is intended to be reasonably self contained, but it assumes that the user has an elementary knowledge of the physics of ground-water flow, finite-difference methods of solving partial differential

equations, matrix algebra, and the FORTRAN IV language.

## Theoretical Development

### Ground-water flow equation

The partial differential equation of ground-water flow in a confined aquifer in two dimensions may be written as

$$\frac{\partial}{\partial x}(T_{xx}\frac{\partial h}{\partial x}) + \frac{\partial}{\partial x}(T_{xy}\frac{\partial h}{\partial y}) + \frac{\partial}{\partial y}(T_{yx}\frac{\partial h}{\partial x}) + \frac{\partial}{\partial y}(T_{yy}\frac{\partial h}{\partial y}) = S\frac{\partial h}{\partial t} + W(x,y,t) \quad (1)$$

in which

$T_{xx}$ ,  $T_{xy}$ ,  $T_{yx}$ ,  $T_{yy}$  are the components of the transmissivity tensor ( $L^2t^{-1}$ );

$h$  is hydraulic head ( $L$ );

$S$  is the storage coefficient (dimensionless);

$W(x, y, t)$  is the volumetric flux of recharge or withdrawal per unit surface area of the aquifer ( $Lt^{-1}$ ).

The reader is referred to Pinder and Bredehoeft (1968) for development and discussion of equation 1. In the simulation model, equation 1 is simplified by assuming that the Cartesian coordinate axes  $x$  and  $y$  are aligned with the principal components of the transmissivity tensor,  $T_{xx}$  and  $T_{yy}$ , giving

$$\frac{\partial}{\partial x}(T_{xx}\frac{\partial h}{\partial x}) + \frac{\partial}{\partial y}(T_{yy}\frac{\partial h}{\partial y}) = S\frac{\partial h}{\partial t} + W(x,y,t). \quad (2)$$

In water-table aquifers, transmissivity is a function of head. Assuming that the coordinate axes are co-linear with the principal components of the hydraulic conductivity tensor, the flow equation may be expressed as (Bredehoeft and Pinder, 1970)

$$\frac{\partial}{\partial x}(K_{xx}b\frac{\partial h}{\partial x}) + \frac{\partial}{\partial y}(K_{yy}b\frac{\partial h}{\partial y}) = S_y\frac{\partial h}{\partial t} + W(x,y,t) \quad (3)$$

in which

$K_{xx}$ ,  $K_{yy}$  are the principal components of the hydraulic conductivity tensor ( $Lt^{-1}$ );

$S_y$  is the specific yield of the aquifer (dimensionless);

$b$  is the saturated thickness of the aquifer ( $L$ ).

### Finite-difference approximations

In order to solve equation 2 or 3 for a heterogeneous aquifer with irregular boundaries, one approach is to subdivide the region into rectangular blocks in which the aquifer properties are assumed to be uniform. The continuous derivatives in equations 2 and 3 are replaced by finite-difference approximations for the derivatives at a point (the node at the center of the block). The result is  $N$  equations in  $N$  unknowns (head values at the nodes) where  $N$  is the number of blocks representing the aquifer.

Utilizing a block-centered, finite-difference grid in which variable grid spacing is permitted (fig. 1), equation 2 may be approximated as

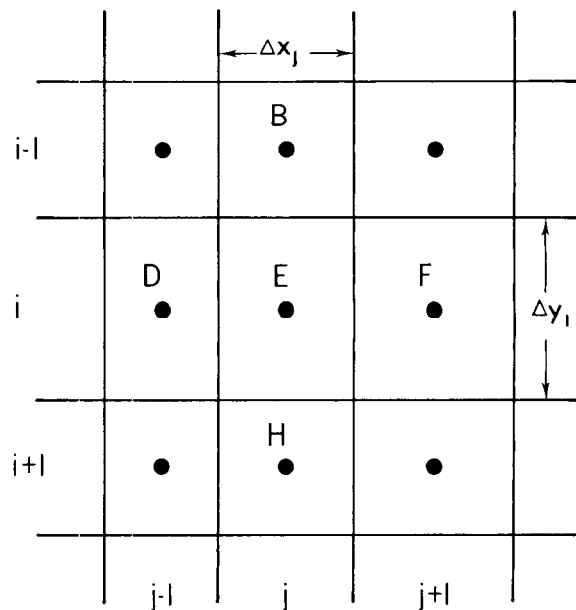


FIGURE 1.—Index scheme for finite-difference grid and coefficients of finite-difference equation written for node  $(i, j)$ .

$$\begin{aligned} & \frac{1}{\Delta x_j} \left[ \left( T_{xx} \frac{\partial h}{\partial x} \right)_{i,j+\frac{1}{2}} - \left( T_{xx} \frac{\partial h}{\partial x} \right)_{i,j-\frac{1}{2}} \right] \\ & + \frac{1}{\Delta y_i} \left[ \left( T_{yy} \frac{\partial h}{\partial y} \right)_{i+\frac{1}{2},j} - \left( T_{yy} \frac{\partial h}{\partial y} \right)_{i-\frac{1}{2},j} \right] \\ & = \frac{S_{i,j}}{\Delta t} (h_{i,j,k} - h_{i,j,k-1}) + W_{i,j,k} \end{aligned} \quad (4)$$

in which

$\Delta x_j$  is the space increment in the  $x$  direc-

tion for column  $j$  as shown in figure 1 ( $L$ );

$\Delta y_i$  is in the space increment in the  $y$  direction for row  $i$  as shown in figure 1 ( $L$ );

$\Delta t$  is the time increment ( $t$ );

$i$  is the index in the  $y$  dimension;

$j$  is the index in the  $x$  dimension;

$k$  is the time index.

Equation 4 may be approximated again as

$$\begin{aligned} & \frac{1}{\Delta x_j} \left\{ \left[ T_{xx(i,j+\frac{1}{2})} \frac{(h_{i,j+1,k} - h_{i,j,k})}{\Delta x_{j+\frac{1}{2}}} \right] - \left[ T_{xx(i,j-\frac{1}{2})} \frac{(h_{i,j,k} - h_{i,j-1,k})}{\Delta x_{j-\frac{1}{2}}} \right] \right\} \\ & + \frac{1}{\Delta y_i} \left\{ \left[ T_{yy(i+\frac{1}{2},j)} \frac{(h_{i+1,j,k} - h_{i,j,k})}{\Delta y_{i+\frac{1}{2}}} \right] - \left[ T_{yy(i-\frac{1}{2},j)} \frac{(h_{i,j,k} - h_{i-1,j,k})}{\Delta y_{i-\frac{1}{2}}} \right] \right\} = \frac{S_{i,j}}{\Delta t} (h_{i,j,k} - h_{i,j,k-1}) + W_{i,j,k} \end{aligned} \quad (5)$$

in which

$T_{xx(i,j+\frac{1}{2})}$  is the transmissivity between node  $(i,j)$  and node  $(i,j+1)$ ;

$\Delta x_{j+\frac{1}{2}}$  is the distance between node  $(i,j)$  and node  $(i,j+1)$ .

Equation 5 is written implicitly, that is, the head values on the left-hand side are at the new ( $k$ ) time level. Following a convention similar to that introduced by Stone (1968), the notation in equation 5 may be simplified by writing

$$\begin{aligned} & F_{i,j} (h_{i,j+1,k} - h_{i,j,k}) - D_{i,j} (h_{i,j,k} - h_{i,j-1,k}) \\ & + H_{i,j} (h_{i+1,j,k} - h_{i,j,k}) - B_{i,j} (h_{i,j,k} - h_{i-1,j,k}) \\ & = \frac{S_{i,j}}{\Delta t} (h_{i,j,k} - h_{i,j,k-1}) + W_{i,j,k} \end{aligned} \quad (6)$$

in which

$$B_{i,j} = \frac{\left[ \frac{2T_{yy[i,j]} T_{yy[i-1,j]}}{T_{yy[i,j]} \Delta y_{i-1} + T_{yy[i-1,j]} \Delta y_i} \right]}{\Delta y_i} \quad (7a)$$

The term in brackets is the harmonic mean of

$$\frac{T_{yy[i,j]}}{\Delta y_i}, \frac{T_{yy[i-1,j]}}{\Delta y_{i-1}}$$

It represents the ratio  $T_{yy(i-\frac{1}{2})}/\Delta y_{i-\frac{1}{2}}$  in equation 5.

Similarly,

$$D_{i,j} = \frac{\left[ \frac{2T_{xx[i,j]} T_{xx[i,j-1]}}{T_{xx[i,j]} \Delta x_{j-1} + T_{xx[i,j-1]} \Delta x_j} \right]}{\Delta x_j}; \quad (7b)$$

$$F_{i,j} = \frac{\left[ \frac{2T_{xx[i,j]} T_{xx[i,j+1]}}{T_{xx[i,j]} \Delta x_{j+1} + T_{xx[i,j+1]} \Delta x_j} \right]}{\Delta x_j}; \quad (7c)$$

$$H_{i,j} = \frac{\left[ \frac{2T_{yy[i+1,j]} T_{yy[i,j]}}{T_{yy[i,j]} \Delta y_{i+1} + T_{yy[i+1,j]} \Delta y_i} \right]}{\Delta y_i}. \quad (7d)$$

Use of the harmonic mean (1) insures continuity across cell boundaries at steady state if a variable grid is used, and (2) makes the appropriate coefficients zero at no-flow boundaries.

Equation 6 is also used to approximate equation 3 by replacing  $S$  with  $Sy$  and defining the transmissivities in equations 7a through 7d as a function of the head from the preceding iteration. As an example,

$$T_{xx(i,j)}^n = K_{xx(i,j)} b_{i,j,k}^{n-1}$$

in which  $n$  is the iteration index.

The notation may be simplified further by omitting subscripts not including a "+1" or "-1" (except where necessary for clarity) and by following the convention that unknown terms are placed on the left-hand side



of the equations. Equation 6 may be rearranged and expressed as

$$Bh_{i-1} + Dh_{j-1} + Eh + Fh_{j+1} + Hh_{i+1} = Q \quad (8)$$

in which

$$E = - \left( B + D + F + H + \frac{S}{\Delta t} \right);$$

$$Q = - \frac{S}{\Delta t} h_{k-1} + W.$$

### Source term

The source term  $W(x, y, t)$  can include well discharge, transient leakage from a confining bed, recharge from precipitation and evapotranspiration. In the model the source term is computed as

$$q'_{i,j,k} \cong (h_{i,j,0} - h_{i,j,k}) \left( \frac{K'_{i,j}}{\pi K'_{i,j} t} \right)^{1/2} \cdot \left\{ 1 + 2 \sum_{n=1}^{\infty} \exp \left[ \frac{-n^2}{\left( \frac{3m_{i,j}^2 S_{s[i,j]}}{K'_{i,j} t} \right)} \right] \right\} + \frac{K'_{i,j}}{m_{i,j}} (\hat{h}_{i,j,0} - h_{i,j,0}) \quad (9)$$

in which

$h_{i,j,0}$

is the hydraulic head in the aquifer at the start of the pumping period ( $L$ );

$\hat{h}_{i,j,0}$

is the hydraulic head on the other side of the confining bed ( $L$ );

$K'_{i,j}$

is the hydraulic conductivity of the confining bed ( $L/t$ );

$m_{i,j}$

is the thickness of the confining bed ( $L$ );

$S_{s[i,j]}$

is the specific storage in the confining layer ( $L^{-1}$ );

$(K'_{i,j} t / m_{i,j}^2 S_{s[i,j]})$

is dimensionless time; see Bredehoeft and Pinder (1970) for a discussion of leakage versus dimensionless time;

$t$

is the elapsed time of the pumping period ( $t$ ).

$$W_{j,j,k} = \frac{Q_{w[i,j,k]}}{\Delta x_j \Delta y_i} - q_{ro[i,j,k]} - q'_{i,j,k} + q_{et[i,j,k]}$$

in which

$Q_{w[i,j,k]}$  is the well discharge ( $L^3 t^{-1}$ );

$q_{ro[i,j,k]}$  is the recharge flux per unit area ( $L t^{-1}$ );

$q'_{i,j,k}$  is the flux per unit area from a confining layer ( $L t^{-1}$ );

$q_{et[i,j,k]}$  is the evapotranspiration flux per unit area ( $L t^{-1}$ ).

### Leakage

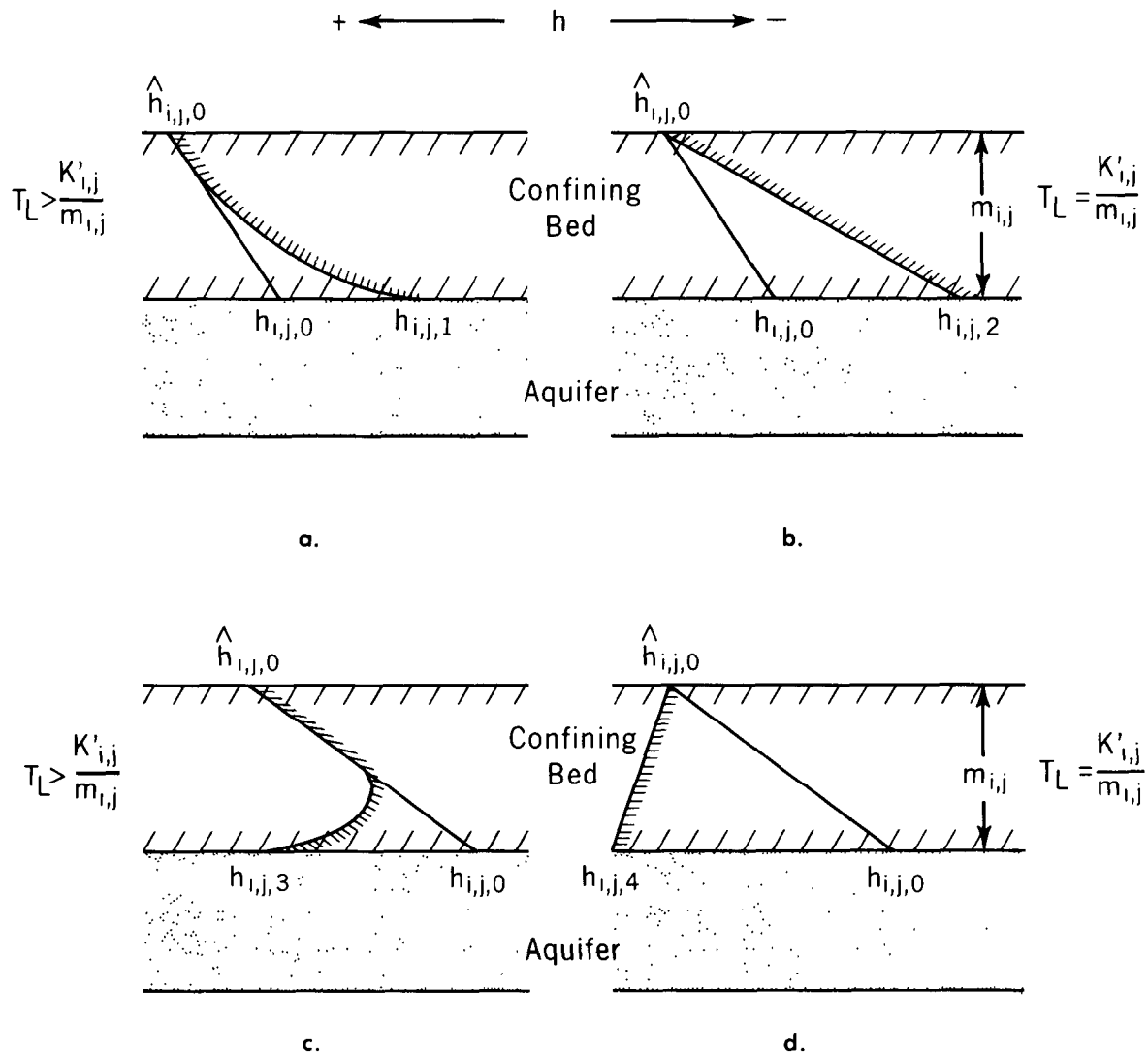
Leakage from a confining layer or streambed in which storage is considered may be approximated by

Equation 9 is modified from Bredehoeft and Pinder (1970, p. 887); note that it is the sum of two terms; the first term on the right-hand side of equation 9 considers transient effects; the second term is steady leakage due to the initial gradient across the confining bed. (See fig. 2.) Figure 2 illustrates the head distribution in the confining layer at any given point in the aquifer system at two different times in each of two successive pumping periods. (The succession of head values in the aquifer is shown by  $h_{i,j,1}, \dots, h_{i,j,4}$ .) The solid line represents the head distribution at the beginning of the pumping period; the gradient  $((\hat{h}_{i,j,0} - h_{i,j,0}) / m_{i,j})$  appears in the second term of equation 9. The hatched line represents the head distribution in the confining bed after stressing the pumped aquifer and is a summation of the initial head distribution and the change in head distribution due to the stresses on the aquifer. The factor  $T_L$  in figure 2 represents the part of the first term in equation 9 independent of head (that is, the transient leakage coefficient).

In figure 2a the confining bed is assumed to have significant storage, pumping has low-

ered the head to  $h_{i,j,1}$  and the net (or total) gradient is for some dimensionless time  $< 0.5$ . After transient effects have dissipated, a uniform gradient across the confining bed is es-

tablished. (See fig. 2b.) Then if the stress on the aquifer is changed by turning off pumping wells and starting recharge wells, the initial head distribution in the confining bed



$$q'_{i,j,k} = T_L (h_{i,j,0} - h_{i,j,k}) + \frac{K'_{i,j}}{m_{i,j}} (\hat{h}_{i,j,0} - h_{i,j,0})$$

#### EXPLANATION

- Initial head in confining bed
- /////// Head in confining bed after stressing the aquifer

FIGURE 2.—In the first pumping period, (a) illustrates the head distribution in the confining bed at one time when transient leakage effects are significant; (b) illustrates a time after transient effects have dissipated; in the second pumping period, (c) is analogous to (a) and (d) is analogous to (b).

for the new conditions is shown in figure 2c and is equal to the final distribution for the first pumping period. The net head distribution in figure 2c is affected by storage in the confining bed and is for some dimensionless time  $< 0.5$  (in the second pumping period). After storage effects have dissipated, the net gradient is shown in figure 2d.

For a simulation of several pumping periods, the program assumes that transient leakage effects from previous pumping periods have dissipated. This is accomplished at the start of each pumping period by initializing  $h_{i,j,0}$  to the head at the end of the previous pumping period and setting  $t$  (and thereby dimensionless time) to zero (note that the parameter storing the cumulative simulation time is not affected). The assumption is reasonable if dimensionless time for previous pumping periods is at least 0.5 (Bredehoeft and Pinder, 1970, fig. 4) and can be checked by noting the value of dimensionless time printed in the output for the end of the previous pumping period. If the assumption is not valid, the code will need to be modified to include transient effects for one or more previous pumping periods.

In the model, equation 9 is used until dimensionless time reaches  $3 \times 10^{-3}$ ; otherwise, the equation

$$q'_{i,j,k} \cong (h_{i,j,0} - h_{i,j,k}) \frac{K'_{i,j}}{m_{i,j}} \left\{ 1 + 2 \sum_{n=1}^{\infty} \exp \left[ -n^2 \pi^2 \left( \frac{K'_{i,j} t}{3m_{i,j}^2 S_{s[i,j]}} \right) \right] \right\} + \frac{K'_{i,j}}{m_{i,j}} (\hat{h}_{i,j,0} - h_{i,j,0}) \quad (10)$$

is used. Equation 10 is computationally more efficient for dimensionless times greater than about  $3 \times 10^{-3}$ .

The transient parts of equations 9 and 10 are based on the analytic solutions for the flux from a confining layer resulting from an instantaneous stepwise change in head in the aquifer. The factor of  $1/3$  appearing in dimensionless time is included in order to approximate the transient flux resulting from the actual drawdown in the aquifer. In effect the transient flux is approximated by applying a step change in head equal to the drawdown from the start of the pumping period at  $1/3$  of the elapsed time in the pumping period. (See fig. 3.)

The results of several numerical experiments indicate that it would be better to use

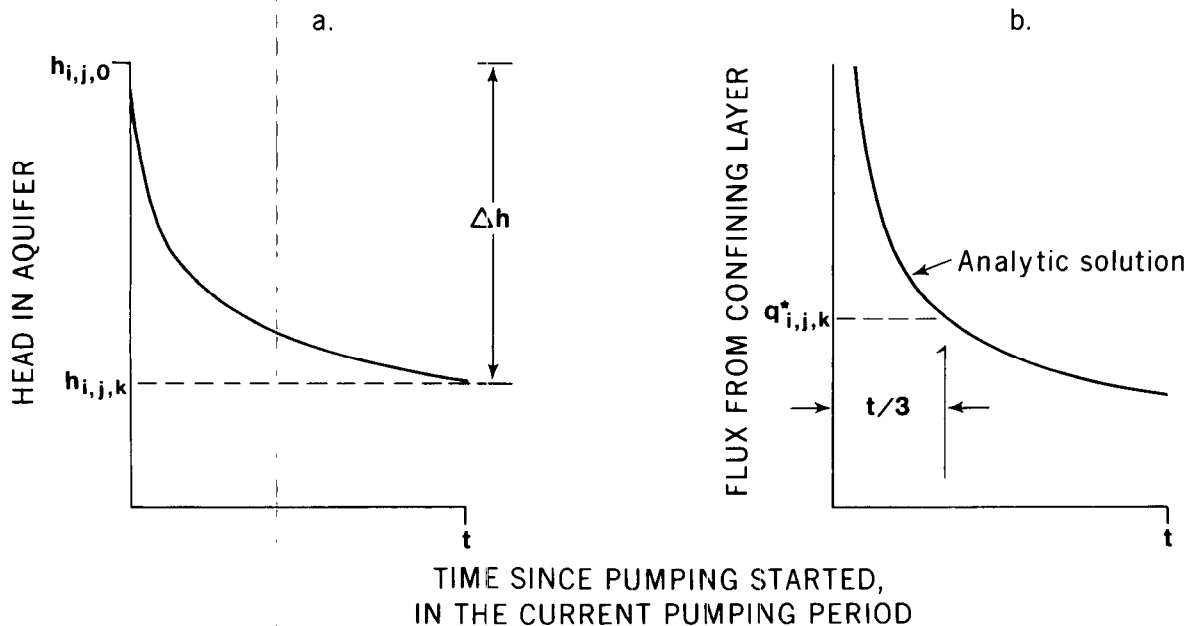


FIGURE 3.—The total drawdown at the elapsed time,  $t$ , in the pumping period (a) is applied at  $t/3$  in equations 9 and 10 to approximate  $q'_{i,j,k}$ , the transient part of  $q'_{i,j,k}$  (b).

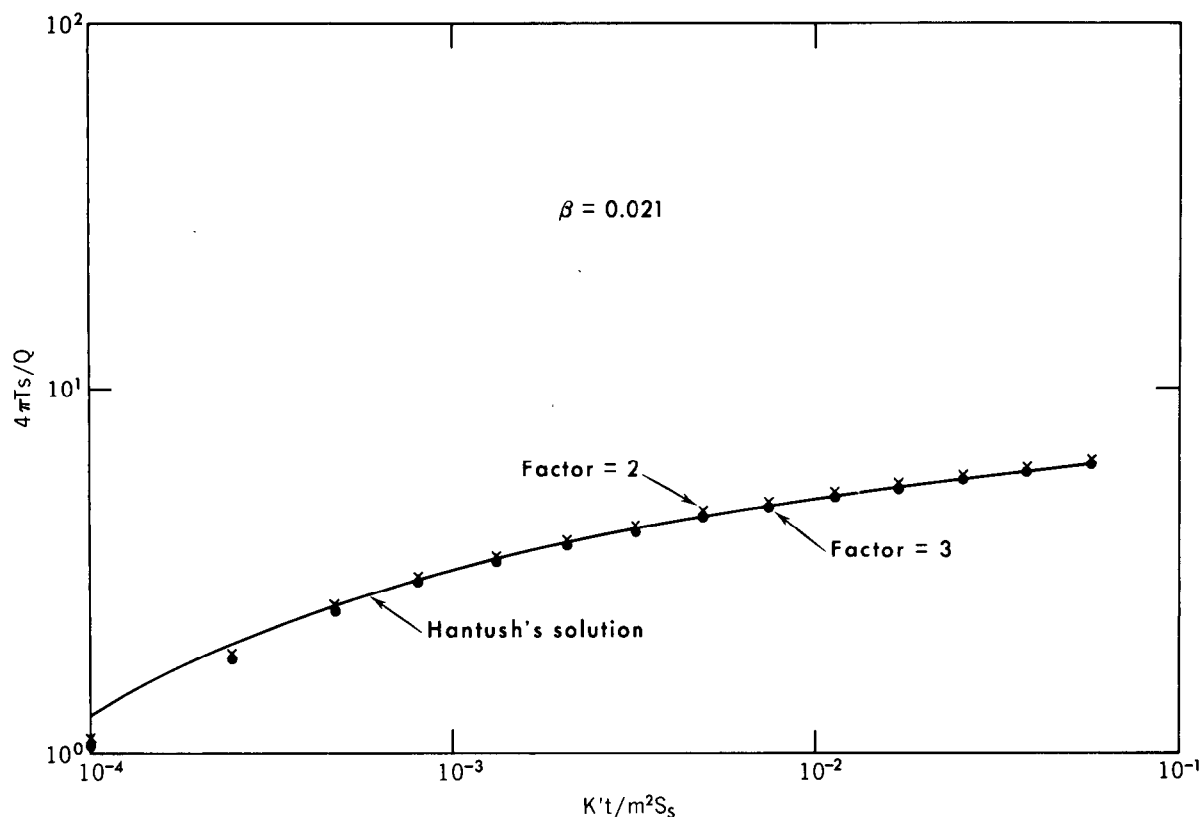


FIGURE 4.—Comparison of analytic solution and numerical results using factors of 2 and 3 in the transient leakage approximation.

a factor of 1/3 rather than the factor of 1/2 used in the approximation by Bredehoeft and Pinder (1970). In figure 4 are plotted numerical results and Hantush's (1960) analytic solution for  $\beta=0.021$  ( $\beta=0.25 r [K'S_s/TS]^{1/2}$  and  $r$  is the radial distance from the center of the pumping well). The drawdown values using a factor of 1/3 are below but very close to the analytic curve after the first few time steps. The results using a factor of 1/2 are close to the analytic solution but are about twice as far above the analytic curve as the factor of 1/3 results are below the curve. In figure 5 are plotted the percent difference between the volume of leakage

computed numerically and the volume determined analytically. Two sets of data are shown: a 14-step simulation between dimensionless times of  $10^{-5}$  and  $5.8 \times 10^{-2}$  and an 11-step simulation between dimensionless times of  $5.8 \times 10^{-3}$  and  $4.4 \times 10^{-1}$ . Based on those experiments, if 4 or 5 time steps are simulated before the period of interest, the volume of leakage and the drawdown computed numerically using a factor 1/3 in equations 9 and 10 are close to the analytic solution.

#### Evapotranspiration

Evapotranspiration as a linear function of depth below the land surface is computed as

$$q_{et[i,j,k]} = \begin{cases} Q_{et} & [h_{i,j,k} \geq G_{i,j}] \\ Q_{et} - \frac{Q_{et}}{ET_z} (G_{i,j} - h_{i,j,k}) & [ET_z > (G_{i,j} - h_{i,j,k}); h_{i,j,k} < G_{i,j}] \\ 0 & [ET_z \leq (G_{i,j} - h_{i,j,k})] \end{cases} \quad (11)$$

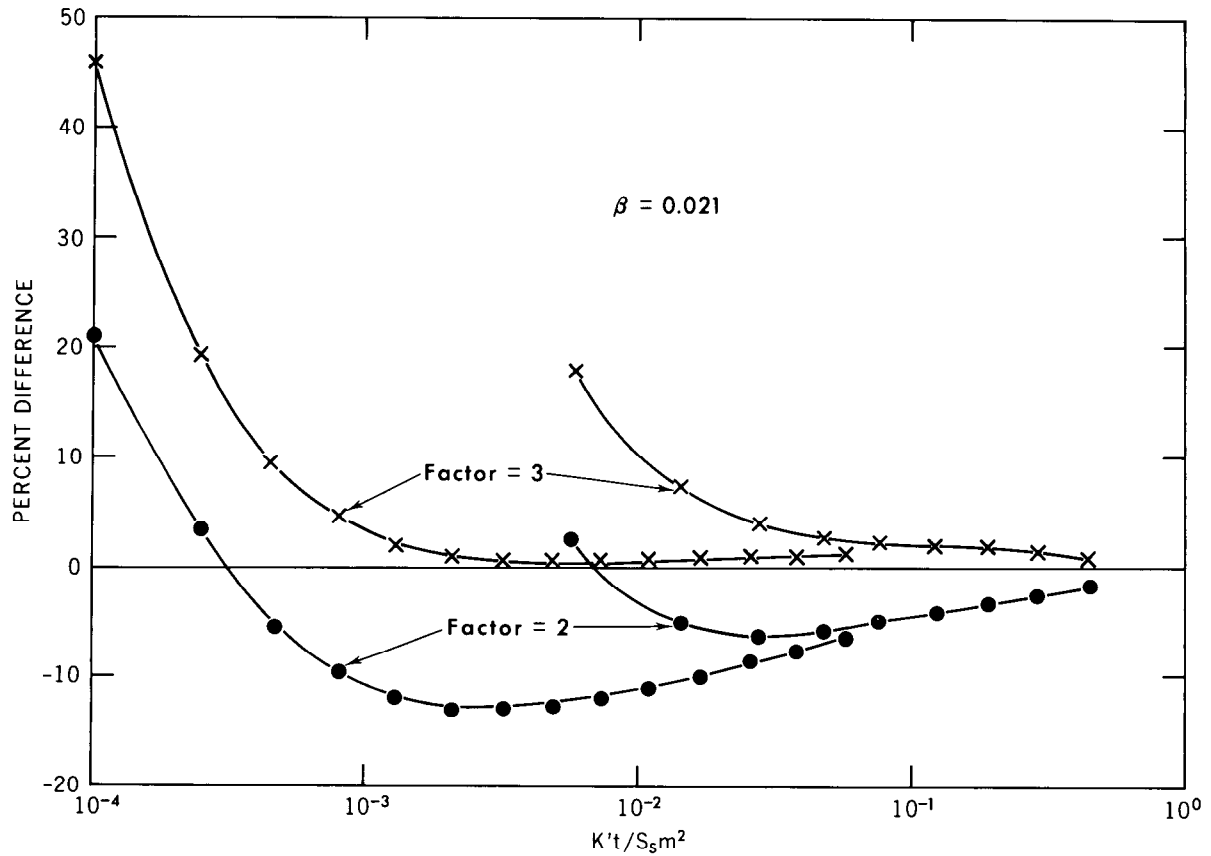


FIGURE 5.—Percent difference between the volume of leakage computed with the model approximation and Hantush's analytical results.

in which

$Q_{et}$  is the maximum evapotranspiration rate ( $Lt^{-1}$ );

$ET_z$  is the depth below land surface at which evapotranspiration ceases ( $L$ );

$G_{i,j}$  is the elevation of the land surface ( $L$ ).

This relationship (illustrated in fig. 6) is treated implicitly by separating the equation into two terms<sup>1</sup>: one term is included with the  $E$  coefficient on the left-hand side of equation 8; the other is a known term included in  $Q$  on the right-hand side of equation 8.

Other functions for evapotranspiration can be defined (for example, decreasing ex-

<sup>1</sup> Some of the methods for implicit treatment of evapotranspiration, storage, and leakage have been adapted from Prickett and Lonquist (1971).

ponentially with depth), but it may be more difficult to treat these relationships numerically. The easiest approach is to make evapotranspiration an explicit function of the head at the previous iteration, but this may cause oscillations and difficulties with convergence. Normally, the oscillations may be dampened by making evapotranspiration a function of the head for the two previous iterations. A more sophisticated approach is to use the Newton-Raphson method, which is a rapidly converging iterative technique for treating systems of non-linear equations. (See, for example, Carnahan, Luther, and Wilkes, 1969, p. 319–329.)

### Computation of head at the radius of a pumping well

The hydraulic head computed for a well node represents an average hydraulic head

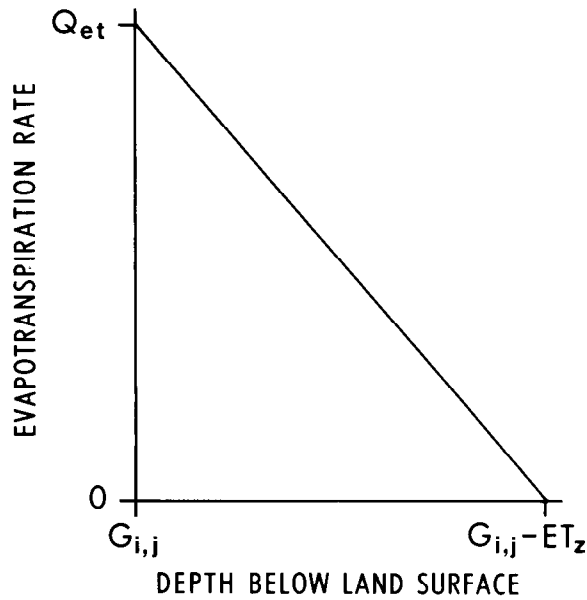


FIGURE 6.—Evapotranspiration decreases linearly from  $Q_{et}$  where the water table is at land surface to zero where the water table is less than or equal to  $G_{i,j} - ET_z$ .

computed for the block and is not the head in a well. An option to compute the head and drawdown at a well is included in the model. This computation uses the radius,  $r_e$ , of a hypothetical well for which the average value of head for the cell applies. An approximating equation is then used to make the extrapolation from  $r_e$  to the radius of a real well.

The radius  $r_e$  can be computed as (Prickett, 1967)

$$r_e = r_1 / 4.81 \quad (12)$$

in which  $r_1 = \Delta x_j = \Delta y_i$  (fig. 7). Equation 12 assumes steady flow, no source term other than well discharge in the well block, and that the area around the well is isotropic and homogeneous. The derivation of equation 12 can be seen with reference to figure 7 in which the four nodes adjacent to node  $i, j$  are assumed to have head values equal to the value at node  $i-1, j$ . In figure 7a one-quarter of the discharge to the well node  $i, j$  is computed by the model as

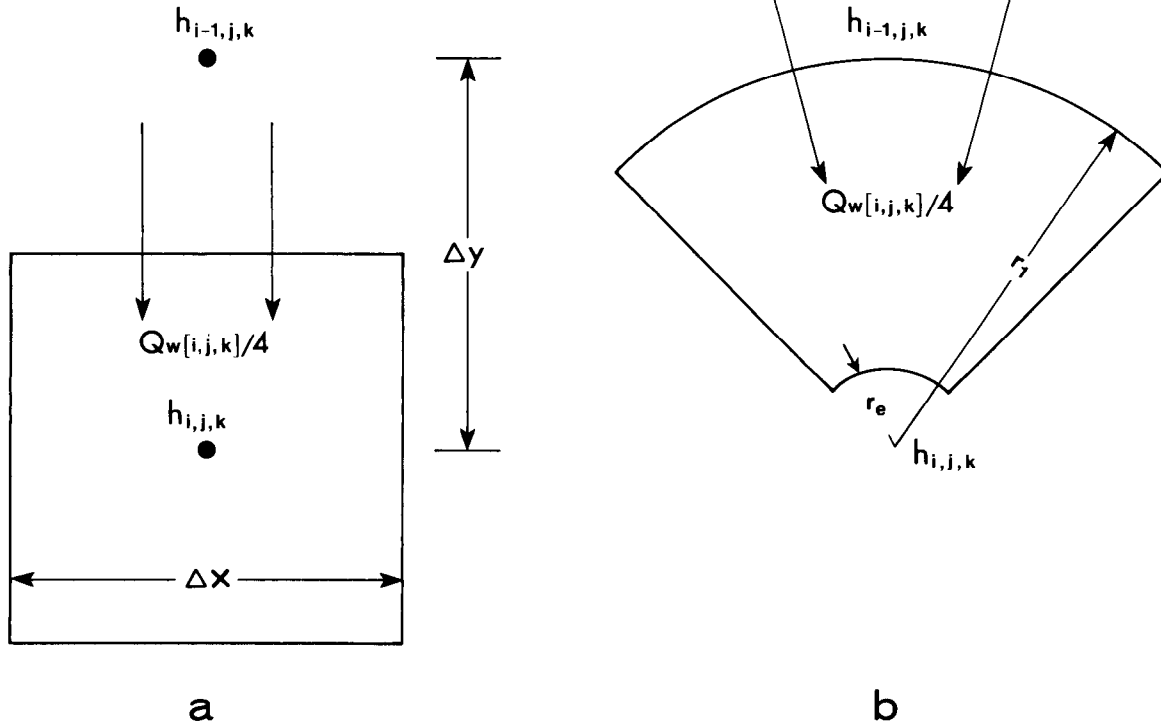


FIGURE 7.—Flow from cell  $(i-1, j)$  to cell  $(i, j)$  (a) and equivalent radial flow to well  $(i, j)$  with radius  $r_e$  (b).

$$\frac{Q_{w[i,j,k]}}{4} = \Delta x_j T_{i,j} \frac{\Delta h}{\Delta y} \quad (13)$$

in which

$$\Delta h = h_{i-1,j,k} - h_{i,j,k};$$

$$T_{i,j} = T_{xx[i,j]} = T_{yy[i,j]}.$$

The equivalent discharge for radial flow to the well is given by the Thiem (1906) equation expressed as (see fig. 7b)

$$\frac{Q_{w[i,j,k]}}{4} = \frac{\pi T_{i,j}}{2} \frac{\Delta h}{\ln(r_1/r_e)}. \quad (14)$$

Equating the discharges in equations 13 and 14 gives equation 12.

The Thiem equation is commonly used to extrapolate from the average hydraulic head for the cell at radius  $r_e$  to the head,  $h_w$ , at the desired well radius,  $r_w$  (Prickett and Lonquist, 1971; Akbar, Arnold, and Harvey, 1974) and is written in the form

$$h_w = h_{i,j,k} - \frac{Q_{w[i,j,k]}}{2\pi T_{i,j}} \ln(r_e/r_w). \quad (15)$$

Equation 15 assumes that: (1) flow is within a square well block and can be described by a steady-state equation with no source term except for the well discharge, (2) the aquifer is isotropic and homogeneous in the well block, (3) only one well is in the block and it fully penetrates the aquifer, (4) flow is laminar, and (5) well loss is negligible.

In an unconfined aquifer, the analogous equation is

$$H_w = \sqrt{H_{i,j,k}^2 - \frac{Q_{w[i,j,k]}}{\pi K_{i,j}} \ln(r_e/r_w)} \quad (16)$$

in which

$H_{i,j,k} = h_{i,j,k}$  - BOTTOM (I,J) is the saturated thickness of the aquifer at radius  $r_e$  (L);

$H_w$  is the saturated thickness of the aquifer at the well (L);

$K_{i,j} = K_{xx[i,j]} = K_{yy[i,j]}$ ;

BOTTOM (I,J) = elevation of the bottom of the aquifer (The uppercase let-

ters indicate that this parameter is identical to that used in the model.)

When the saturated thickness computed with equation 16 is negative, the message, 'X,Y WELL IS DRY' is generated. This situation has no effect on the computations, but should stimulate careful consideration of the value of results for subsequent time steps in the simulation.

The conditions when the Thiem equation or equation 16 will be accurate can be computed. Table 1 was prepared to give a few examples of the head values computed by the model with the Thiem equation for a well with a radius of 1.25 feet in an infinite leaky artesian aquifer and in an infinite nonleaky artesian aquifer. The analytic solutions for these conditions are included for comparison. A variable grid was used in the model but the dimensions of the well block were  $\Delta x = \Delta y = 1,000$  feet. For conditions which depart significantly from the assumptions given above (for example, a well in a rectangular block with anisotropic transmissivity or a well in a large block that has a significant amount of leakage) the results using equations 15 and 16 should be checked with a more rigorous analysis. Additional drawdown due to the effects of partial penetration and well loss can be computed separately or added to the code as needed.

Table 1.—Comparison of drawdowns computed with equation 15 and the analytic values

Aquifer	Time step	Dimensionless time	Drawdown	
			Approximation	Analytic
Nonleaky artesian -----	3	$Tt/r^2S$		
	14	$3.0 \times 10^5$ $3.7 \times 10^7$	41.1	42.7
Leaky artesian -----	3	$K't/m^2S_e$	58.3	58.1
	9	0.028 .44	51.8	52.1
			57.1	57.3

### Combined artesian—water-table simulation

Simulation of an aquifer that is partly confined and elsewhere has a free surface requires special computations for the transmissivity, storage coefficient, and leakage

term. The following paragraphs describe the computations required. Some of the methods of coding these procedures have been adapted from Prickett and Lonquist (1971).

### Transmissivity

The transmissivity is computed as the saturated thickness of the aquifer times the hydraulic conductivity. This computation requires that the elevations of the top and bottom of the aquifer be specified. Where the aquifer crops out, the top of the aquifer is assigned a fictitious value greater than or equal to the elevation of the land surface.

### Storage

The storage term requires special treatment at nodes where a conversion from artesian to water-table conditions, or vice versa, occurs during a time step. The program first checks for a change at a node during the last iteration. If there has been a change from artesian to water-table conditions, the storage term is

$$\frac{S_{y[i,j]}}{\Delta t} (h_{i,j,k}^n - h_{i,j,k-1}) - \text{SUBS}$$

in which

$$\text{SUBS} = (h_{i,j,k-1} - \text{TOP}(I,J)) (S_{y[i,j]} - S_{y[i,j]}) / \Delta t;$$

$\text{TOP}(I,J)$  = elevation of the top of the aquifer.

The purpose of SUBS is to correctly apportion the storage coefficient and specific yield according to the relationship in figure 8a.

For a change from water-table to artesian conditions, the storage term is

$$\frac{S_{i,j}}{\Delta t} (h_{i,j,k}^n - h_{i,j,k-1}) - \text{SUBS}$$

in which

$$\text{SUBS} = (h_{i,j,k-1} - \text{TOP}(I,J)) (S_{y[i,j]} - S_{i,j}) / \Delta t.$$

SUBS subtracts the storage coefficient and adds the specific yield for the distance  $B$  illustrated in figure 8b.

### Leakage

To treat leakage more realistically if parts of an artesian aquifer change to water-table conditions, the maximum head difference across the confining bed is limited to  $\hat{h}_{i,j,0} - \text{TOP}(I,J)$ .

Two examples illustrate the calculation of leakage in conversion simulations. In figure 9a the head at the start of the pumping period,  $h_{i,j,0}$  is below the water-table head,  $\hat{h}_{i,j,0}$ , but above the top of the aquifer; the current pumping level is below the top of the aquifer. The applicable equation is

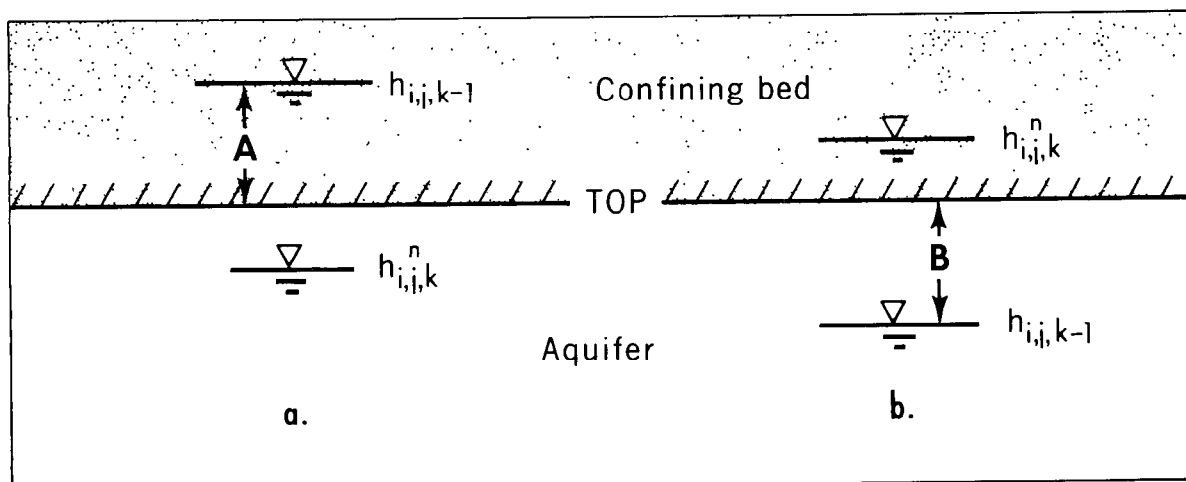


FIGURE 8.—Storage adjustment is applied to distance A in conversion from artesian to water-table conditions (a) and to distance B in conversion from water-table to artesian conditions (b).



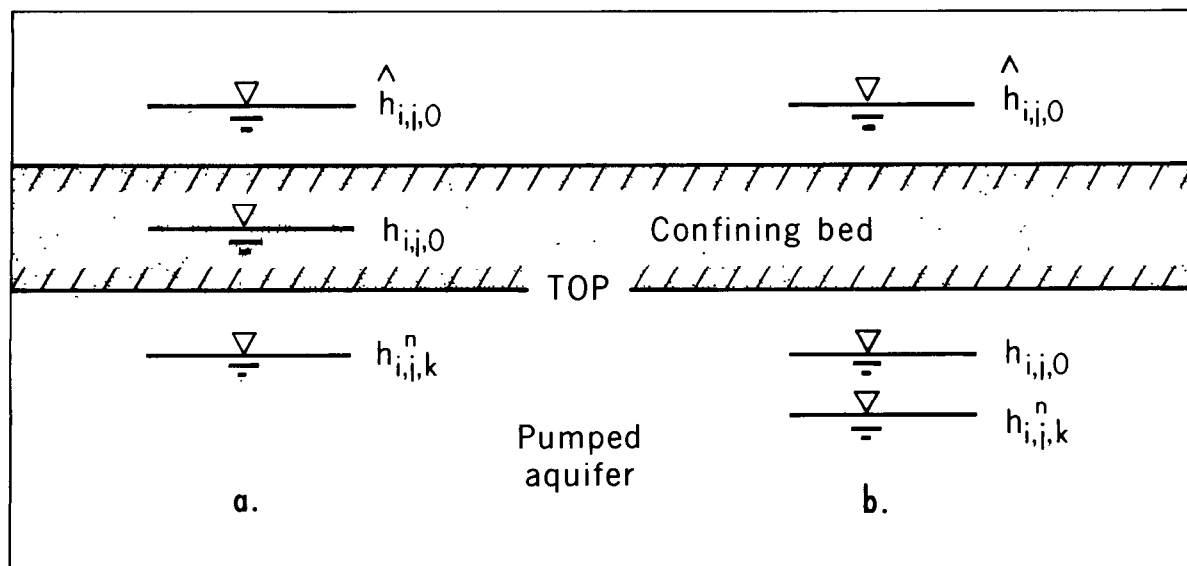


FIGURE 9.—Two of the possible situations in which leakage is restricted in artesian-water-table simulations.

$$q'_{i,j,k} = \frac{K'_{i,j}}{m_{i,j}} (\hat{h}_{i,j,0} - h_{i,j,0}) + T_L (h_{i,j,0} - \text{TOP}(I,J)).$$

For this situation  $q'_{i,j,k}$  appears on the right-hand side of the difference equation and is treated explicitly. Only if both  $h_{i,j,0}$  and  $h_{i,j,k}^n$  are above the top of the aquifer is the leakage term treated implicitly by including  $T_L$  in the  $E$  coefficient. This is accomplished in the code by setting  $U=1$ .

In the second example (fig. 9b), both  $h_{i,j,0}$  and  $h_{i,j,k}^n$  are below the top of the aquifer and the equation for leakage reduces to

$$q'_{i,j,k} = \frac{K'_{i,j}}{m_{i,j}} (\hat{h}_{i,j,0} - \text{TOP}(I,J)).$$

If leakage across a subjacent confining bed is significant, it will be necessary to add a second leakage term. The flux described by this term will not be restricted where water-table conditions occur.

## Test Problems

In a subsequent section the computational work required for solution of four test problems by the numerical techniques available in the model is analyzed. It is appropriate, however, to introduce the test problems here because they are used in the discussion of iteration parameters in the section on numerical

techniques. The problems are for steady-state conditions since the resulting set of simultaneous equations are more difficult to solve than are the set of equations for transient problems which generally involve smaller head changes.

For each of these problems a closure criterion was chosen to decide when a solution is obtained to the set of finite-difference equations. (See Remson, Hornberger, and Molz, 1971, p. 185–186.) Normally, in this model, a solution is assumed if:

$$\text{Max} | h^n - h^{n-1} | \leq \epsilon$$

where  $\epsilon$  is an arbitrary closure criterion ( $L$ ). For the purpose of the numerical comparisons given later in this documentation, the absolute value of the maximum residual (defined by equation 28) is used to compare methods.

The first problem is a square aquifer with uniform properties and grid spacing (fig. 10). The finite-different grid is  $20 \times 20$ , but only 18 rows and columns are inside the aquifer because the model requires that the first and last rows and columns be outside the aquifer boundaries. Two discharging wells and one recharge well are the stress on the system; boundaries are no flux except for part of one side which is a constant-head

## PROBLEM CHARACTERISTICS

Transmissivity:  $T_{xx} = T_{yy} = 0.1 \text{ ft}^2/\text{s}$  ( $0.009 \text{ m}^2/\text{s}$ )  
 Grid spacing:  $\Delta x = \Delta y = 5000 \text{ ft}$  ( $1500 \text{ m}$ )  
 Dimensions of grid.  $18 \times 18$

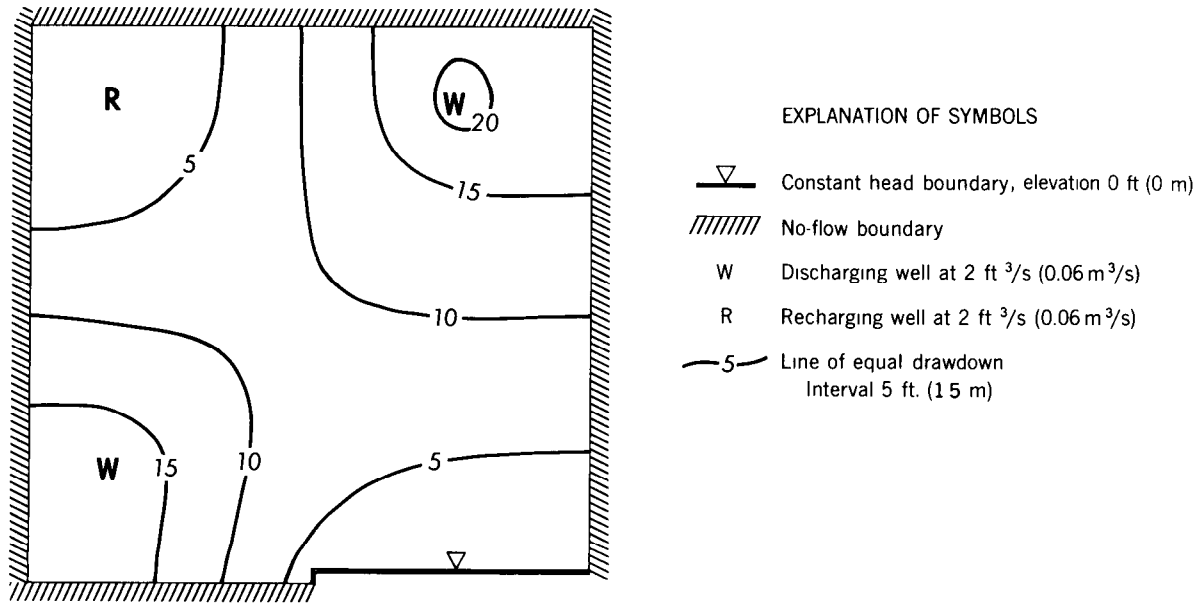


FIGURE 10.—Characteristics of test problem 1.

boundary. A closure criterion of 0.001 foot (0.0003 metre) was used.

Konikow (1974) designed the second problem in his analysis of ground-water pollution at the Rocky Mountain Arsenal northeast of Denver, Colo. It is included as one of the test problems because it is typical of many field problems and because there is some difficulty in obtaining a steady-state solution with the alternating-direction implicit procedure. The transmissivity distribution is shown in figure 11; note the extensive areas where the transmissivity is zero because the surficial deposits are unsaturated. The finite-difference grid representing this aquifer is  $25 \times 38$  with square blocks 1,000 feet (300 metres) on a side. The model has constant-head boundaries at the South Platte River and where the aquifer extends beyond the limits of the model; elsewhere no-flux boundaries are employed. Although this is a water-table aquifer, it is assumed for problem 2 that

transmissivity is independent of head. The model includes 49 irrigation wells and recharge from canals and irrigation. In figure 11 the observed water-table configuration is shown, and it is used as the initial surface for the simulation; the computed water table is generally within a few feet of the observed. For this problem the closure criterion is 0.001 foot (0.0003 metre).

The third problem is a cross-section with three horizontal layers and other characteristics shown in figure 12. Transmissivity equals hydraulic conductivity for this problem because it is conceived as a slice one unit wide. The values for transmissivity are arbitrary. Note in particular that the horizontal conductivity is 100 times the vertical conductivity in all layers and that the middle layer acts as a confining layer between the upper and lower layers. The coefficients  $B_{i,j}$  and  $H_{i,j}$ , however, are 100 times greater than the horizontal coefficients  $D_{i,j}$  and  $F_{i,j}$  because of

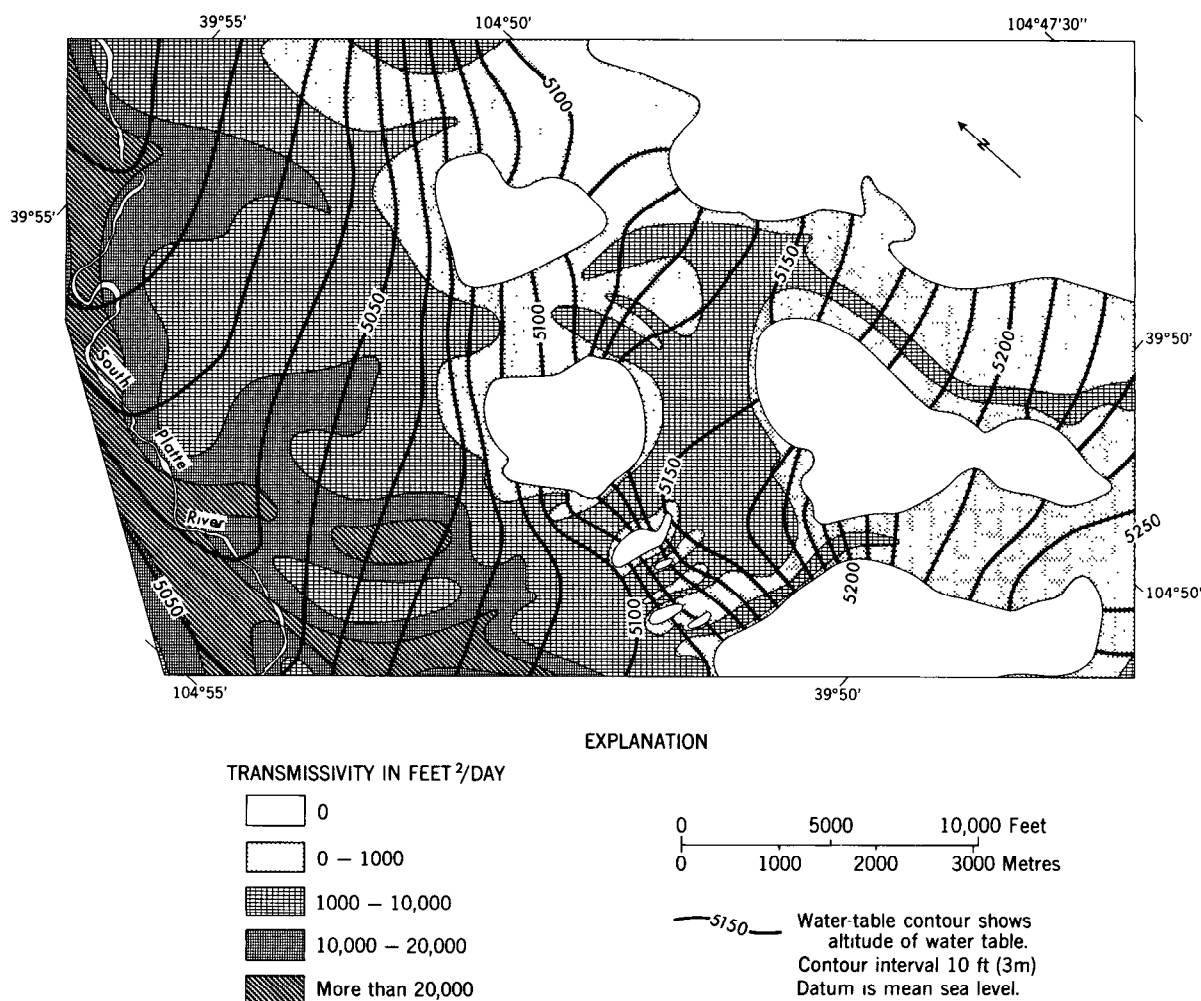


FIGURE 11.—Transmissivity and observed water-table configuration for test problem 2 (fieldwork and model design by Konikow, 1975).

the grid spacing used. For this problem, the closure criterion is 0.0001 feet (0.00003 metre).

In the third problem the upper boundary (the water table) is fixed as a constant-head boundary. It could also be treated as a no-flow boundary which would effectively confine the system. This model was not designed specifically for simulation of cross sections, and consequently it does not have provision for a moving boundary. Rather than modifying this one-phase model for a moving-boundary problem, it would be better to design a model specifically for this purpose. The two-phase model described by Freeze (1971) is a good example.

The fourth problem is to consider the water-table case of the second problem. The only difference from problem 2 is that transmissivity is dependent upon (1) head in the aquifer, (2) aquifer base elevation, and (3) hydraulic conductivity of the aquifer.

## Numerical Solution

In Pinder (1969) and Trescott (1973) the iterative, alternating-direction implicit procedure (ADI) was the only option available for numerical solution. For many field problems ADI is convergent and competitive, in terms of the computational work required,

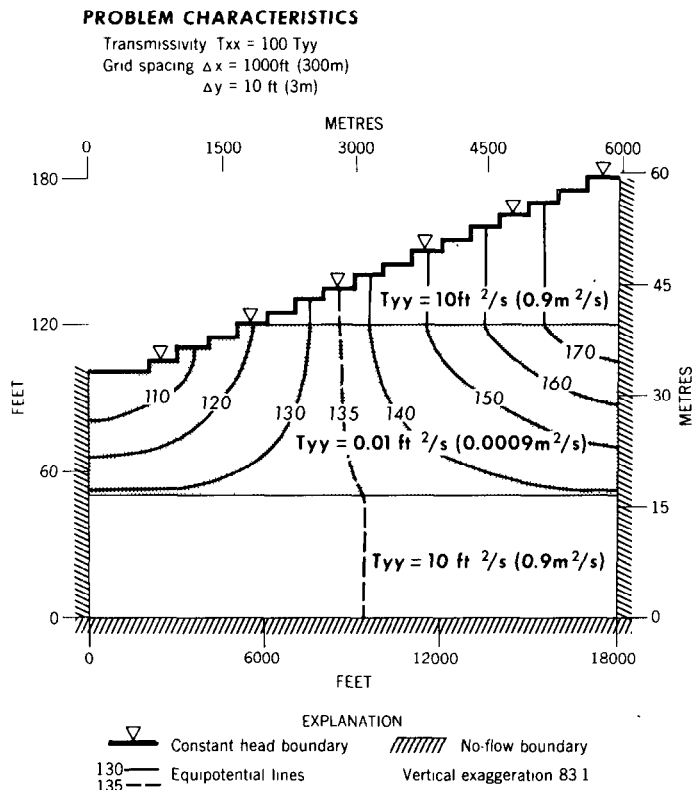


FIGURE 12.—Characteristics of test problem 3.

with other iterative techniques available. It may be difficult, however, to obtain a solution for some problems with ADI (for example, steady-state simulations involving extremely variable coefficients). Consequently, it is convenient to have available other numerical techniques that may be more suited than ADI to particular problems. The three numerical methods available with this model are ADI, the strongly implicit procedure (SIP), and line successive overrelaxation (LSOR).

The following sections outline the computational algorithms for the three numerical methods. More details are given in the discussion on SIP, because that method is more complex.

For additional details on the theory behind the methods and rigorous analysis of convergence rates, see for example, Varga (1962) and Remson, Hornberger, and Molz (1971). The methods are presented in order of increasing complexity. In general, the more complex methods converge more rapidly and are applicable to more types of problems than the simpler methods such as LSOR. For clari-

ty, the numerical treatment of the source term is left to other sections.

### Line successive overrelaxation

Line successive overrelaxation (LSOR) improves head values one row (or column) at a time. Whether the solution is oriented along rows or columns is generally immaterial for isotropic problems but has a significant affect on the convergence rate in anisotropic problems. The solution should be oriented in the direction of the larger coefficients, either  $B_{i,j}$  and  $H_{i,j}$  or  $D_{i,j}$  and  $F_{i,j}$  (Breitenbach, Thurnau, and van Poolen, 1969, p. 159). Differences in the magnitude of the coefficients may result from anisotropic transmissivity or from a large difference in grid spacing between the  $x$  and  $y$  directions. In problem 3 the largest transmissivity is in the horizontal direction in each layer, but the small grid spacing in the vertical direction makes the coefficients  $B_{i,j}$  and  $H_{i,j} \gg D_{i,j}$  and  $F_{i,j}$ .

With the solution oriented along rows, an



The values of  $\bar{\xi}^\dagger$  for row  $i$  are then computed by backward substitution as

$$\xi_{i,j,k}^\dagger = G_j - BE_j \xi_{i,j+1,k}^\dagger$$

where

$$\xi_{i,N_x,k}^\dagger = G_{N_x}$$

since

$$BE_{N_x} = 0.$$

The head values for row 1 are then computed by the equation

$$h_{i,j}^n = h_{i,j}^{n-1} + \omega \xi_{i,j}^\dagger, \quad j = 1, \dots, N_x$$

If  $\omega$  is 1, the solution is by the line Gauss-Seidel formula, but convergence is slow in general. The convergence rate is improved significantly by "overrelaxation" with  $1 < \omega < 2$ . Discussion of the acceleration parameter is deferred until after the following section on two-dimensional correction.

#### Two-dimensional correction to LSOR

In certain problems, the rate of convergence of LSOR can be improved by applying a one-dimensional correction (1DC) procedure introduced by Watts (1971) or the extended two-dimensional correction (2DC) method described by Aziz and Settari (1972). These methods remove the components of certain eigenvectors in the LSOR iteration matrix from the solution vector. If the eigenvalues associated with these eigenvectors dominate the problem, particularly those including anisotropy, the convergence rate is greatly improved.

The 2DC method is applied after one or more LSOR iterations. The corrected head values are used as an improved starting point for the next iteration and the process is repeated until convergence is achieved.

The two-dimensional correction for the head at  $(i,j)$  is defined as

$$h_{i,j,k}^{n*} = h_{i,j,k}^n + \alpha_i + \hat{\beta}_j, \quad \begin{matrix} i = 1, \dots, N_y \\ j = 1, \dots, N_x \end{matrix}$$

in which

- $h_{i,j,k}^{n*}$  is the corrected head at iteration  $n$ ;
- $\alpha_i$  is the correction for row  $i$ ;
- $\hat{\beta}_j$  is the correction for column  $j$ .
- $N_y$  is the number of nodes in a column.

An approximate equation for  $\bar{\alpha}$  is

$$B'_i \alpha_{i-1} + E'_i \alpha_i + H'_i \alpha_{i+1} = R'_i, \quad i = 1, 2, \dots, N_y \quad (18)$$

in which

$$B'_i = -\sum_j B_{i,j};$$

$$E'_i = \sum_j (B_{i,j} + H_{i,j} + \frac{S_{i,j}}{\Delta t});$$

$$H'_i = -\sum_j H_{i,j};$$

$$R'_i = \sum_j R_{i,j}^n;$$

$$R_{i,j}^n = B_{i,j} h_{i-1,j,k}^n + D_{i,j} h_{i,j-1,k}^n + E_{i,j} h_{i,j,k}^n + F_{i,j} h_{i,j+1,k}^n + H_{i,j} h_{i+1,j,k}^n + \frac{S_{i,j}}{\Delta t} h_{i,j,k-1}^n - W_{i,j,k};$$

An approximate equation for  $\bar{\beta}$  is

$$D'_j \hat{\beta}_{j-1} + E'_j \hat{\beta}_j + F'_j \hat{\beta}_{j+1} = R'_j, \quad j = 1, 2, \dots, N_x \quad (19)$$

in which

$$D'_j = -\sum_i D_{i,j};$$

$$E'_j = \sum_i (D_{i,j} + F_{i,j} + \frac{S_{i,j}}{\Delta t});$$

$$F'_j = -\sum_i F_{i,j};$$

$$R'_j = \sum_i R_{i,j}^n$$

Equations 18 and 19 are derived with the following equations

$$\sum_{j=1}^{N_x} R_{j,i}^{n*} = 0, \quad i = 1, 2, \dots, N_y$$

and

$$\sum_{i=1}^{N_y} R_{i,j}^{n*} = 0, \quad j = 1, 2, \dots, N_x$$

which force the sum of residuals for each row and each column to zero when the vector  $\bar{h}^{n*}$  is substituted into equation 8. Aziz and Settari (1972) give the exact equations for  $\bar{\alpha}$  and  $\bar{\beta}$  but point out that equations 18 and 19 are good approximations and, in practice, are easier to solve. For example, equation 19, which used alone is Watts' 1DC method, is written in matrix form as

$$\begin{bmatrix} E'_1 & F'_1 \\ D'_2 & E'_2 & F'_2 \\ D'_3 & E'_3 \end{bmatrix} \begin{bmatrix} \hat{\beta}_1 \\ \hat{\beta}_2 \\ \hat{\beta}_3 \end{bmatrix} = \begin{bmatrix} R'_1 \\ R'_2 \\ R'_3 \end{bmatrix}$$

for the problem in figure 13. Equation 18 has an analogous form and both are easily solved by the Thomas algorithm.

Note that  $\bar{\alpha}$  and  $\bar{\beta}$  in the model are zero for those rows and columns in which one or more constant-head nodes are located. If  $\bar{\alpha}$  and  $\bar{\beta}$  were not zero it would not be possible to maintain a constant value at the appropriate nodes. As Watts (1973) points out, therefore, the procedure is most useful in simulations dominated by no-flow boundaries. For those simulations in which 2DC is useful, it is generally better to apply the corrections after several rather than after each LSOR iteration. After experimenting with a few problems, we have found it practical to apply 2DC after every 5 LSOR iterations.

#### LSOR acceleration parameter

The optimum value of  $\omega$  for maximum rate of convergence lies between 1 and 2 and is commonly between 1.6 and 1.9. If only one or two runs will be made on a problem, it is probably best to choose an  $\omega$  based on experience. If many runs will be made, it will be worthwhile to use an  $\omega$  close to the optimum value. For simple problems  $\omega_{opt}$  can be computed as explained, for example, by Remson, Hornberger, and Molz (1971, p. 188-199) using the equation

$$\omega = \frac{2}{1 + \sqrt{1 - \rho(G)}} \quad (20)$$

in which

$$\rho(G) \cong \left| \frac{\xi_{\max}^{\dagger(n)}}{\xi_{\max}^{\dagger(n-1)}} \right|$$

$\rho(G)$  is the spectral radius (dominant eigenvalue) of the Gauss-Seidel iteration matrix. For typical field problems it is possible to use equation 20 to estimate  $\omega_{opt}$  in an iterative process if 2DC is not used. In the first simulation of the problem, set  $\omega = 1.0$  and allow at least 100 iterations. In applying this method

to problems 1, 2, and 3 it took 25 iterations to arrive at  $\omega_{opt}$  for problem 2, but about 100 iterations to obtain  $\omega_{opt}$  for problem 1 and 3. Obviously this method may involve a lot of computational effort to obtain  $\omega_{opt}$ . More efficient methods using equation 20 have been devised to update  $\omega$  during the iteration process. For example, Breitenbach, Thurnau, and van Poolen (1969) use a modified form of Varga's (1962) "power method," Carré's (1961) method is described by Remson, Hornberger, and Molz (1971, p. 199-203), and Cooley (1974) has a simple method for improving  $\omega$  for transient problems.

Figure 14 illustrates the rate of convergence of LSOR and LSOR+2DC for test problems 1, 2, and 3 using different acceleration parameters chosen by trial and error. The values exceeding 100 iterations for problem 1 were estimated by using a plot, which is nearly a straight line, of the absolute value of the log of the maximum residual (defined by equation 28) versus the number of iterations. This plot was extrapolated to the value of maximum residual that corresponded roughly to the closure criterion chosen for the problem. The same procedure was used on problem 3 for values exceeding 200 iterations.

For problem 1 the optimum acceleration parameter is 1.87 for LSOR. Two-dimensional correction significantly improves the convergence rate of LSOR for this problem with an optimum acceleration parameter of 1.7. In problem 2, 2DC had no effect on the rate of convergence of LSOR because of the numerous constant-head nodes in the problem. Consequently, the optimum acceleration parameter is 1.6 with or without the application of 2DC. In problem 3, with LSOR oriented across the bedding,  $\omega_{opt}$  is 1.88 for LSOR and about 1.70 for LSOR+2DC. Note in problems 1 and 3 that finding  $\omega_{opt}$  for LSOR is more critical than with LSOR+2DC. LSOR is poorly suited for problem 4 because too many nodes drop out in the iteration process if  $1 < \omega < 2$ . Satisfactory results for problem 4 at the expense of slow convergence are obtained if  $\omega = 0.5$  (See fig. 23.)

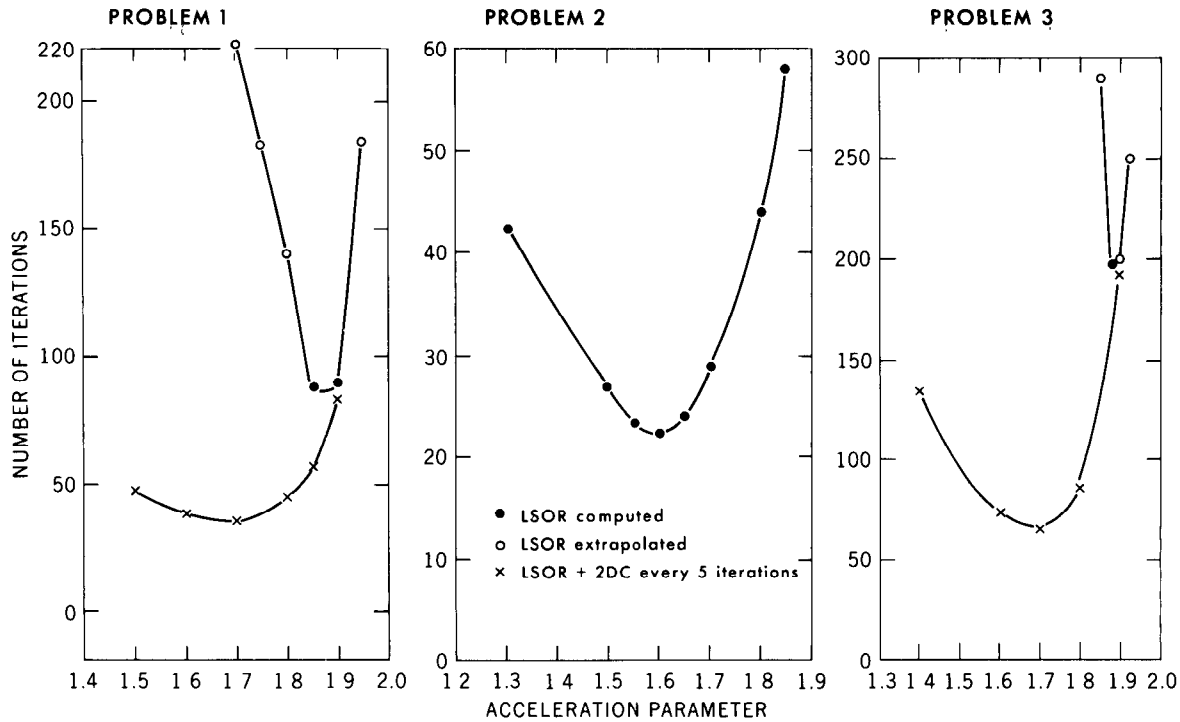


FIGURE 14.—Number of iterations required for solution by LSOR and LSOR + 2DC using different acceleration parameters.

### Alternating-direction implicit procedure

Peaceman and Rachford (1955) described the iterative, alternating-direction implicit procedure for solution of a steady-state (Laplace) equation in two space dimensions. This procedure, however, is equally applicable to transient problems where it has the advantage of allowing larger time steps than can be used with non-iterative ADI. (Non-iterative ADI was used by Pinder and Bredehoeft, 1968.) In the ADI technique, two sets of matrix equations are solved each iteration. The equations for rows in which head values along rows are computed implicitly and those along columns are obtained from the previous column computations are defined as

$$Dh_{j-1}^{n-\frac{1}{2}} + E_r h^{n-\frac{1}{2}} + Fh_{j+1}^{n-\frac{1}{2}} = Q_r, j=1,2,\dots,N_x \quad (21a)$$

in which

$$E_r = -\left(D + F + \frac{S}{\Delta t} + M_l\right);$$

$$Q_r = -Bh_{i-1}^{n-1} + (B+H-M_l)h^{n-1} - Hh_{i+1}^{n-1} - \frac{S}{\Delta t}h_{k-1} + W;$$

$M_l$  is the iteration parameter;

$l$  is the iteration parameter index.

In matrix form equation 21a is

$$\bar{A}_r \bar{h}^{n-\frac{1}{2}} = \bar{Q}_r. \quad (21b)$$

To put equation 21b in residual form, add and subtract  $\bar{A}_r \bar{h}^{n-1}$  to the right-hand side giving

$$\bar{A}_r \bar{h}^{n-\frac{1}{2}} = \bar{Q}_r - \bar{A}_r \bar{h}^{n-1} + \bar{A}_r \bar{h}^{n-1} \quad (21c)$$

Rearrange equation 21c to read:

$$\bar{A}_r \bar{\xi}^{n-\frac{1}{2}} = \bar{R}_r^{n-1} \quad (21d)$$

in which

$$\bar{\xi}^{n-\frac{1}{2}} = \bar{h}^{n-\frac{1}{2}} - \bar{h}^{n-1};$$

$$\bar{R}_r^{n-1} = \bar{Q}_r - \bar{A}_r \bar{h}^{n-1}.$$

Equation 21d is the ADI row formula in residual form. Its matrix form is the same as that for equation 17d and is solved for each row by the Thomas algorithm. To complete the first half of the ADI iteration,  $\bar{h}^{n-\frac{1}{2}}$  is computed by



$$\bar{h}^{n-\frac{1}{2}} = \bar{h}^{n-1} + \bar{\xi}^{n-\frac{1}{2}}.$$

The equations in which head values along columns are considered implicitly and those along rows explicitly are written as:

$$Bh_{i-1}^n + E_c h_i^n + Hh_{i+1}^n = Q_c, \quad i=1, 2, \dots, N_y \quad (22a)$$

in which

$$E_c = -\left(B + H + \frac{S}{\Delta t} + M_l\right);$$

$$Q_c = -Dh_{j-1}^{n-\frac{1}{2}} + (D + F - M_l)h^{n-\frac{1}{2}} - Fh_{j+1}^{n-\frac{1}{2}} - \frac{S}{\Delta t}h_{k-1} + W.$$

Equation 22a in matrix form is

$$\bar{A}_c \bar{h}^n = \bar{Q}_c. \quad (22b)$$

By adding and subtracting  $\bar{A}_c \bar{h}^{n-\frac{1}{2}}$  to the right-hand side of equation 22b, it can be put in the residual form

$$\bar{A}_c \bar{\xi}^n = \bar{R}_c^{n-\frac{1}{2}}; \quad (22c)$$

in which

$$\bar{\xi}^n = \bar{h}^n - \bar{h}^{n-\frac{1}{2}};$$

$$\bar{R}_c^{n-\frac{1}{2}} = \bar{Q}_c - \bar{A}_c \bar{h}^{n-\frac{1}{2}}.$$

Equation 22c is solved for each column by the Thomas algorithm, and the vector  $\bar{h}^n$  for each row is obtained by the equation

$$\bar{h}^n = \bar{h}^{n-\frac{1}{2}} + \bar{\xi}^n.$$

A set of iteration parameters is computed by the equation

$$M_l = \omega_l (B + D + F + H)$$

in which  $\omega$  ranges between a minimum defined by

$$\omega_{\min} = (\text{over grid}) \left[ \frac{\pi^2}{2N_x^2} \frac{1}{1 + \left( \frac{T_{yy[i,j]} (\Delta x_i)^2}{T_{xx[i,j]} (\Delta y_i)^2} \right)} \right. \\ \left. \frac{\pi^2}{2N_y^2} \frac{1}{1 + \left( \frac{T_{xx[i,j]} (\Delta y_i)^2}{T_{yy[i,j]} (\Delta x_i)^2} \right)} \right] \quad (23a)$$

and a maximum given by

$$\omega_{\max} = \begin{cases} 1 & [T_{xx} \cong T_{yy}]; \\ 2 & [T_{xx} >> T_{yy} \text{ or } T_{yy} >> T_{xx}]. \end{cases}$$

The set of parameters are spaced in a geometric sequence given by

$$\omega_{l+1} = \gamma \omega_l \quad (23b)$$

in which

$$\ln \gamma = \frac{\ln(\omega_{\max}/\omega_{\min})}{L-1}. \quad (23c)$$

$L$  = the number of iteration parameters used.

The iteration parameters starting with  $\omega_{\min}$  are cycled until convergence is achieved.

Equation 23a is based on a von Neuman error analysis of the normalized flow equations. (See, for example, Weinstein, Stone, and Kwan, 1969.) It will compute the optimum  $\omega_{\min}$  only for simple problems. For general problems  $\omega_{\min}$  computed by equation 23a may or may not be close to the optimum  $\omega_{\min}$  for the problem. This is illustrated in figure 15 in which the rate of reduction in the maximum residual for arbitrarily chosen minimum parameters is compared with that for  $\omega_{\min}$  computed with equation 23a. Ten parameters were used in problems 1 and 2, and four parameters were used in problem 3. The lines on figure 15 are meant to show the general trend only. The convergence rate using the best  $\omega_{\min}$  in figure 15 is nearly the same as that computed with equation 23a for problem 1, but there is a significant difference in rates for problems 2 and 3. (See figs. 21 and 22.)

The other factor that may be critical in determining the rate of convergence using ADI is the number of parameters. In general, the number of parameters is chosen as 5 if  $\omega_{\max} - \omega_{\min}$  is about two orders of magnitude; if  $\omega_{\max} - \omega_{\min}$  is three or more orders of magnitude, 7 or more parameters are chosen.

For the test problems, the number of iteration parameters were varied from 4 to 10 (fig. 16). The minimum parameter was calculated by equation 23a; the maximum parameter was 1 for problems 1 and 2 and was 2 for problem 3. The number of parameters had a relatively small effect in determining the rate of convergence for problems 1 and 3. For problem 2, however, the computations do not converge using 4 or 5 parameters. Problem 2 can be solved with ADI using 6 to 10 parameters with 10 parameters giving the most rapid convergence. ADI did not give satisfactory solutions for problem 4 (an ex-

cessive number of nodes always drop out of the solution) and, consequently, no results for problem 4 are shown in figure 16.

When difficulties occur with ADI in steady-state simulations, rather than experimenting with the critical minimum parameter or the number of parameters, it may be worthwhile to make the simulation a transient problem. In effect,  $S/\Delta t$  is used as an additional iteration parameter. If the storage coefficient is not made too large or the time step too small,

steady state should be achieved within a reasonable number of time steps with rapid convergence at each time step.

### Strongly implicit procedure

The set of equations (corresponding to equation 8) for the  $3 \times 3$  problem in figure 13 may be expressed in matrix form as

$$\bar{A} \bar{h} = \bar{Q} \quad (24)$$

$$\begin{array}{c}
 \underbrace{\hspace{10em}}_{N_x + 1 \text{ elements}} \\
 \left. \begin{array}{c} N_x + 1 \text{ elements} \\ \left[ \begin{array}{cccccccc} E_1 & F_1 & 0 & H_1 & & & & \\ D_2 & E_2 & F_2 & 0 & H_2 & & & \\ 0 & D_3 & E_3 & 0 & 0 & H_3 & & \\ B_4 & 0 & 0 & E_4 & F_4 & 0 & & \\ & B_5 & 0 & 0 & E_5 & F_5 & H_4 & \\ & & B_6 & 0 & D_6 & E_6 & 0 & H_6 \\ & & & B_7 & 0 & 0 & E_7 & F_7 & 0 \\ & & & & B_8 & 0 & D_8 & E_8 & F_8 \\ & & & & & B_9 & 0 & D_9 & E_9 \end{array} \right] \begin{bmatrix} h_1 \\ h_2 \\ h_3 \\ h_4 \\ h_5 \\ h_6 \\ h_7 \\ h_8 \\ h_9 \end{bmatrix} = \begin{bmatrix} Q_1 \\ Q_2 \\ Q_3 \\ Q_4 \\ Q_5 \\ Q_6 \\ Q_7 \\ Q_8 \\ Q_9 \end{bmatrix} \end{array} \right\}
 \end{array}$$

Direct solution of equation 24 by Gaussian elimination usually requires more work and computer storage than iterative methods for problems of practical size because  $\bar{A}$  decomposes into a lower triangular matrix with non-zero elements from  $B$  to  $E$  in each row and an upper triangular matrix with non-zero elements from  $E$  to  $H$  in each row. All of these intermediate coefficients must be computed during Gaussian elimination, and the coefficients in the upper triangular matrix must be saved for backward substitution.

To reduce the computation time and storage requirements of direct Gaussian elimination, Stone (1968) developed an iterative method using approximate factorization. In this approach a modifying matrix  $\bar{B}$  is added to  $\bar{A}$  forming  $(\bar{A} + \bar{B})$  so that equation 24 becomes

$$(\bar{A} + \bar{B}) \bar{h} = \bar{Q} + \bar{B} \bar{h}. \quad (25)$$

$(\bar{A} + \bar{B})$  can be made close to  $\bar{A}$  but can be factored into the product of a lower triangular matrix  $\bar{L}$  and an upper triangular matrix  $\bar{U}$ , each of which has no more than three non-zero elements in each row, regardless of the size of  $N_x$  and  $N_y$ . Therefore, if the right-hand side of equation 25 is known, simple

recursion formulas can be derived, resulting in a considerable savings in computer time and storage. This leads to the iteration scheme

$$(\bar{A} + \bar{B}) \bar{h}^n = \bar{Q} + \bar{B} \bar{h}^{n-1}. \quad (26)$$

In order to transform equation 26 into a residual form,  $\bar{A} \bar{h}^{n-1}$  is subtracted from both sides giving

$$(\bar{A} + \bar{B}) \bar{\xi}^n = \bar{R}^{n-1} \quad (27)$$

in which

$$\begin{aligned}
 \bar{\xi}^n &= \bar{h}^n - \bar{h}^{n-1}; \\
 \bar{R}^{n-1} &= \bar{Q} - \bar{A} \bar{h}^{n-1}.
 \end{aligned} \quad (28)$$

The iterative scheme defined by equation 26 or 27 is closer to direct methods of solution (more implicit) than ADI (hence the term strongly implicit procedure or SIP). The SIP algorithm requires (1) relationships among the elements of  $\bar{L}$ ,  $\bar{U}$  and  $(\bar{A} + \bar{B})$  defined by rules of matrix multiplication for the equation

$$\bar{L} \bar{U} = (\bar{A} + \bar{B}), \quad (29)$$

and (2) relationships among the elements of  $\bar{A}$  and  $(\bar{A} + \bar{B})$ .

$\bar{L}$  and  $\bar{U}$  have the following form for a general  $3 \times 3$  problem (much of the notation is adapted from Remson, Hornberger, and Molz, 1971);



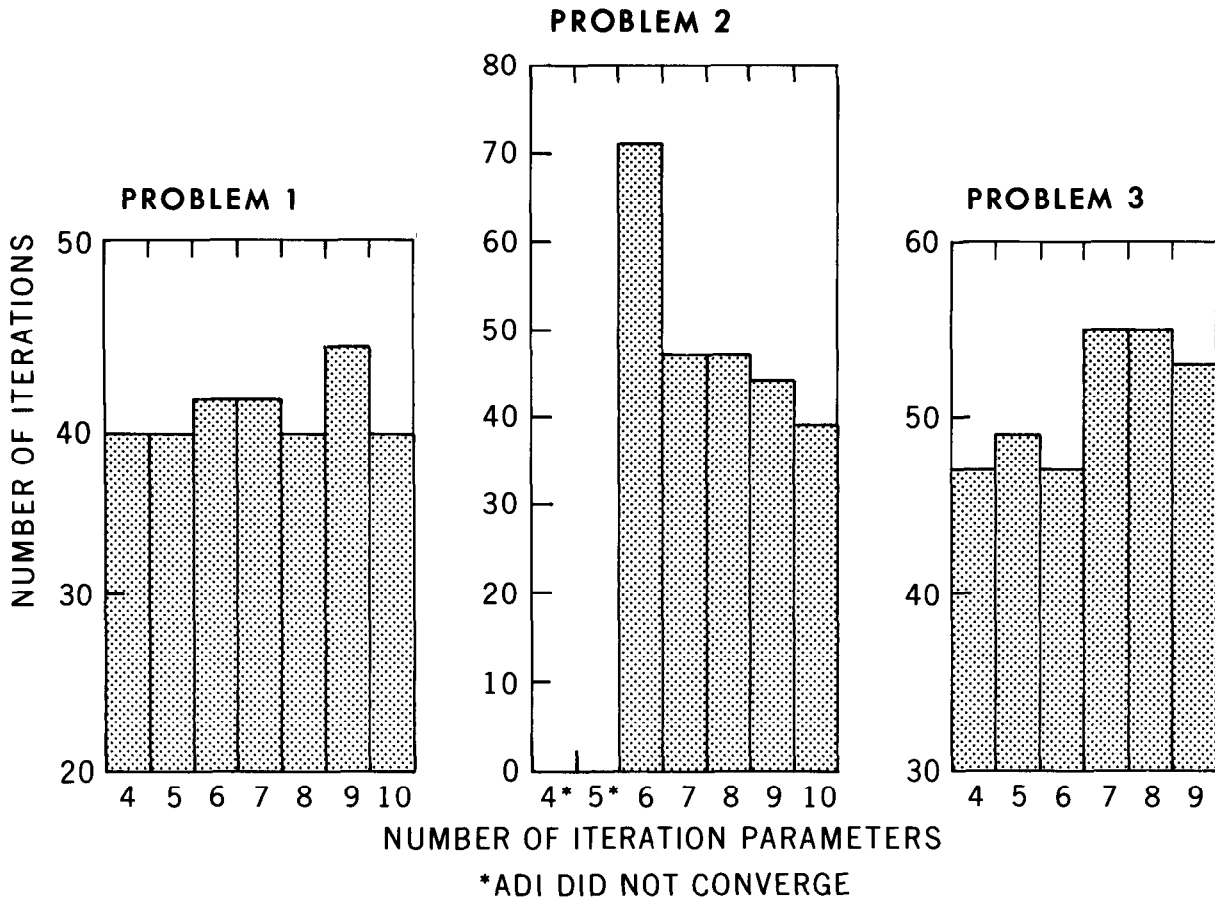


FIGURE 16.—Number of iterations required for solution of the test problems with ADI using different numbers of parameters.

The product  $\overline{L\overline{U}} = (\overline{A+B})$  is

$$(\overline{A+B}) = \begin{bmatrix} \hat{E}_1 & \hat{F}_1 & 0 & \hat{H}_1 & & & & & & & \\ \hat{D}_2 & \hat{E}_2 & \hat{F}_2 & \hat{G}_2 & \hat{H}_2 & & & & & & \\ 0 & \hat{D}_3 & \hat{E}_3 & \boxed{\hat{F}_3} & \hat{G}_3 & \hat{H}_3 & & & & & \\ \hat{B}_4 & \hat{C}_4 & \boxed{\hat{D}_4} & \hat{E}_4 & \hat{F}_4 & \boxed{\hat{G}_4} & \hat{H}_4 & & & & \\ & \hat{B}_5 & \hat{C}_5 & \hat{D}_5 & \hat{E}_5 & \hat{F}_5 & \hat{G}_5 & \hat{H}_5 & & & \\ & & \hat{B}_6 & \boxed{\hat{C}_6} & \hat{D}_6 & \hat{E}_6 & \boxed{\hat{F}_6} & \hat{G}_6 & \hat{H}_6 & & \\ & & & \hat{B}_7 & \hat{C}_7 & \boxed{\hat{D}_7} & \hat{E}_7 & \hat{F}_7 & \boxed{\hat{G}_7} & & \\ & & & & \hat{B}_8 & \hat{C}_8 & \hat{D}_8 & \hat{E}_8 & \hat{F}_8 & & \\ & & & & & \hat{B}_9 & \boxed{\hat{C}_9} & \hat{D}_9 & \hat{E}_9 & & \end{bmatrix}$$

Because of the boundary conditions, the elements of  $(\overline{A+B})$  inside squares will be zero for the  $3 \times 3$  problem illustrated in figure 13.

The relationships among the elements of  $\overline{L}$ ,  $\overline{U}$ , and  $(\overline{A+B})$  are

$$\alpha = \hat{B} \quad (30a)$$

$$\alpha\delta_{i-1} = \hat{C} \quad (30b)$$

$$\beta = \hat{D} \quad (30c)$$

$$\gamma + \alpha\eta_{i-1} + \beta\delta_{j-1} = \hat{E} \quad (30d)$$

$$\gamma\delta = \hat{F} \quad (30e)$$

$$\beta\eta_{j-1} = \hat{G} \quad (30f)$$

$$\gamma\eta = \hat{H} \quad (30g)$$

where the  $i$  and  $j$  subscripts refer to the location on the model grid, not in matrix  $(\overline{A+B})$ .

In order to use equations 30a–30g as the basis of a numerical technique for solving equation 24 efficiently by elimination, relationships between the elements of  $\overline{A}$  and  $(\overline{A+B})$  must be defined. One possibility is to let the elements correspond exactly and ignore the  $\hat{C}$  and  $\hat{G}$  diagonal in  $(\overline{A+B})$ . Stone (1968), however, found that this could not be used as the basis of a rapidly convergent iterative procedure. Instead, he defined a family of modified matrices starting with 30b and 30f.

Then the other elements of  $(\overline{A+B})$  can be defined as equal to the corresponding elements in  $\overline{A}$  plus a linear combination of  $\hat{C}$  and  $\hat{G}$ . For example

$$\hat{B} = B + \phi_1\hat{C} + \phi_2\hat{G}$$

in which  $\phi_1$  and  $\phi_2$  are constants depending on the problem being solved.

What are appropriate linear combinations of  $\hat{C}$  and  $\hat{G}$  with the elements of  $\overline{A}$ ? If equation 27 is written for node  $(i,j)$ , non-zero coefficients appear not only for the unknowns in the original difference equation but also for  $\xi_{i-1,j+1}^n$  and  $\xi_{i+1,j-1}^n$ . This is illustrated in figure 17. To minimize the effects of the terms introduced in forming the modified matrix equation,  $\overline{B}\xi^n$  for the node  $(i,j)$  is defined as

$$\hat{C}[\xi_{i-1,j+1}^n - \omega(\xi_{i-1}^n + \xi_{j+1}^n - \xi^n)] + \hat{G}[\xi_{i+1,j-1}^n - \omega(\xi_{i+1}^n + \xi_{j-1}^n - \xi^n)] \quad (31)$$

where the terms in parentheses are second-order correct approximations for  $\xi_{i-1,j+1}$ , and  $\xi_{i+1,j-1}$ , respectively. (See Remson, Hornberger, and Molz, 1971, p. 226, for derivation of these approximations.) To consider these terms good approximations to  $\xi_{i-1,j+1}$  and

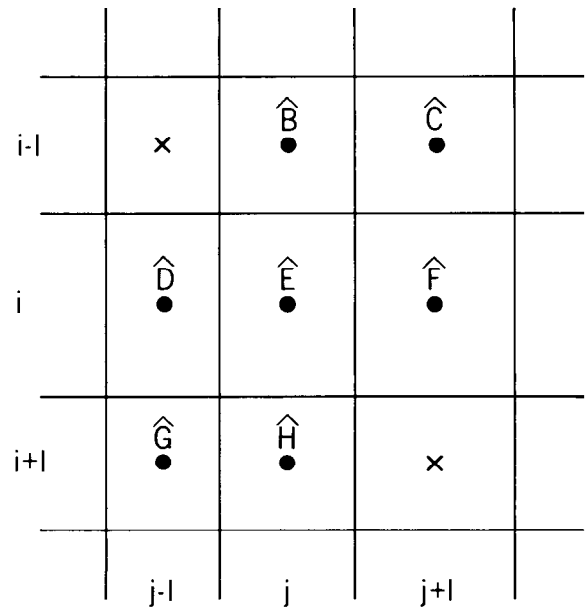


FIGURE 17.—Coefficients of unknowns in equation 27.

$\xi_{i+1,j-1}$  an iteration parameter,  $\omega$ , is added. The value of  $\omega$  ranges between 0 and 1, and its computation is discussed at the end of this section.

With the definition of  $\overline{B}$  (31), the iteration scheme (equation 27) becomes

$$B\xi_{i-1}^n + D\xi_{j-1}^n + E\xi^n + F\xi_{j+1}^n + H\xi_{i+1}^n + \hat{C}[\xi_{i-1,j+1}^n - \omega(\xi_{i-1}^n + \xi_{j+1}^n - \xi^n)] + \hat{G}[\xi_{i+1,j-1}^n - \omega(\xi_{i+1}^n + \xi_{j-1}^n - \xi^n)] = R^{n-1} \quad (32)$$

Collecting coefficients in equation 32 associated with the nodal positions in the original difference equation gives the desired linear combinations of  $\hat{C}$  and  $\hat{G}$  with the elements of  $\overline{A}$  that define the remaining elements of  $(\overline{A+B})$ :

$$\hat{B} = B - \omega\hat{C} \quad (33a)$$

$$\hat{D} = D - \omega\hat{G} \quad (33b)$$

$$\hat{E} = E + \omega\hat{C} + \omega\hat{G} \quad (33c)$$

$$\hat{F} = F - \omega\hat{C} \quad (33d)$$

$$\hat{H} = H - \omega\hat{G} \quad (33e)$$

The coefficient  $\hat{C}$  is obtained explicitly by combining equations 33a, 30a, and 30b as

$$\hat{C} = \frac{\delta_{i-1}B}{1 + \omega\delta_{i-1}} \quad (34a)$$

Finally combining equation 33b and equations 30c and 30f gives

$$\hat{G} = \frac{\eta_{j-1}D}{1 + \omega\eta_{j-1}} \quad (34b)$$

Equations 34, 33 and 30 (in that order) are the first part of the SIP algorithm.

Equation 28 written for node  $(i,j)$  is

$$R^{n-1} = Q - (Bh_{i-1}^{n-1} + Dh_{j-1}^{n-1} + Eh^{n-1} + Fh_{j+1}^{n-1} + Hh_{i+1}^{n-1}).$$

As in the Thomas algorithm, the vector  $\bar{\xi}^n$  is obtained by a process of forward and backward substitution. Combining equations 27 and 29 gives

$$\bar{L}\bar{U}\bar{\xi}^n = \bar{R}^{n-1} \quad (35)$$

Define an intermediate vector  $\bar{V}^n$  by

$$\bar{U}\bar{\xi}^n = \bar{V}^n. \quad (36)$$

Then equation 35 becomes

$$\bar{L}\bar{V}^n = \bar{R}^{n-1}. \quad (37)$$

$\bar{V}^n$  is first computed by forward substitution. This can be seen by writing equation 37 for node  $(i,j)$ :

$$\alpha V_{i-1}^n + \beta V_{j-1}^n + \gamma V^n = R^{n-1}$$

or

$$V^n = (R^{n-1} - \alpha V_{i-1}^n - \beta V_{j-1}^n) / \gamma.$$

The vector  $\bar{\xi}^n$  may then be computed by backward substitution. Equation 36 for node  $(i,j)$  is

$$\xi^n + \delta \xi_{i+1}^n + \eta \xi_{i+1}^n = V^n$$

or

$$\xi^n = V^n - \delta \xi_{i+1}^n - \eta \xi_{i+1}^n.$$

Stone (1968) recommends an alternating computational procedure. On odd iterations, the equations are ordered in a "normal" manner as shown in figure 13. On even iterations, the numbering scheme is changed to that illustrated in figure 18. This has the effect of making non-zero coefficients appear for the heads  $h_{i-1,j-1}$  and  $h_{i+1,j+1}$  (the X's in fig. 17) instead of  $h_{i-1,j+1}$  and  $h_{i+1,j-1}$  and significantly improves the convergence rate. Note that some of the recursion equations are modified by reordering the grid points in the "reverse" manner. The modifications required for the reverse algorithm are

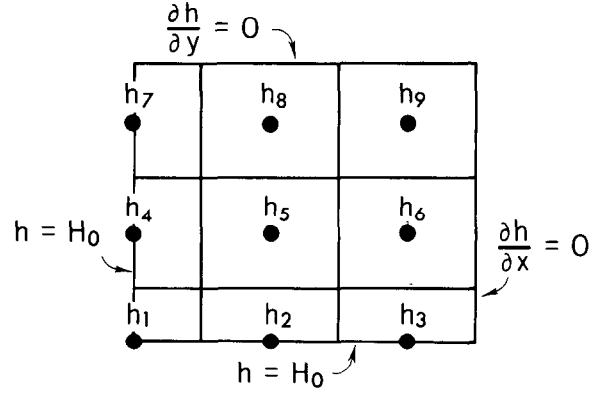


FIGURE 18.—Reverse numbering scheme for  $3 \times 3$  problem.

$$\hat{C} = \frac{\delta_{i+1}H}{1 + \omega\delta_{i+1}};$$

$$\hat{B} = H - \omega\hat{C};$$

$$\hat{H} = B - \omega\hat{G};$$

$$\gamma = E - \alpha\eta_{i+1} - \beta\delta_{j-1};$$

$$V^n = (R^{n-1} - \alpha V_{i+1}^n - \beta V_{j-1}^n) / \gamma;$$

$$\xi^n = V^n - \delta \xi_{j+1}^n - \eta \xi_{i-1}^n.$$

The iteration parameters are computed by equations given in Stone (1968). For variable transmissivity and grid spacing, Stone's equation is

$$(1 - \omega_{\max}) = \sum_{i=1}^{N_y} \sum_{j=1}^{N_x} \text{Min} \left[ \frac{2(\delta x_j)^2}{1 + \left( \frac{T_{yy}[i,j](\delta x_i)^2}{T_{xx}[i,j](\delta y_i)^2} \right)}, \frac{2(\delta y_i)^2}{1 + \left( \frac{T_{xx}[i,j](\delta y_i)^2}{T_{yy}[i,j](\delta x_j)^2} \right)} \right] + (N_x \times N_y) \quad (38)$$

in which

$$\delta x = \Delta x_j / \text{width of model}$$

$$\delta y = \Delta y_i / \text{length of model}$$

Equation 38 computes an arithmetic average of  $\omega_{\max}$  for the algorithm.

The remaining iteration parameters are computed by

$$1 - \omega_{l+1} = (1 - \omega_{\max})^{1/(L-1)}, l = 0, 1, \dots, L-1$$

in which  $L$  is the number of parameters in a cycle.

Stone (1968) recommends using a minimum of four parameters, each used twice in

succession, starting with the largest first. Weinstein, Stone, and Kwan (1969), however, indicate that it is not necessary to start with the largest parameter first or to repeat them.

The results using different numbers and sequences of parameters for the three test problems are shown in figure 19. Except for the sequence 4, 3, 2, 1 in problem 1 the number of iterations required for solution varies up to a maximum of 50 percent for the parameter sequences tested. Several parameter sequences (for example, 1, 2, 3, 4, 5) give convergence near the maximum observed rate for all problems. This result suggests that conducting numerical experiments to determine the best sequence of parameters for a particular problem is generally not justified.

Weinstein, Stone, and Kwan (1969) have a slightly different definition of the maximum parameter ( $1 - \omega_{\max} = \text{ADI minimum parameter}$

ter). Their definition of the maximum parameter (which is the maximum over the model, not the arithmetic average of values computed for each node) was used in solving several test problems. In every case convergence was faster using equation 38 to compute the maximum parameter.

Stone (1968) states that a more general form of equation 27 includes another iteration parameter,  $\beta'$ , to multiply the term  $\bar{R}^{n-1}$ . His experience indicated, however, that values of  $\beta'$  other than unity did not generally improve the method. In contrast, the use of  $\beta'$  other than unity has proven to be effective for some of the test problems. In fact, for the fourth problem, a value of  $\beta'$  less than unity is required to obtain a reasonable solution using SIP. Results for problem 4 are not shown in figure 19 because the best sequence of parameters (No. 3) for problem 2 was used in experimenting with the parameter  $\beta'$ .

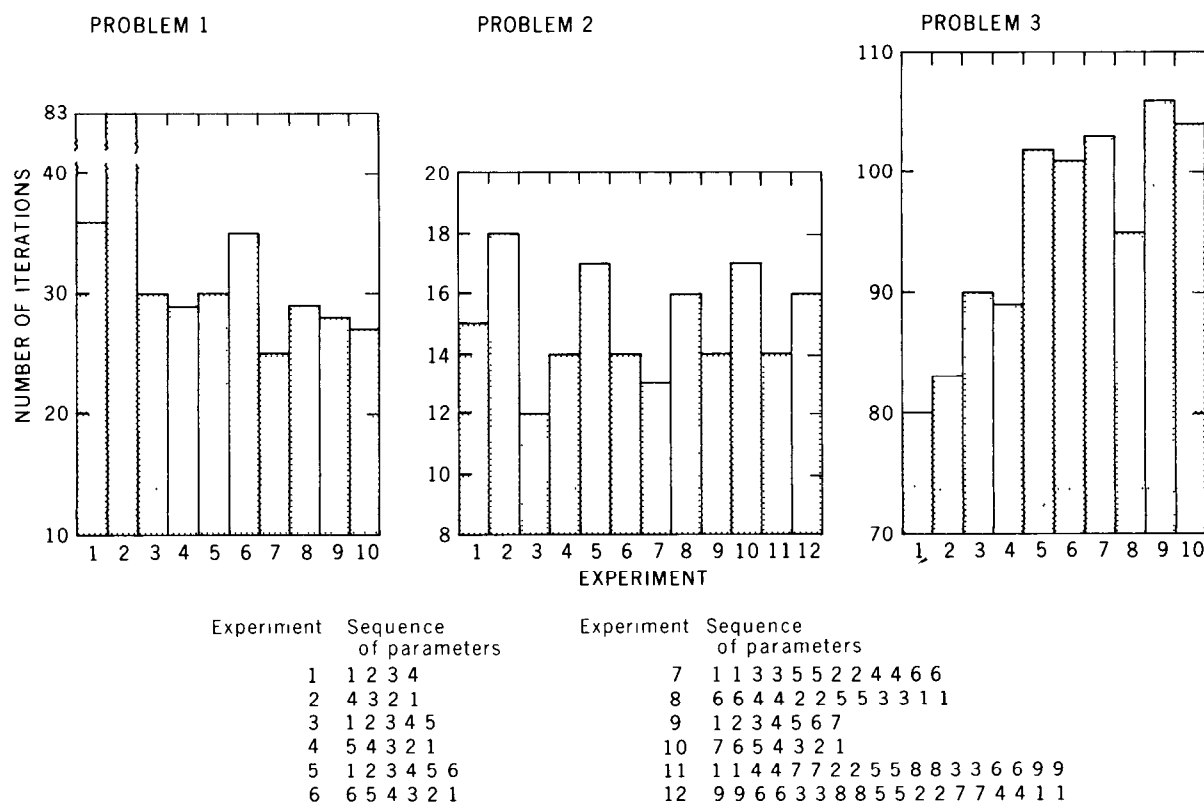


FIGURE 19.—Iterations required for solution of the test problems by SIP using different numbers and sequences of parameters.

## Comparison of Numerical Results

The rate of convergence using different numerical techniques for solving the test problems is compared in figures 20 to 23. The best results from the experiments with each iterative technique are used in the comparisons. Two curves (except for fig. 23) are shown for SIP: one with the parameter  $\beta' = 1$  and the other with the best rate of convergence for  $\beta' \neq 1$ . The sequence of  $\omega$  parameters is the same for both curves. Two curves are also shown for ADI: one in which the minimum parameter was calculated with equation 23a (indicated by an asterisk in the figures); the other with the best minimum parameter shown on figure 15.

In figures 20 to 23 the absolute value of the maximum residual for each iteration is plotted versus computation time where one unit of work is equal to the time required to complete one SIP iteration. Relative work per iteration is about 1 for ADI, 0.6 for LSOR, and 0.8 for LSOR+2DC. The maximum residual for SIP and ADI fluctuates from a maximum to a minimum over each cycle of parameters. For clarity, the curves connect the local minima for these two methods. Comparisons in figures 20-23 should be made on the basis of the horizontal displacement of the curves, not on the basis of the termination of the curves. This is similar to the type of comparisons made by Stone (1968).

Figure 20 shows the results for problem 1 (10 parameters for ADI,  $\omega = 1.87$  for LSOR,  $\omega = 1.7$  for LSOR+2DC, parameter sequence, 1,1,3,3,5,5,2,2,4,4,6,6, for SIP). Of the sequence of  $\beta'$  parameters tried, the minimum work required to reduce the residual is obtained with  $\beta' = 1.4$ , but this is only moderately better than using  $\beta' = 1.0$ . ADI converges as rapidly as SIP for the first cycles of iteration, but from that point on converges slower than the other iterative techniques. The two ADI curves show about the same rate of convergence for this problem. Next to SIP, LSOR+2DC is most attractive for this problem.

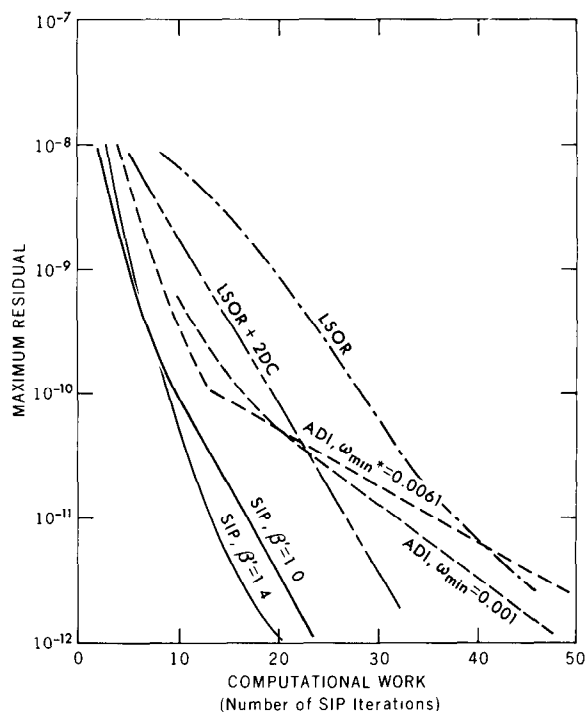


FIGURE 20.—Computational work required by different iterative techniques for problem 1.

The results for problem 2 are shown in figure 21 (10 parameters for ADI,  $\omega = 1.6$  for LSOR and LSOR+2DC, parameter sequence 1,2,3,4,5 for SIP). SIP requires the least amount of work for this problem (using  $\beta' \neq 1.0$  does not significantly reduce the work required). LSOR and ADI using the best  $\omega_{\min}$  from figure 15 are competitive with SIP. ADI using  $\omega_{\min}$  computed with equation 23a requires about twice as much computational work. LSOR and LSOR+2DC take the same number of LSOR iterations so that the extra work required for 2DC is wasted for this problem.

In figure 22, the results using 4 parameters for ADI, the parameter sequence 1,2,3,4 for SIP,  $\omega = 1.88$  for LSOR and  $\omega = 1.70$  for LSOR+2DC are plotted for problem 3. In this problem LSOR (with solution lines oriented along columns), ADI with  $\omega_{\min}$  computed with equation 23a, and SIP with  $\beta' = 1$  are competitive. Convergence is significantly improved by adding 2DC to LSOR, choosing the best  $\omega_{\min}$  from figure 15 for ADI and letting  $\beta' = 1.5$  with SIP.



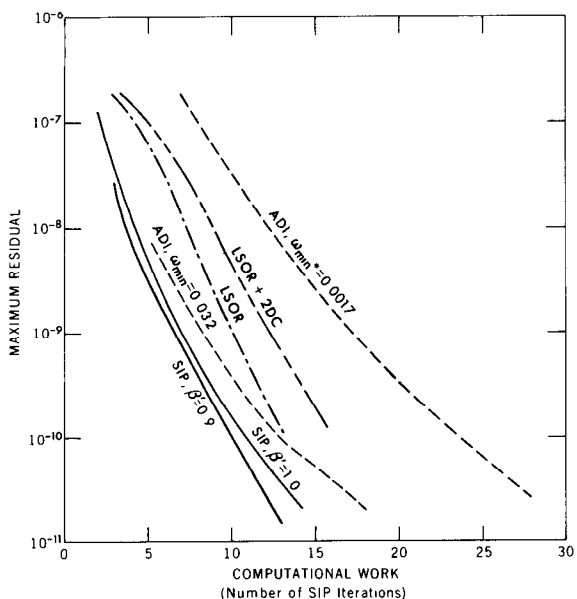


FIGURE 21.—Computational work required by different iterative techniques for problem 2.

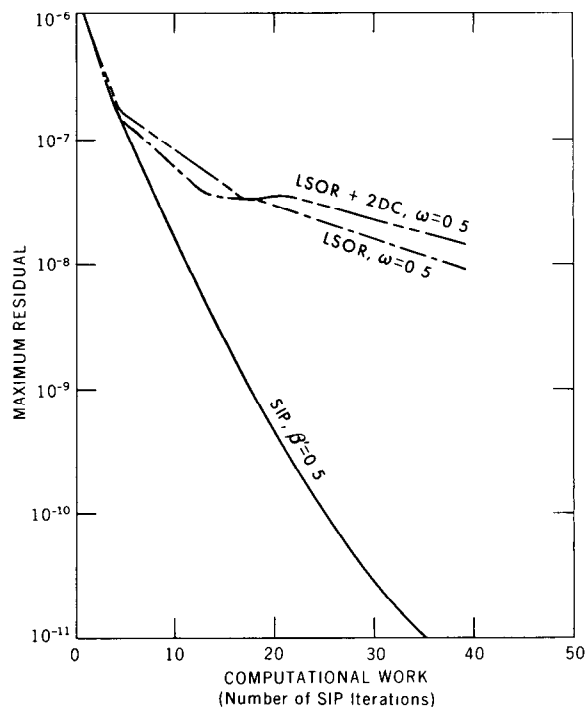


FIGURE 23.—Computational work required by different iterative techniques for problem 4.

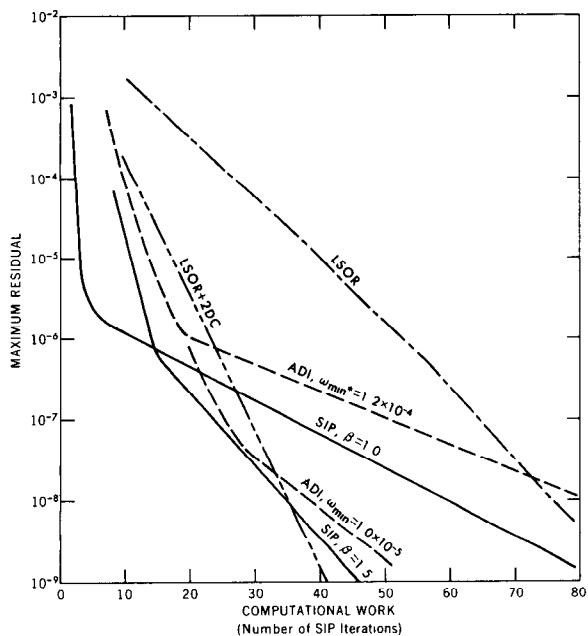


FIGURE 22.—Computational work required by different iterative techniques for problem 3.

The results for problem 4 are shown in figures 23 and 24. The  $\omega$  iteration parameter sequence for SIP is 1,2,3,4,5, and the two-dimensional correction is applied every fifth iteration for LSOR+2DC. Konikow (oral

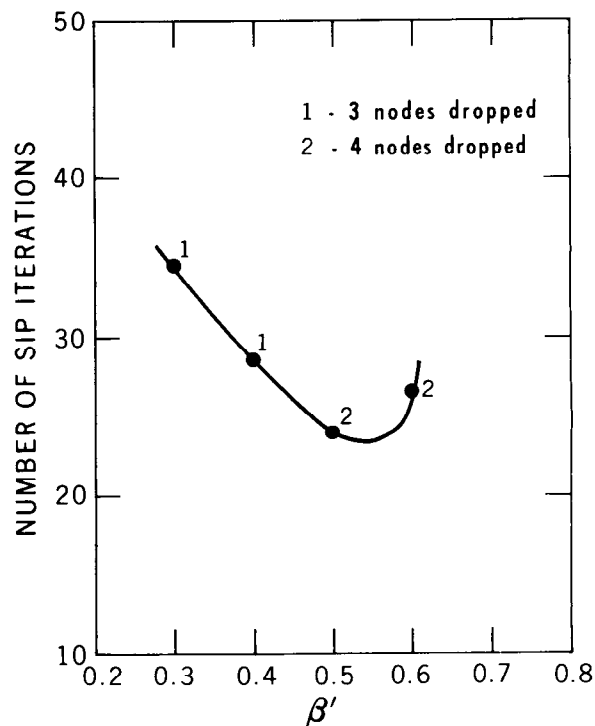


FIGURE 24.—Number of iterations required for solution of problem 4 by SIP using different values of  $\beta'$ .

commun., 1975) was unable to obtain a solution to problem 4 using ADI due to oscillations that eliminated nodes that should have been in the solution. This problem occurred not only with ADI but also with LSOR and LSOR+2DC with  $\omega > 0.6$  and with SIP with  $\beta' > 0.6$ . The oscillations are apparently caused in part by the nonlinearities of the water-table problem and the necessity to calculate transmissivity at the known iteration level. In a water-table simulation the transmissivity is set to zero and nodes are dropped from the aquifer if the computed head is below the base of the aquifer. For problem 4, at least 3 nodes should be dropped with the initial conditions used.

A solution to problem 4 in which 3 to 4 nodes are dropped is obtained with LSOR and LSOR+2DC when  $\omega = 0.5$  at the expense of slow convergence. Clearly the most suitable method for this problem is SIP with  $\beta' \leq 0.6$  (fig. 23). In effect the use of  $\beta' < 1$  for SIP and  $\omega < 1$  for LSOR represents "under-relaxation" and has the effect of dampening oscillations of head from one iteration to the next. This reduces the tendency for incorrect deletion of nodes from the solution.

Solution of problem 4 emphasizes the advantage of the extra SIP iteration parameter. The optimum value of  $\beta'$  inferred from figure 24 is about 0.5. Note in figure 24 that an additional node is dropped for  $\beta' = 0.5$  and 0.6. However, the effect of this node on the remainder of the solution is negligible. For  $\beta' > 0.6$ , either convergence was not obtained or excessive numbers of nodes were dropped for those cases that did converge.

The numerical experiments included in this report support the general conclusions of Stone (1968) and Weinstein, Stone, and Kwan (1969) that SIP is a more powerful iterative technique than ADI for most problems. SIP is attractive, not only because of its relatively high convergence rates but because it is generally not necessary to conduct numerical experiments to select a suitable sequence of parameters. SIP has the disadvantage of requiring 3 additional  $N_x \times N_y$  arrays.

For the first three problems examined here, ADI is a slightly better technique than LSOR when  $\omega_{\min}$  near the optimum is used. Although this result agrees with Bjordammen and Coats (1969) who concluded that ADI is superior to LSOR for the oil reservoir problems they investigated, it is deceptive because less work is required to obtain  $\omega_{\text{opt}}$  for LSOR than is required to find the best  $\omega_{\min}$  for ADI by trial and error. Furthermore, LSOR is clearly superior to ADI in application to problem 4 where a solution was not possible with ADI as used in this simulator.

LSOR+2DC seems to be particularly useful with problems dominated by no-flux boundaries. The correction procedure can significantly improve the rate of convergence of LSOR even in problems such as problem 3 where all  $\beta_j$  are zero and non-zero  $\alpha_i$  occur for the lower half of the model only.

## Considerations in Designing an Aquifer Model

### Boundary conditions

An aquifer system is usually larger than the project area. Nevertheless the physical boundaries of the aquifer should be included in the model if it is feasible. Where it is impractical to include one or more physical boundaries (for example, in an alluvial valley that may be several hundred miles long) the finite-difference grid can be expanded and the boundaries located far enough from the project area so that they will have negligible effect in the area of interest during the simulation period. The influence of an artificial boundary can be checked by comparing the results of two simulation runs using different artificial boundary conditions.

Boundaries that can be treated by the model are of two types: constant head and constant flux. Constant-head boundaries are specified by assigning a negative storage coefficient to the nodes that define the constant-head boundary. This indicates to the program that these nodes are to be skipped in the computations.

A constant flux may be zero (impermeable boundaries) or have a finite value. A zero-flux boundary is treated by assigning a value of zero transmissivity to nodes outside the boundary. The harmonic mean of the transmissivity at the cell boundary is zero, and consequently, the flux across the boundary is zero. A no-flow boundary is inserted around the border of the model as a computational expediency, and constant-head or finite-flux boundaries are placed inside this border. A finite-flux boundary is treated by assigning recharge (or discharge) wells to the appropriate nodes. Figure 25 illustrates various types of boundary conditions.

The type of boundaries appropriate to the field problem may require careful consideration. In particular, should streams be treated as constant-head boundaries or are they more realistically treated as partially penetrating with a leaky streambed? If a leaky streambed is used, note that the leakage occurs over the area of the blocks assigned to the stream. If the area of the streambed is less than the area of the blocks, the ratio of streambed hydraulic conductivity to thickness can be proportionately reduced to make the amount of leakage realistic.

### Initial conditions

In many simulations, the important results are not the computed head but the changes in head caused by a stress such as pumping wells. For this objective in a confined aquifer for which the equations are linear, there is no need to impose the natural flow system as the initial condition since the computed draw-down can be superimposed on the natural flow system, if desired.

If initial conditions are specified so that transient flow is occurring in the system at the start of the simulation, it should be recognized that water levels will change during the simulation, not only in response to the new pumping stress, but also due to the initial conditions. This may or may not be the intent of the user.

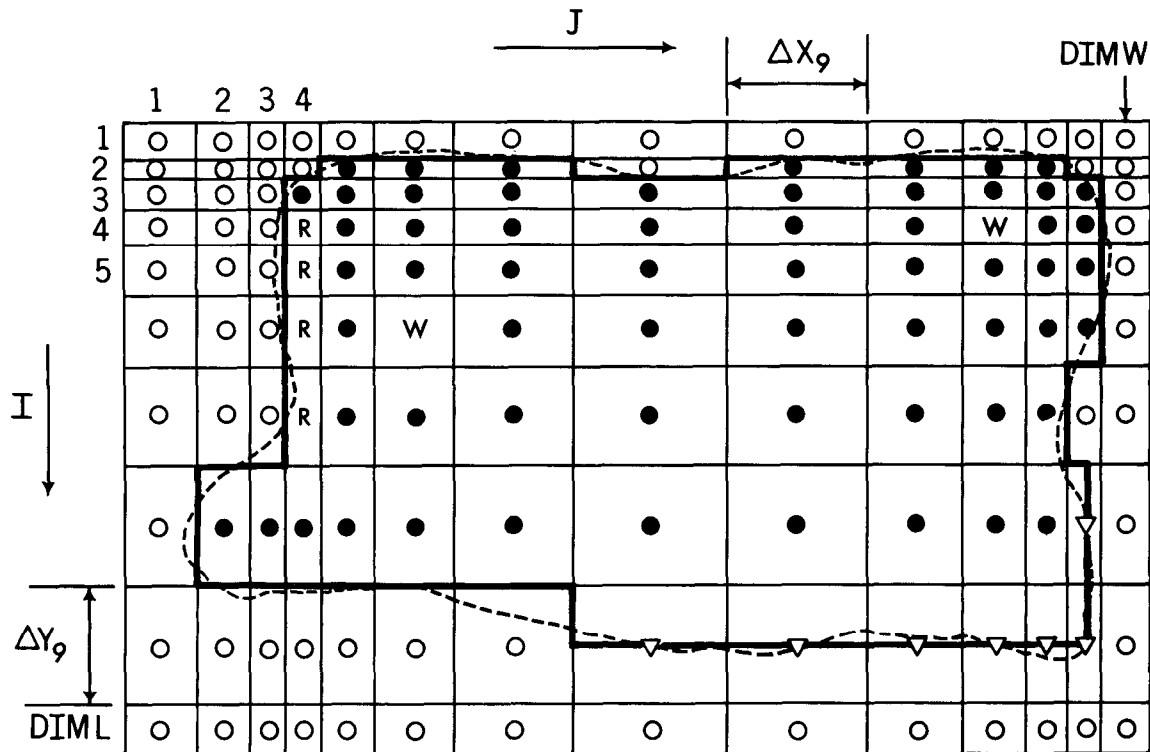
To start from steady-state conditions in which flow is occurring, the model can be used

to compute the initial head by leaving out the new stress (for example, wells) and setting all storage terms to zero. This is also a useful calibration procedure to compute unknown terms such as the ratio of hydraulic conductivity to thickness for leakage.

### Designing the finite-difference grid

In designing a finite-difference grid, the following considerations should be kept in mind:

1. Nodes representing pumping and observation wells should be close to their respective positions to facilitate calibration. If several pumping wells are close together, their discharge may be lumped and assigned to one node since discharge is distributed over the area of the cell.
2. Boundaries within the project area should be located accurately. Distant boundaries can be located approximately and with fewer nodes by expanding the grid. In expanding a finite-difference grid in the positive  $X$  direction, experience has shown that restricting the ratio  $\Delta X_j / \Delta X_{j-1} \leq 1.5$  will avoid large truncation errors and possible convergence problems.
3. Nodes should be placed close together in areas where there are spatial changes in transmissivity. For example, in cross-sectional problems with aquifers separated by confining beds, many layers of nodes are required in the confining bed to obtain a good approximation of the head distribution (and consequently the flux) during transient conditions.
4. The grid should be oriented so that a minimum of nodes are outside the aquifer. The orientation of the grid with respect to latitude and longitude or some other geographic grid system would be a secondary consideration. However, if the aquifer is anisotropic, the grid should be oriented with its axes parallel to the principal directions of the transmissivity tensor. Otherwise,



## EXPLANATION

## Node symbols

Inside aquifer (transmissivity > 0)

w Discharge well

R Recharge well

▽ Constant head

● Node without wells or specified head

## Outside aquifer

○ Transmissivity = 0

--- Aquifer boundary

— Mathematical boundary

DIML Number of rows

DIMW Number of columns

## Boundary conditions



Constant head

Constant flux



$\frac{\partial h}{\partial x} = 0$



$\frac{\partial h}{\partial x} = C$

FIGURE 25.—Variable, block-centered grid with mixed boundary conditions.

the flow equation would include cross-product terms and the solution would be restricted to ADI and LSOR because additional diagonals appear in the coefficient matrix and SIP, in its usual form, cannot be used.

5. The rows should be numbered in the short dimension for the alphameric plot on the line printer or for plotting data with an X-Y plotter. On these plots, the X-direction is vertical and, for practical purposes, this dimension is unlimited. The Y direction is across the page which limits this dimension to the maximum width of the page. (See fig. 26.)
6. The core requirements and computation time are proportional to the number of nodes representing the aquifer.

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**COMPUTER PROGRAM AND RELATED DATA**

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## Attachment I

## Notation

$\bar{A}$	coefficient matrix;	$q'$	transient part of $q'$ ( $Lt^{-1}$ );
$\bar{A}_c, \bar{A}_r$	coefficient matrices for ADI column and row equations;	$q_{0i}$	evapotranspiration flux ( $Lt^{-1}$ );
$\bar{A}_\lambda$	LSOR coefficient matrix;	$q_{re}$	recharge flux ( $Lt^{-1}$ );
$b$	saturated thickness of the aquifer ( $L$ );	$Q$	known term in difference equation;
$\bar{B}$	modifying matrix for SIP;	$Q_{0i}$	maximum evapotranspiration rate ( $Lt^{-1}$ );
$B, D, E, F, H$	coefficients in difference equation;	$Q_n, Q_r$	known terms in equations defining ADI;
$\hat{B}, \hat{C}, \hat{D}, \hat{E}, \hat{F}, \hat{G}, \hat{H}$	coefficients of $(\bar{A} + \bar{B})$ ;	$Q_w$	well discharge ( $L^3t^{-1}$ );
$B', D', E', F', H'$	coefficients of equations defining 2DC;	$Q_\lambda$	known term in equation defining LSOR;
$BE, G, W$	notation used in Thomas algorithm;	$r$	radial distance from center of pumping well ( $L$ );
$E_r, E_o$	coefficients in equations defining ADI;	$r_e$	effective radius for a well block ( $L$ );
$ET_s$	depth below land surface at which evapotranspiration ceases ( $L$ );	$r_w$	well radius ( $L$ );
$G$	elevation of the land surface ( $L$ );	$r_1$	radius equivalent to the average grid spacing for the well block ( $L$ );
$h$	hydraulic head ( $L$ );	$R$	residual;
$h^t$	intermediate head value ( $L$ );	$R^{n*}$	residual computed for 2DC;
$h^{n*}$	corrected head at iteration $n$ ( $L$ );	$R'$	sum of residuals for 2DC;
$h_{i,j,0}$	initial head in the aquifer ( $L$ );	$R_o, R_r$	residuals for ADI column and row computations;
$\hat{h}_{i,j,0}$	hydraulic head on the other side of the confining bed ( $L$ );	$R_\lambda$	LSOR residual;
$h_w$	hydraulic head in a well ( $L$ );	$S$	storage coefficient (dimensionless);
$H_{i,j}$	saturated thickness of the aquifer at radius $r_e$ ( $L$ );	$S_c$	specific storage of the confining bed ( $L^{-1}$ );
$H_w$	saturated thickness of the aquifer at radius $r_w$ ( $L$ );	$S_y$	specific yield (dimensionless);
$i$	index in the $y$ dimension;	$t$	elapsed time of the pumping period ( $t$ );
$j$	index in the $x$ direction;	$T$	transmissivity ( $L^2t^{-1}$ );
$k$	time index;	$T_L$	transient leakage coefficient ( $t^{-1}$ );
$K_{xx}, K_{yy}$	principal components of the hydraulic conductivity tensor ( $Lt^{-1}$ );	$T_{xx}, T_{xy}, T_{yx}, T_{yy}$	components of the transmissivity tensor ( $L^2t^{-1}$ );
$K'$	hydraulic conductivity of the confining bed ( $L/t$ );	$\bar{U}$	upper triangular factor of $(\bar{A} + \bar{B})$ ;
$l$	iteration parameter index;	$\bar{V}$	intermediate vector in SIP algorithm;
$L$	number of iteration parameters in a cycle;	$W(x,y,t)$	volume flux per unit area ( $Lt^{-1}$ );
$\bar{L}$	lower triangular factor of $(\bar{A} + \bar{B})$ ;	$a$	row correction for LSOR;
$m$	thickness of the confining bed ( $L$ );	$\alpha, \beta, \gamma, \delta, \eta$	elements of factors of $(\bar{A} + \bar{B})$ ;
$M$	vector of ADI parameters;	$\beta$	parameter in Hantush (1960) solution;
$n$	iteration index;	$\hat{\beta}$	column correction for LSOR;
$N_s$	number of arrays required for the options;	$\beta'$	iteration parameter for SIP;
$N_r$	number of nodes in a row;	$\gamma$	constant used in calculating ADI parameters;
$N_c$	number of nodes in a column;	$\delta_x, \delta_y$	normalized grid spacing;
$q'$	flux from a confining bed ( $Lt^{-1}$ );	$\Delta h$	head change between adjacent nodes ( $L$ );
		$\Delta t$	time increment ( $t$ );
		$\Delta x$	space increment in the $x$ direction ( $L$ );
		$\Delta y$	space increment in the $y$ direction ( $L$ );
		$\epsilon$	closure criterion ( $L$ );



$\bar{y}$	vector of change in head over an iteration;
$\rho(G)$	spectral radius of Gauss-Seidel iteration matrix;
$\phi_1, \phi_2$	constants in definition of coefficients of $(\bar{A} + \bar{B})$ ;
$\omega$	acceleration parameter;
$\omega_i$	iteration parameter;
$\omega_{max}$	maximum iteration parameter;
$\omega_{min}$	minimum iteration parameter;
$\omega_{opt}$	optimum acceleration parameter.

## Attachment II, Computer Program

### Main program

The first function of the main program is to dimension the arrays for the field problem being simulated. The algorithm allocates storage space reserved in a vector, Y. Some arrays are required for every simulation; others are needed only if certain options are specified. The information needed to allocate space to the arrays is contained in the Group I data cards which are read by the main program (see Attachment III).

Once the model is compiled, it does not need to be recompiled for a new field problem unless (1) the logic is changed or (2) the vector Y is not dimensioned large enough for the new problem. The minimum dimension of the vector Y (YDIM) can be computed by

$$YDIM \cong (15 + N_a) N_x N_y \quad (39)$$

in which  $N_a$  is the total number of arrays required for the options (from table 2).

Equation 39 is approximate, but normally will give a value that is sufficient for the simulation. The exact dimension required is

Table 2.—Number of arrays required for the options

Option	Number of arrays
Water Table -----	3
Conversion <sup>1</sup> -----	1
Leakage -----	3
Evapotranspiration -----	1
SIP -----	4

<sup>1</sup> Conversion also requires the arrays for the water table option.

printed on the first page of the output as 'WORDS OF VECTOR Y USED=XXXX'.

In the second part of the main program, the location of the initial addresses of the arrays are passed to the subroutines. (See table 3 for details.) The variables in table 3 defining the dimensions of the arrays are defined in Attachment VI; the first four arrays and XII are double precision.

The last part of the main program controls the sequence of computations illustrated by the generalized flow chart (Appendix V). In the flow chart, the routines are lettered in sequence starting with the main program. Entry points for the routines are numbered in sequence along the left side of the chart. Exits from a routine are indicated by circles containing the entry point of the routine to which control passes. A break occurs in the flow chart following an unconditional exit. Variables used in the flow chart are defined in Attachment VI.

### Subroutine DATAI

Instructions for the preparation of the data deck are given in Attachment III. Data may be input to the model in any consistent set of units in which second is the time unit. It is organized into four groups: Data in groups I and II are the simulation options and scalar parameters: group III cards are used to initialize the arrays. These three groups are required for each new simulation. Group IV contains data that varies with each new pumping period. The program permits changing well discharge and the time parameters each pumping period, but the program can be modified to read other data (for example, recharge rate) with this set of cards.

### Time parameters

The time parameters include the initial time step, DELT; a multiplication factor for increasing the size of the time step, CDLT; the number of time steps, NUMT; and the simulation period, TMAX. Since the rate of water-level decline decreases during a pumping period, the time step is increased by the factor CDLT each step (commonly 1.5). For

Table 3.—Arrays passed to the subroutines and their relative location in the vector Y

Array	Sequence number in vector Y	Subroutine						Dimensions
		DATAI	STEP	SOLVEI	COEF	CHECKI	PRNTAI	
PHI -----	1	×	×	×	×	×	×	(IZ, JZ) '8
BE -----	2	--	--	×	--	--	--	IMAX '8
G -----	3	--	--	×	--	--	--	IMAX '8
TEMP -----	4	--	--	×	--	--	--	IMAX '8
KEEP -----	5	--	×	×	×	×	--	IZ, JZ
PHE -----	6	--	--	×	×	×	--	IZ, JZ
STRT -----	7	×	×	×	×	×	--	IZ, JZ
SURI -----	8	×	×	--	×	--	×	IZ, JZ
T -----	9	×	×	×	×	×	--	IZ, JZ
TR -----	10	×	--	--	×	×	--	IZ, JZ
TC -----	11	×	--	--	×	×	--	IZ, JZ
S -----	12	×	--	×	×	×	×	IZ, JZ
QRE -----	13	×	--	×	--	×	--	IZ, JZ
WELL -----	14	×	×	×	×	×	×	IZ, JZ
TL -----	15	×	--	×	×	×	--	IZ, JZ
SL -----	16	×	--	×	×	--	--	IZ, JZ
PERM -----	17	×	×	--	×	×	--	IP, JP
BOTTOM -----	18	×	×	--	×	×	--	IP, JP
SY -----	19	×	--	--	×	×	--	IP, JP
RATE -----	20	×	--	--	×	×	--	IR, JR
RIVER -----	21	×	--	--	×	×	--	IR, JR
M -----	22	×	--	--	×	×	--	IR, JR
TOP -----	23	×	×	--	×	×	--	IC, JC
GRND -----	24	×	--	--	×	×	--	IL, JL
DEL -----	25	--	--	×	--	--	--	IS, JS
ETA -----	26	--	--	×	--	--	--	IS, JS
V -----	27	--	--	×	--	--	--	IS, JS
XI -----	28	--	--	×	--	--	--	IS, JS
DELX -----	29	×	×	×	×	×	×	JZ
DDN -----	30	--	×	--	--	--	--	JZ
BETA -----	31	--	--	×	--	--	--	JZ
DELY -----	32	×	×	×	×	×	×	IZ
ALFA -----	33	--	--	×	--	--	--	IZ
WR -----	34	×	×	--	--	--	--	IH
NWR -----	35	×	×	--	--	--	--	IH, 2
XII -----	36	--	--	×	--	--	--	IMAX '8
TEST 3 -----	37	--	×	×	--	--	--	IMX1

any time step ( $k$ ) the time increment is given by

$$\text{DEL}T_k = \text{CDLT} * \text{DEL}T_{k-1}.$$

$\text{DEL}T_0$  is the time step recorded on the data card.

The program has two options for selecting the time parameters:

1. To simulate a given period of time, select CDLT and an appropriate  $\text{DEL}T_0$ , and set NUMT greater than the expected number of time steps. The program computes the required initial  $\text{DEL}T_0$  (which will not exceed the value of

$\text{DEL}T_0$  coded on card 1 of group IV) and NUMT to arrive exactly at TMAX on the final time step. In a simulation of one pumping period in which results are required at several specific times, the simulation can be broken into several "pumping periods." Each period will have the same pumpage, and TMAX is used to specify the appropriate times for display of results.

2. To simulate a given number of the time steps, set TMAX greater than the expected simulation period and the program will use  $\text{DEL}T_0$ , CDLT, and

NUMT as specified on the time parameter card.

To minimize the error due to approximation of the time derivative, several time steps should be simulated before the first step at which results are displayed. This suggestion should be followed unless the system is nearly steady-state before the results are needed. In this case a one-step simulation may be satisfactory, but this approach should be checked by making one run as a multistep simulation so that the results can be compared.

For steady-state simulations, set the storage coefficient and (or) specific yield of the aquifer and the specific storage of the confining bed to zero. Compute for one time step of any length (for example, set TMAX=1, NUMT=1, CDLT=1, DELT=24) and the program should iterate to a solution. The maximum permitted number of iterations (ITMAX) should be larger for steady-state than for transient simulations. If the calculations do not converge to a solution within a reasonable number of iterations, it may be necessary to use a transient simulation for enough steps to attain steady state (see also the discussion of ADI iteration parameters) or use another numerical technique.

#### Initialization

In addition to reading data and computing the time parameter, this routine initializes other arrays and scalar parameters. In particular, note that the leakage coefficient,  $TL$ , will equal  $K'_{i,j}/m_{i,j}$  and can be computed once for the entire simulation if the specific storage of the confining bed is zero. The computation of the steady leakage term,  $SL$ , and the division of well discharge by the area of the cell need to be done only once for each pumping period. At the beginning of each pumping period the starting head (STRT) and the simulation time (SUMP) used in computing transient leakage are initialized.

#### Subroutine STEP

Subroutine STEP initializes variables for a new time step, checks for steady-state conditions after a solution is obtained for the

time step, and controls the printing and punching of results and the writing of results on disk. If head values are punched at the end of the simulation or are written on disk, they can be used to extend the simulation or as input to plotting routines. (See the program by Cosner and Horwich, 1974.) Currently, a general program is being written to display results in various forms on the line printer and plotters; it is described in detail in another section of this report.

In the check for steady state during transient simulations, the head change over a time step is computed. If the absolute value of change at all nodes is less than EROR, the message 'STEADY STATE AT TIME STEP X' is printed. The program then prints all desired output for the final time step (X) and proceeds to read data for the next pumping period, if any.

#### Maximum head change for each iteration

The printed results are explained in the section on theory and in the discussions of subroutines COEF, CHECKI, and PRNTAI or are self explanatory, except for the listing of the absolute value of the maximum head change for each iteration. This information is useful if convergence is slow with ADI or SIP because it may indicate that a slightly larger error criterion will give a satisfactory solution with considerably fewer iterations.

#### Subroutine SOLVE

The three SOLVE routines, SOLVE1, SOLVE2, and SOLVE3 are, respectively, SIP, LSOR and ADI. They have been described in previous sections, but a few additional comments are necessary.

In these routines and in subroutine COEF, the usual (I,J) notation has been replaced in favor of single-subscript notation. Less time is involved in finding the value of a variable with a single subscript than in finding the value of one with a double subscript and, as a consequence, computational efficiency is improved. The five variables used as subscripts in this notation are defined in Attachment VI.

### SIP iteration parameters

The algorithm in ITER1 permits computation of the iteration parameters in increasing or decreasing order and repeat of parameters depending on the initialization of the vector IORDER. Note that LENGTH is twice the number of different parameters and that the DATA statement that initializes IORDER assumes LENGTH=10. Replace the DATA statement with a READ statement if additional flexibility is desired in choosing the order of parameters without recompiling the subroutine.

### Exceeding permitted iterations

If the permitted number of iterations for a time step is exceeded, the message 'EXCEEDED PERMITTED NUMBER OF ITERATIONS' is printed. Following the message the mass balance, head matrix, etc., as specified in the options are printed for the final iteration. This information is useful in determining the cause of the nonconvergence. Before terminating the run, the mass balance and head values will be punched if PUNC was specified in the options or written on disk if IDK2 was specified. With punched output or results on disk, the user has the option to extend the number of iterations if it appears that a solution can be obtained. If iterations are exceeded on the first time step, the head values saved (punched or written on disk) were computed in the last iteration. If iterations are exceeded on a subsequent time step, KT, the head values and mass-balance parameters saved are the results for time step KT-1.

### Subroutine COEF

Most of the calculations for coefficients used in the solution of the numerical schemes are done in this routine. The more extensive computations except those described in the section on theory are discussed in the following paragraphs.

#### Transient leakage coefficients

The algorithm for the transient parts of equations 9 and 10 is the same except for two

conditional statements that recompute PPT and DENOM if dimensionless time is in the range for applying equation 9. In performing the infinite summation, the code checks for the significance of additional terms, but in any case limits the summation to a maximum of 200 terms. The minimum and maximum values of dimensionless time, TMIN and TT, are retained and printed with the results for the time step so that the user will know whether or not transient leakage effects are significant.

#### Transmissivity as a function of head

The transmissivity for water-table or combined water-table-artesian aquifers is computed as a function of the saturated thickness of the aquifer. If a cell (except a cell with well discharge) goes dry, a message 'NODE I, J GOES DRY' is printed, the transmissivity for the cell is set to zero, and the head is set to the initial surface (so that the location of the cell will show up in the output). No provision is made to permit the cell to resaturate in subsequent pumping periods because the additional code necessary to accommodate this special situation is not warranted in a general program.

When a cell with well discharge goes dry (that is, a hypothetical well with radius  $r_e$  goes dry), the program terminates the computation with printed output, and, if specified in the options, saves the results. Printed output is headed by 'WELL I, J GOES DRY' followed by drawdown when the well went dry. If results for the previous time step were not printed, drawdown and a mass balance (if specified in the program options) for the previous time step are printed. Finally, if specified in the options, mass-balance parameters and head values for the previous time step are punched or written on disk so that the user has the option of continuing the simulation after modifying the well discharge.

#### TR and TC coefficients

The TR and TC arrays save values that are used repeatedly in the algorithm. They are computed once for artesian problems and

each iteration for water-table and combined artesian-water-table simulations. TR (I,J) is the harmonic mean of  $T_{xx}(I,J)/DELX(J)$ ,  $T_{xx}(I,J+1)/DELX(J+1)$ ; TC (I,J) is the harmonic mean of  $T_{yy}(I,J)/DELY(I)$ ,  $T_{yy}(I+1,J)/DELY(I+1)$ .

### Subroutine CHECKI

A mass balance is computed in this routine. The results are expressed in two ways: (1) as a cumulative volume of water from each source and each type of discharge and (2) as rates for the current time step.

In the cumulative mass balance, storage is treated as a source of water. Flow to and from constant-head boundaries is computed with Darcy's law using the gradients from constant-head nodes to adjacent nodes inside the aquifer. Other computations in the algorithm are self explanatory.

The difference between the sum of sources and sum of discharges from the system is usually less than 1 percent. A larger error, however, does not necessarily mean that the results are poor; it may be due to lack of precision in calculating the mass balance. This has been observed, for example, if a leaky streambed is given a large  $K'/m$  ratio so that it is effectively a constant-head boundary. The leakage computation is inaccurate if the head values at a stream node are identical to 6 or 7 significant figures and they are stored as single precision variables.

To the right of the cumulative mass balance are printed the flow rates for the current time step. They are self explanatory except for leakage. "Leakage from previous pumping period" is the leakage resulting from gradients across the confining bed at the start of the current pumping period. The "total" leakage is the sum of leakage due to the initial gradients plus leakage induced by head changes during the current pumping period.

### Subroutine PRNTAI

This routine prints a map of drawdown and hydraulic head. Up to three characters

are plotted for each cell with the rightmost character as close to the location of the node as the printer will allow. An option to permit the printing of results at different scales in the  $x$  and  $y$  dimensions is useful for cross sections. This routine is useful for displaying results during calibration runs. More elegant graphical displays for final results are described in another section.

The user specifies XSCALE and YSCALE, the multiplication factors required to change from units used in the model to units used on the map; DINCH, the number of map units per inch; FACT1 and FACT2, the multiplication factors for adjusting the values of drawdown and head to be plotted, respectively; and MESUR, the name of the unit used on the map. As an example, assume that the length unit used in the model is feet, the map is to be scaled at 3 miles per inch and drawdown values at 1 foot increments and head values at 10 foot increments are to be plotted. Then XSCALE=YSCALE=5280, DINCH=3, FACT1=1, FACT2=0.1; and MESUR=MILES.

To print a map of maximum possible size, number the rows in the short dimension to take advantage of the orientation of the map on the computer page where the X direction is vertical and the Y direction horizontal. (See fig. 26.) The origin is the upper left-hand corner of the block for row 2, column 2. Orienting the map with the origin in the upper left-hand corner, the right and bottom sides of the map include the node locations for the second to last column and row, respectively. The border is located to the nearest inch outside these node locations and may or may not fall on the cell boundaries depending on the scaling. The map is automatically centered on the page and is limited to a maximum of 12 inches (300 mm) in the Y direction. If the parameters for a map are specified such that the Y dimension is more than 12 inches (300 mm) adjustments are automatically made to fit the map within this limit. A common mistake is to specify a value for Y scale that is less than 1.0. This generates the message 'NOTE: GENERALLY SCALE SHOULD BE >OR = 1.0,' and a suit-

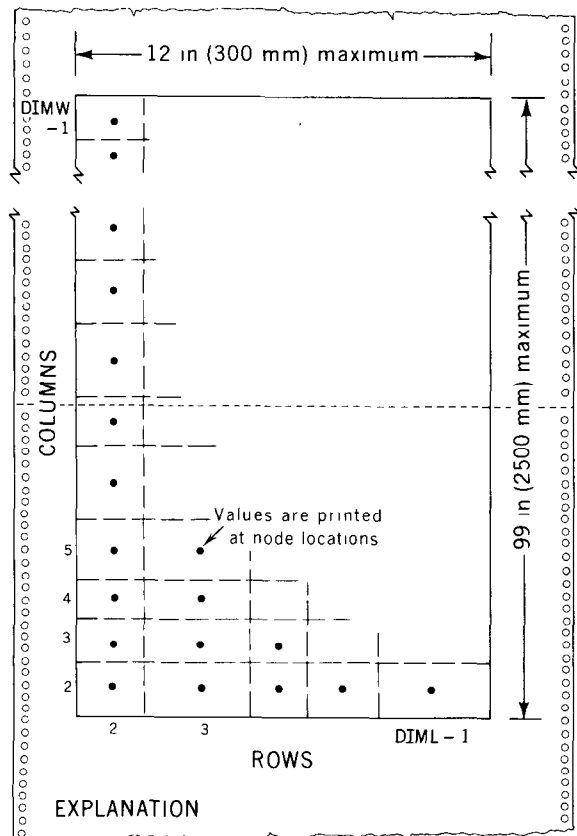


FIGURE 26.—Orientation of map on computer page.

able adjustment is made to DINCH. In the X direction, the map is limited only by the dimension of the NX vector. (For example, when the dimension of NX is 100, the map is limited in the X direction to  $100-1=99$  inches (2500 mm).) Several parameters (PRNT, BLANK, N1, N2, N3, and XN1) are initialized in the BLOCK DATA routine to values that assume the line printer prints 6 lines per inch, 10 characters per inch, and 132 characters per line. These parameter values may need to be changed for a line printer with other specifications.

The PRNTAI subroutine can be modified to cycle a set of alphameric symbols for drawdown. If this type of map is desired, remove the C from column 1 of statements PRN1060 and PRN1230. This will cycle the symbols 1,2,3,4,5,6,7,8,9,0 for drawdown. To plot a different set of symbols will require modification of the initialization of SYM in BLOCK DATA. To cycle more than 10 sym-

bols will require more extensive changes to the initialization of SYM and modifications to the code in ENTRY PRNTA.

## BLOCK DATA routine

The BLOCK DATA routine initializes scalar parameters and arrays used in PRNTAI and other subroutines. The unit numbers for card reader, line printer, and card punch are commonly 5, 6 and 7, respectively. At computer installations where other numbers are used, change the initialization of P, R, and PU.

## Technical information

### Storage requirements

Using the FORTRAN G, Level 21 compiler, the source code and fixed-dimension arrays require 100K bytes of memory (88K bytes if only one SOLVE routine is compiled). The storage requirements including all options but not including storage requirements for reading and writing on disk are  $(100+X/256)$  K bytes where X is the dimension of the vector Y in the main program. Subtract 14K bytes from the values if the FORTRAN H, OPT=2 compiler is used. The FORTRAN G compile step requires 120K bytes of memory and the FORTRAN H, OPT=2 compiler requires 218K bytes of memory.

### Computation time

Computation time is a function of so many variables that no general rule can be stated. For example, the simulation of a nonlinear water-table problem requires many more computations per time step than does the simulation of a linear artesian-aquifer problem.

As an example, the simulation of a linear aquifer system (problem 2) with a grid of  $25 \times 38$  required 45 seconds for 40 iterations with the program compiled under FORTRAN G. This is about 0.002 seconds for each node inside the aquifer each iteration on the IBM 370/155. A significant reduction (about 1/3)

in execution time can be achieved by using the FORTRAN H compiler which generates a more efficient code than the FORTRAN G compiler.

Further significant reductions in execution time can be achieved if the model is designed for a specific problem. Problem 3, for example, does not require computation of leakage, storage, or evapotranspiration terms.

#### Use of disk facilities for storage of array data and interim results

In an effort to expedite use of the program on remote terminals connected to the IBM 370/155, options are included to utilize disk storage facilities. These options enable storage and retrieval of array data (STRT, PERM, and so forth) and the saving of interim head values without punching them on cards.

Use of these options can be particularly beneficial at remote terminals with low speed data transmission or without punch output capability. Also, the type of read statements used afford more efficient data transmission from disk than from cards.

*Storage of array data* is accomplished via a direct access data set that is defined by a DEFINE FILE statement in the main program (card MAN0480) and by a DD statement in the JCL string used to execute the program. To establish the data set, the DEFINE FILE statement and the DD statement must indicate the amount of space that is required. The DEFINE FILE statement takes the following form:

```
DEFINE FILE 2(14,???,U,KKK)
                                MAN0480
```

where ??? is the number of nodes for the problem being solved (DIMLxDIMW). Parameters U and KKK are indicators and do not vary.

The DD statement contains information, such as account number, that will be different for each user. Also, the first reference to the data set is somewhat different from subsequent references. To utilize one of the disk packs provided by the system (IBM 370/

155) for semipermanent storage of user data, the first reference to the data set will take the following general form if the FORTGCG procedure is used to compile and execute the program.

```
//GØ. FT02F001 DD DSN=Azzzzzz.AZbbb.
                                cxxwwwww.aaaaaaa,
//   UNIT=ØNLINE,DISP=(NEW,
                                KEEP),
//   SPACE=(????,(14)),DCB=
                                (RECFM=F)
```

where

zzzzzz	are the first six digits of a nine digit account number;
bbb	are the last three digits of a nine digit account number;
c	is the center code (same as column 3 on job card);
xx	is the two digit organization code (same as columns 4 and 5 on job card);
wwwww	is the four or five digit program number (same as the program number beginning in column 24 of the job card);
aaaaaaa	is any 1 to 8 character name used to designate the name of the data set;
????	is the number of bytes per record that are to be reserved and should be set equal to DIMLxDIMWx4.

The instructions for the DSN parameter are also given in the CCD users manual, chapter 5, pages 3 and 4. When this initial allocation is processed the system will indicate in the HASP system log, JCL string output, the volume on which the data set was established (for example, SYS011 or SYS015). Subsequent use of the data set must indicate this information by modifying the underlined parameters in the initial reference to the data set. Thus the DD statement will read:

```
//GØ. FT02F001 DD DSN=Azzzzzz.Azbbb.
                                cxxwwwww.aaaaaaa,
```

```
// UNIT=ØNLINE, DISP=SHR,VØL
   =SER=yyyyyy
```

where the DSN parameter is the same as the initial run and yyyyyy indicates the volume (for example, SYS015) on which the data set was established by the initial run. The individual data arrays that are to be stored and later retrieved from this data set are specified on the parameter card for each array. These specifications will be discussed completely in the section on Data Deck Instructions (Attachment III).

If use of this option is selected, space for buffers must be reserved via the REGION parameter on the EXEC card. The amount of space needed is approximately equal to two times the number of bytes per record (indicated in the SPACE parameter on the DD card defined above).

*Interim results* (head values, cumulative simulation time, and mass-balance parameters) can be punched on cards or can be stored and retrieved from data sets on disk in much the same manner as array data. Use of storage on disk is initiated by parameters on the simulation options card. (See attachment III, card 3.)

Definition of the sequential data set on disk where the information will be stored is accomplished by a DD statement in the JCL string used to execute the program. If one of the system disk packs is used to store the data set, the first reference to the data will be different from subsequent references as in the case of array data sets. The first reference will take the following form if the FORTGCG procedure is used.

```
//GØ. FT04F001 DD DSN=Azzzzz.AZbbb.
   cxxwwwww.aaaaaaaa,
// UNIT=ØNLINE, DISP=(NEW,
   KEEP),SPACE=(TRK,(1,1),
   RLSE),
// DCB=(RECFM=VBS,LRECL
   =dddd, BLKSIZE=eeee)
```

The DSN parameter is defined in the same manner as previously discussed for the direct access (array) data sets and:

dddd—equals DIMLxDIMWx8 + 48 ( $\leq 6440$ )  
 eeee—equals DIMLxDIMWx8 + 52 ( $\leq 6440$ )

If BLKSIZE (eeee) exceeds 6444, code 6444 for (eeee) and 6440 for (dddd). Also, additional core equal to about two times the value of BLKSIZE must be reserved for buffers via the REGION parameter on the EXEC card.

Once the initial reference to the data set has been successfully processed, the system will indicate (via the JCL printout) on what volume the data set has been established (for example, SYS011 or SYS015) and, subsequent references to the data set will appear as follows:

```
//GØ. FT4F001 DD DSN=Azzzzz.AZbbb.
   cxxwwwww.aaaaaaaa,
// UNIT=ØNLINE, VØL=SER=
   yyyyyy,DISP=SHR
```

where yyyyyy is the name of the disk pack (for example, SYS011) that contains the data set and DSN is as previously described.

To destroy (erase) an array data set or an interim results data set, simply execute the following job.

```
// EXEC PGM=IEFBR14
//X DD DSN=Azzzzz.AZbbb.cxxwwwww.
   aaaaaaaaa,
// UNIT=ØNLINE, VØL=SER=yyyyyy,
   DISP=(ØLD,DELETE)
```

Use of the disk facilities is illustrated in Appendix IV.

## Graphical display package

A series of computer programs are currently being written and assembled that will enable graphical display of results of computer models. Components of this graphical display package will include:

1. time-series plots of model results on the printer,
2. time-series plots on pen plotters (CALCOMP),
3. contour maps of model results at selected time steps on the printer,
4. contour maps utilizing pen plotters, and
5. other graphical displays, such as perspective (three-dimensional) drawings.



The FORTRAN code shown in figure 27 can be inserted into the program to produce output that can be used in the graphical display package. The changes to MAIN and STEP are required after statements MAN2600 and STP1000, respectively. Statement MAN2600 is deleted. In subroutine PRNTAI, the REAL\*8 specification and the DIMENSION statement must be added and the remaining code inserted after statement PRN1650. Also, unit numbers 10 and 11 must be specified on DD statements when the program is executed. Unit 10 is used only for temporary storage and the following DD statement will generally suffice.

```
//GØ. FT10F001 DD DSN=&&DATA,DISP
    = (NEW,DELETE),UNIT
    = ØNLINE,
//    SPACE = (TRK, (10,5)),DCB =
    (RECFM = VBS,LRECL = 6440,
    BLKSIZE = 6444)
```

Unit 11 points to the data set that is used to store the data required by the graphical display package and must be semipermanent in nature. That is, it must not be deleted upon completion of your job. The DD statement will generally take the following form.

```
//GØ. FT11F001 DD DSN = Azzzzzz.AZbbb.
    cxwwwww.aaaaaaa,
//    DISP = (NEW,KEEP),UNIT =
    ØNLINE,SPACE = (TRK, (10,5),
    RLSE),
//    DCB = (RECFM = VBS,LRECL = 6440,
    BLKSIZE = 6444)
```

The data set name parameter (DSN) was discussed in the previous section. The SPACE and DCB parameters shown above should generally be adequate. Recall that once the data set is established, it will be assigned to a certain volume (disk pack) by the IBM operating system. Subsequent references to the data set must include this volume number in the DD statement, that is, VØL = SER = ??.

Results of using a preliminary version of the graphical display package are shown in figures 28 and 29. The time-series plot shown in figure 28 was made on the line printer and the contour map shown in figure

29 was made on a CALCOMP plotter. Documentation on the use of the graphical display package is currently being written.

## Modification of program logic

Some users may wish to compile only one or two numerical options with the program. This is done by removing the SOLVE routine(s) not needed from the card deck and modifying the main program in either of the following ways, assuming for this example that SIP is being removed: (1) remove the three IF statements that call SOLVE1, ITER1, and NEWITA, or (2) punch a C in column 1 of these statements and leave them in the main program.

Other modifications to the program logic will be required for certain applications. Modifications will range from changing a few statements to adding a subroutine or deleting options not used. In any case the changes should be made by a programmer familiar with the computational scheme because almost any change has an unanticipated effect on another part of the program requiring several debugging runs.

Reasonably simple modifications to the program include changing format statements and shifting data sets (for example, recharge rate) from GROUP III to GROUP IV so they can be modified for each pumping period.

Adding a second confining bed would be a more complex modification because it may require additional arrays, and ENTRY CLAY in subroutine COEF would have to be made general to accept confining-bed parameters for either bed.

## FORTRAN IV

The program includes several FORTRAN IV features that are not in ANS FORTRAN (for example, ENTRY, END parameter in read statement, mixed-mode expressions, G format code, literal enclosed in apostrophes). If the program is used at a computer center where the FORTRAN compiler does not include these extensions, programmers at the

## MAIN

```

300 CALL GRAPH (Y(L(3)),Y(L(4)),Y(L(5)),Y(L(6)),Y(L(7)),
1 YY(1),Y(L(8)))
READ (R, 320,END=310) NEXT

```

## STEP

```

WRITE(10) PHI,SUM

```

## PRINTAI

```

REAL*8 HD

DIMENSION NN(1),SUMX(1),SUMY(1),X(1),Y(1),ZZ(1),HD(1)

C*****
ENTRY GRAPH (SUMX,SUMY,X,Y,ZZ,HD,NN)
C*****
C COMPUTE X AND Y COORDINATES OF ROWS AND COLUMNS
SUMX(1)=DELX(1)/2.
SUMY(1)=DELY(1)/2.
DO 325 I=2,DIML
325 SUMY(I)=SUMY(I-1)+(DELY(I)+DELY(I-1))/2.
DO 330 I=2,DIMW
330 SUMX(I)=SUMX(I-1)+(DELX(I)+DELX(I-1))/2.
C DETERMINE NUMBER OF ACTIVE NODES, THEIR STORAGE LOCATION,
C AND THEIR X AND Y COORDINATES
N=0
DO 340 I=2,IN01
DO 340 J=2,JN01
IF (T(I,J).EQ.0.) GO TO 340
N=N+1
NN(N)=I+DIML*(J-1)
X(N)=SUMX(J)
Y(N)=SUMY(I)
340 CONTINUE
C WRITE X AND Y COORDINATES ON UNIT 11
WRITE(11) (X(I),I=1,N)
WRITE(11) (Y(I),I=1,N)
C REWIND UNIT 10 AND REPROCESS PHI MATRIX AT EACH TIME STEP
C PLACING PHI VALUES AT ACTIVE NODES IN THE ZZ ARRAY (REAL*4)
REWIND 10
DO 380 I=1,KT
READ(10) PHI,SUM
DO 350 J=1,N
NIJ=NN(J)
350 ZZ(J)=HD(NIJ)
C WRITE PHI VALUES AT ACTIVE NODES AND ELAPSED SIMULATION TIME
C ON UNIT 11
WRITE(11) (ZZ(J),J=1,N),SUM
380 CONTINUE
WRITE(6,390) N,KT,SUMX(DIMW),SUMY(DIML)
390 FORMAT(/,' GRAPHICS OUTPUT FOR ',I6,' ACTIVE NODES AND ',I4,
1 ' TIME STEPS HAS BEEN WRITTEN ON UNIT 11',/,
2 ' MAXIMUM X,Y COORDINATE PAIR IS ',F10.2,' ',F10.2)
RETURN

```

FIGURE 27.—Additional FORTRAN code required to produce output for graphical display.

```

SAMPLE GRAPHICS OUTPUT
X,Y IS (11475. 1350.0
NML = 5
NSBH = 6
NVL = 12
NSBV = 10
NSCALE = 1
NSCALE = 0
NSCALE = 2
NSCALE = 0
NSCALE = 0
YMAX=104.00
YMTA =92.000
XMAX =.0
XMIN =.18000F 06
1 REPRESENTS X,Y COORDINATES (11475. 1350.0 )
2 REPRESENTS X,Y COORDINATES (3825.0 2750.0 )
3 REPRESENTS X,Y COORDINATES (11475. 1950.0 )
4 REPRESENTS X,Y COORDINATES (11475. 625.00 )

```

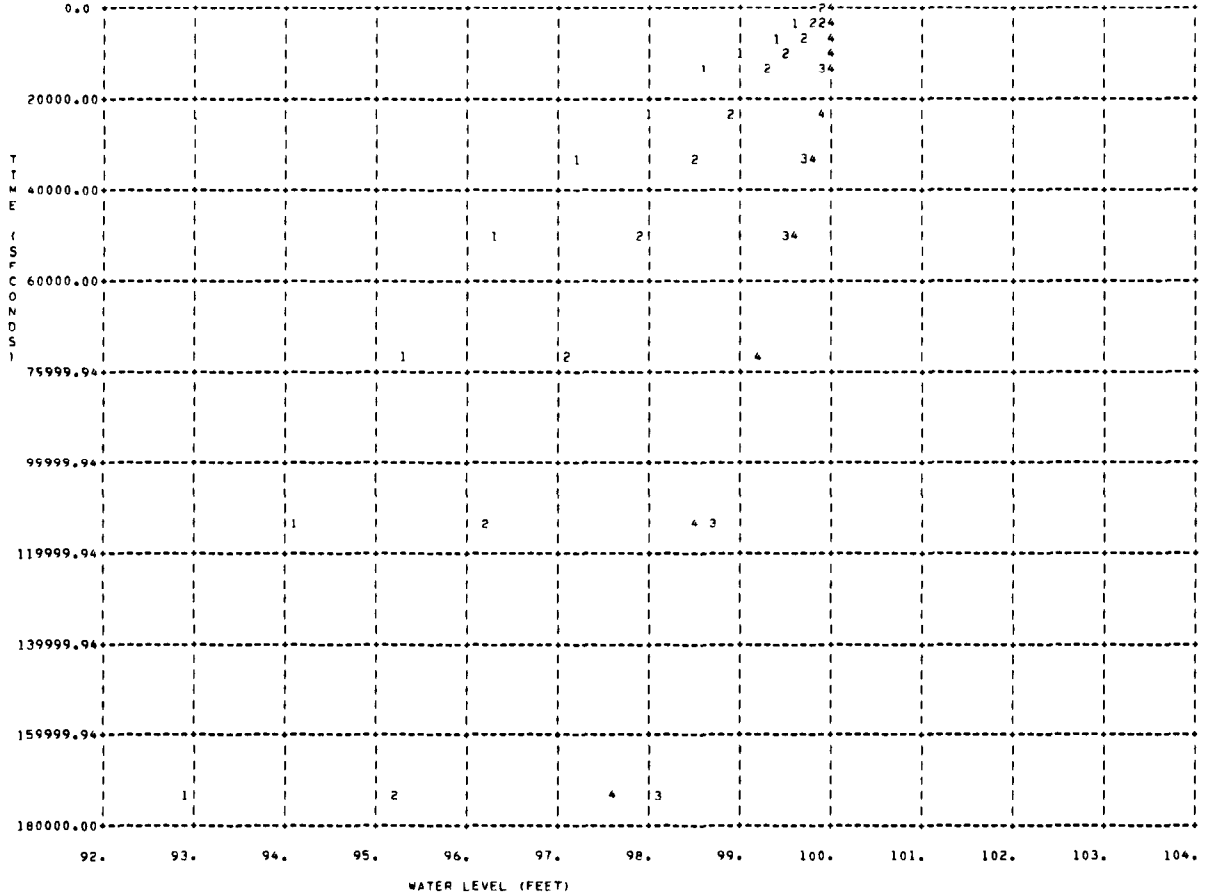


FIGURE 28.—Water level versus time at various nodes of the sample aquifer problem produced by the graphical display package.

selected installation may be available to modify the computer code as necessary.

### Limitations of program

The model documented in this report is reasonably free of errors and has been used successfully to simulate a variety of aquifer systems in two dimensions. Undiscovered er-

rors in the logic, however, may appear as the model is applied to a variety of new problems.

The user is cautioned against using this model to make more than a crude simulation of three-dimensional problems. A rigorous analysis of three-dimensional aquifer systems can be made only with the appropriate analog or digital simulators.

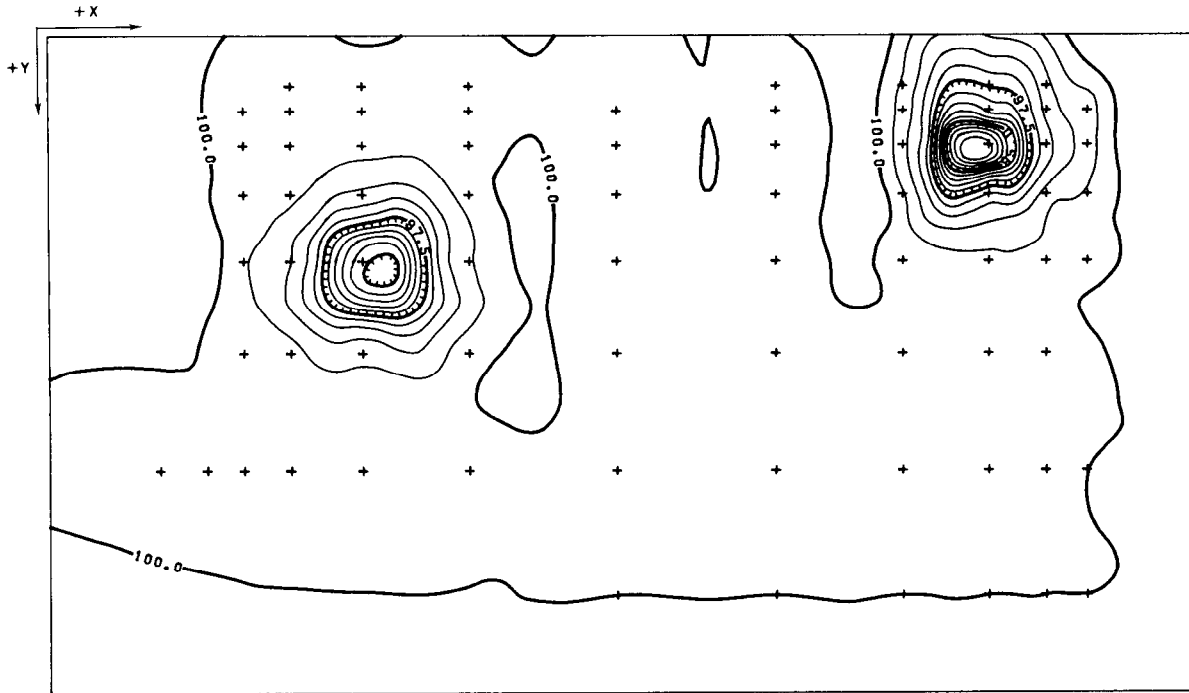


FIGURE 29.—Contour map of water level (in feet) for sample aquifer problem produced by graphical display package. Contour interval is 0.5 ft.

## Attachment III

### Data Deck Instructions

#### Group I: Title, simulation options, and problem dimensions

This group of cards, which are read by the main program, contains data required to dimension the model. To specify an option on card 3, punch the characters underlined in the definition, starting in the first column of the field. For any option not used, leave the appropriate columns blank.

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
1	1-80	20A4	HEADNG	Any title the user wishes to print on one line at the start of output.
2	1-48	12A4		
3	1-5	A4,1X		
	6-10	A4,1X	LEAK	<u>LEAK</u> for an aquifer system including leakage from a stream or confining bed.
	11-15	A4,1X	CONVRT	<u>CONV</u> for combined artesian-water-table aquifer.
	16-20	A4,1X	EVAP	<u>EVAP</u> to permit discharge by evapotranspiration.
	21-25	A4,1X	RECH	<u>RECH</u> to include a constant recharge rate.

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
	26-30	A4,1X	NUMS	<u>SIP</u> or <u>LSOR</u> or <u>ADI</u> to designate the equation-solving scheme.
	31-35	A4,1X	CHCK	<u>CHEC</u> to compute a mass balance.
	36-40	A4,1X	PNCH	<u>PUNC</u> for punched output at the end of the simulation.
	41-45	A4,1X	IDK1	<u>DK1</u> to read initial head and mass balance parameters from disk (unit 4).
	46-50	A4,1X	IDK2	<u>DK2</u> to save (write) computed head, elapsed time, and mass balance parameters on disk (unit 4).
	51-55	A4,1X	NUM	<u>NUME</u> to print drawdown in numeric form.
	56-60	A4,1X	HEAD	<u>HEAD</u> to print the head matrix.
(All variables on card 4 are integers)				
4	1-10	I10	DIML	Number of rows.
	11-20	I10	DIMW	Number of columns.
	21-30	I10	NW	Number of pumping wells for which drawdown is to be computed at a "real" well radius.
	31-40	I10	ITMAX	Maximum number of iterations per time step.

NOTE.—Steady-state simulations often require more than 50 iterations. Transient time steps usually require less than 30 iterations.

#### Group II: Scalar parameters

The parameters required in every problem are underlined. The other parameters are required as noted; when not required, their location on the card can be left blank. The G format is used to read E, F and I data. Minimize mistakes by always right-justifying data in the field. If F format data do not contain significant figures to the right of the decimal point, the decimal point can be omitted. *Default typing of variables applies.*

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
1	1-4	A4	CONTR	<u>CONT</u> to generate a map of drawdown and (or) hydraulic head; <i>for no maps</i> insert a blank card.
	11-20	G10.0	XSCALE	Factor to convert model length unit to unit used in X direction on maps (that is, to convert from feet to miles, XSCALE=5280).
	21-30	G10.0	YSCALE	Factor to convert model length unit to unit used in Y direction on maps.
	31-40	G10.0	DINCH	Number of map units per inch.
	41-50	G10.0	FACT1	Factor to adjust value of drawdown printed*.
	51-60	G10.0	FACT2	Factor to adjust value of head printed*.

*Value of drawdown or head	FACT 1 or FACT 2	Printed value
52.57	.01	0
	.1	5
	1	52
	10	525
	100	***

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
	61-68	A8	MESUR	Name of map length unit.
2	1-10	G10.0	<u>NPER</u>	Number of pumping periods for this simulation.
	11-20	G10.0	<u>KTH</u>	Number of time steps between printouts.
NOTE.—To print only the results for the final time step in a pumping period, make KTH greater than the expected number of time steps. The program always prints the results for the final time step.				
	21-30	G10.0	<u>ERR</u>	Error criterion for closure ( $L$ ).
NOTE.—When the head change at all nodes on subsequent iterations is less than this value (for example, 0.01 foot), the program has reached a solution for the time step.				
	31-40	G10.0	EROR	Steady-state error criterion ( $L$ ).
NOTE.—If the head change between time steps in transient simulations is less than this amount, the pumping period is terminated.				
	41-50	G10.0	SS	Specific storage of confining bed ( $1/L$ ).
NOTE.—SS has a finite value only in transient simulations where leakage is a function of storage in the confining bed.				
	51-60	G10.0	QET	Maximum evapotranspiration rate ( $L/T$ ).
	61-70	G10.0	ETDIST	Depth at which ET ceases below land surface ( $L$ ).
NOTE.—QET and ETDIST required only for simulations including evapotranspiration.				
	71-80	G10.0	<u>LENGTH</u>	Definition depends on the numerical solution used: <i>LSOR</i> : number of LSOR iterations between 2-D corrections. <i>ADI and SIP</i> : Number of iteration parameters; unless the program is modified, code 10 for SIP.
3	1-10	G10.0	<u>HMAX</u>	Definition depends on numerical solution used: <i>LSOR</i> : acceleration parameter. <i>ADI</i> : maximum iteration parameter. <i>SIP</i> : value of $\beta'$ .
NOTE.—See the discussion of the numerical methods in the text for information on iteration parameters.				
	11-20	G10.0	<u>FACTX</u>	Multiplication factor for transmissivity in X direction.
	21-30	G10.0	<u>FACTY</u>	Multiplication factor for transmissivity in Y direction.

NOTE.—FACTX = FACTY = 1 for isotropic aquifers.

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
4	1-20	G20.10	SUM	Parameters in which elapsed time and cumulative volumes for mass balance are stored. For the start of a simulation insert three blank cards. For continuation of a previous run from punched output, remove the three blank cards and insert the first three cards of the punched output from the previous run. If continuation is from interim storage on disk, the three blank cards should remain.
	21-40	G20.10	SUMP	
	41-60	G20.10	PUMPT	
	61-80	G20.10	CFLUXT	
5	1-20	G20.10	QRET	
	21-40	G20.10	CHST	
	41-60	G20.10	CHDT	
	61-80	G20.10	FLUXT	
6	1-20	G20.10	STORT	
	21-40	G20.10	ETFLXT	
	41-60	G20.10	FLXNT	

### Group III: Array data

Each of the following data sets, except the first one (PHI), consists of a parameter card and, if the data set contains variable data, may include a set of data cards. *Default typing applies except for M(I,J) which is a real array.* Each parameter card contains five variables defined as follows:

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
Every parameter card.	1-10	G10.0	FACT	If IVAR=0, FACT is the value assigned to every element of the matrix; If IVAR=1, FACT is the multiplication factor for the following set of data cards.
	11-20	G10.0	IVAR	=0 if no data cards are to be read for this matrix; =1 if data cards for this matrix follow.
	21-30	G10.0	IPRN	=0 if input data for this matrix are to be printed; =1 if input data for the matrix are <i>not</i> to be printed.
	31-40	G10.0	IRECS	=0 if the matrix is being read from cards or if each element is being set equal to FACT. =1 if the matrix is to be read from disk (unit 2).
	41-50	G10.0	IRECD	=0 if the matrix is <i>not</i> to be stored on disk. =1 if the matrix being read from cards or set equal to FACT is to be stored on disk (unit 2) for later retrieval.

Refer to the examples in figures 31-33, Attachment IV. Figure 33 illustrates data for the sample problem without using disk files.

For the uniform starting head=100, FACT=100, IVAR=IPRN=IRECS=IRECD=0 and no data cards are required. The storage coefficient matrix is used to locate a constant-head boundary; therefore, FACT=-1, IVAR=1, IPRN=IRECS=IRECD=0 and a set of data cards with the location of the boundary nodes follows.

To save the storage coefficient matrix on disk (provided unit 2 has been defined on a DD statement; see technical information), set FACT=1, IVAR=1, IPRN=IRECS=0, IRECD=1, and include the set of data cards (figure 31). After this has been processed successfully, subsequent runs need only include a parameter card with the following: FACT=IVAR=IPRN=0, IRECS=1, IRECD=0. The set of data cards are not included and the storage coefficient matrix is input via unit 2 from disk storage. (See figure 32.)

When data cards are included, start each row on a new card. To prepare a set of data cards for an array that is a function of space, the general procedure is to overlay the finite-difference grid on a contoured map of the parameter and record the average value of the parameter for each finite-difference block on coding forms according to the appropriate format. In general, record only significant digits and no decimal points (except for data set 2); use the multiplication factor to convert the data to their appropriate values. For example, if vertical conductivity of the confining bed (RATE) ranges from  $2 \times 10^{-9}$  to  $9 \times 10^{-8}$  ft/sec, coded values should range from 2 to 90; the multiplication factor (FACT) would be  $1.0 \text{ E}-9$ .

Arrays needed in every simulation are underlined. Omit parameter cards and data cards not used in the simulation (however, see the footnote for the S matrix).

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
1	1-80	8F10.4	PHI(I,J)	Head values for continuation of a previous run ( <i>L</i> ).

NOTE.—For a new simulation this data set is omitted. Do not include a parameter card with this data set.

2	1-80	8F10.4	<u>STRT(I,J)</u>	Starting head matrix ( <i>L</i> ).
3	1-80	20F4.0	<u>S(I,J)</u>	Storage coefficient (dimensionless).

NOTE.—Always required. In addition to specifying storage coefficient values for artesian aquifers, this matrix is used to locate constant-head boundaries by coding a negative number at constant-head nodes. At these nodes T or PERM must be greater than zero. For a problem with no constant-head nodes and that does not require S values, insert a blank parameter card.

4	1-80	20F4.0	T(I,J)	Transmissivity ( $L^2/T$ ).
---	------	--------	--------	-----------------------------

NOTE.—(1) Required for artesian aquifer simulation only.

(2) Zero values must be placed around the perimeter of the T or PERM matrix for reasons inherent in the computational scheme. If IVAR=0, zero values are automatically inserted around the border of the model.

5	1-80	20F4.0	PERM(I,J)	Hydraulic conductivity ( $L/T$ ) (see note 2 for data set 4).
6	1-80	20F4.0	BOTTOM(I,J)	Elevation of bottom of aquifer ( <i>L</i> ).
7	1-80	20F4.0	SY(I,J)	Specific yield (dimensionless).

NOTE.—Data sets 5, 6, and 7 are required for water table or combined artesian-water table simulations.

8	1-80	20F4.0	TOP (I,J)	Elevation of top of aquifer ( <i>L</i> ).
---	------	--------	-----------	---

NOTE.—Required only in combined artesian-water-table simulations.



DATA SET	COLUMNS	FORMAT	VARIABLE	DEFINITION
9	1-80	20F4.0	RATE (I,J)	Hydraulic conductivity of confining bed ( $L/T$ ).
10	1-80	20F4.0	RIVER (I,J)	Head on the other side of confining bed ( $L$ ).
11	1-80	20F4.0	M (I,J)	Thickness of confining bed ( $L$ ).
NOTE.—Data sets 9, 10, and 11 are required in simulations with leakage. If the confining bed or streambed does not extend over the entire aquifer use the M matrix to locate the confining bed. If RATE and RIVER do not vary over the extent of the confining bed they can be initialized to a uniform value.				
12	1-80	20F4.0	GRND (I,J)	Land elevation ( $L$ ).
NOTE.—Required for simulations with evapotranspiration.				
13	1-80	20F4.0	QRE (I,J)	Recharge rate ( $L/T$ ).
NOTE.—Omit if not used.				
14	1-80	8G10.0	DELX (J)	Grid spacing in X direction ( $L$ ).
15	1-80	8G10.0	DELY (I)	Grid spacing in Y direction ( $L$ ).

#### Group IV: Parameters that change with the pumping period

The program has two options for the simulation period:

1. To simulate a given number of time steps, set TMAX to a value larger than the expected simulation period. The program will use NUMT, CDLT, and DELT as coded.
2. To simulate a given pumping period, set NUMT larger than the number required for the simulation period (for example, 100). The program will compute the exact DELT (which will be  $\leq$  DELT coded) and NUMT to arrive exactly at TMAX on the last time step.

*Default typing applies.*

CARD	COLUMNS	FORMAT	VARIABLE	DEFINITION
1	1-10	G10.0	KP	Number of the pumping period.
	11-20	G10.0	KPM1	Number of the previous pumping period.
NOTE.—In general KPM1=0 if KP=1 KPM1=1 if KP=2, etc.				
This causes the time parameter used in ENTRY CLAY to be set to zero and STRT to be initialized to PHI. However, for continuation of a previous pumping period KPM1=KP, and STRT and the time parameter are not affected.				
	21-30	G10.0	NWEL	Number of wells for this pumping period.
	31-40	G10.0	TMAX	Number of days in this pumping period.
	41-50	G10.0	NUMT	Number of time steps.
	51-60	G10.0	CDLT	Multiplying factor for DELT.
NOTE.—1.5 is commonly used.				
	61-70	G10.0	DELT	Initial time step in hours.

If  $NWEL=0$  the following set of cards is omitted.

DATA SET 1		(NWEL cards)	
COLUMNS	FORMAT	VARIABLE	DEFINITION
1-10	G10.0	I	Row location of well.
11-20	G10.0	J	Column location of well.
21-30	G10.0	WELL (I,J)	Pumping rate ( $L^3/T$ ), negative for a pumping well.
31-40	G10.0	RADIUS	Real well radius ( $L$ ).

NOTE.—Radius is required only for those wells, if any, where computation of drawdown at a real well radius is to be made.

For each additional pumping period, another set of group IV cards is required (that is, NPER sets of group IV cards are required).

If another simulation is included in the same job, insert a blank card before the next group I cards.

## Attachment IV

### Sample Aquifer Simulation And Job Control Language

This appendix includes examples of job control language (JCL) for several different runs and an example problem designed to illustrate many of the options in the program. The grid and boundary conditions for the problem are given in figure 25. Figure 30 illustrates in cross section the type of problem being simulated, but note that it is not to scale.

The listing of data with the JCL examples is not on a coding form, but it should not be

difficult to determine the proper location of the numbers since the fields are either 4 or 10 spaces. Zero values have not been coded on the data cards to avoid unnecessary punching.

Figures 31 and 32 illustrate the JCL and data decks for two successive simulations of the sample problem. They are designed to show the use of disk facilities to store array data and interim results. The first run (fig. 31) is terminated after 5 iterations and interim results are stored on the data set specified by the FT04F001 DD statement. Note that arrays S, PERM, DELX, and DELY have been stored in the array data set specified by the FT02F001 DD statement (a 1 appears in column 40 of the parameter card for these arrays). The second run (fig. 32) continues computations from the previous stopping point and calculates a solution. Note that PHI, S, PERM, DELX, and DELY are read from disk storage. The final example (fig. 33) illustrates the JCL and data deck for a run without using the disk files. Following figure 33 is the output for the sample prob-

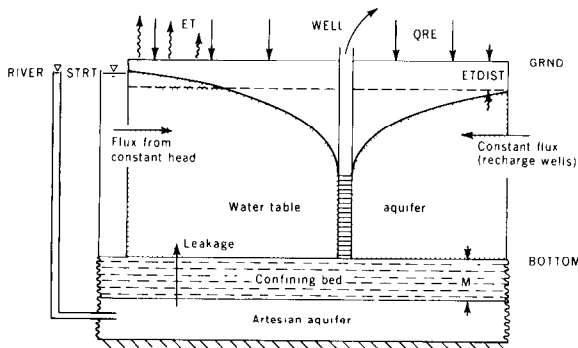


FIGURE 30.—Cross section illustrates several options included in the sample problem and identifies the meaning of several program parameters.

```
// EXEC FORTGCG
//FORT.SYSIN DD *
Model
source
cards
/*
//GO.FT02F001 DD DSN=A442702.AZ100.AG4W0000.MATRIX,
// UNIT=ONLINE,DISP=(NEW,KEEP),
// SPACE=(560,(14)),DCB=(RECFM=F)
//GO.FT04F001 DD DSN=A442702.AZ100.AG4W0000.HEAD,
// UNIT=ONLINE,DISP=(NEW,KEEP),SPACE=(TRK,(1,1),RLSE),
// DCB=(RECFM=VBS,LRECL=1168,BLKSIZE=1172)
//GO.SYSIN DD *
```

		---- SAMPLE AQUIFER PROBLEM ----									
Group I	WATE LEAK	EVAP	RECH	SIP	CHEC	DK2	NUME	HEAD			
	10	14	1	1	5	1	.1	FEET	10	10	
Group II	CONT	1	1	.003	.01	0	.4E-06				
	1	1	1								
STRT	100										
	-1	1				1					
S											
					1	1	1	1	1	1	
Group III		.002	1					1			
			1	1	1	2	2	2	2	2	
			1	1	1	2	2	2	2	2	
PERM			1	1	2	2	2	2	2	2	
			2	2	2	2	2	2	2	2	
			2	2	3	3	3	3	3	3	
			3	3	3	4	4	4	4	4	
	4	4	4	4	4	4	3	3	3	3	
					3	3	3	3	3	3	
BOTTOM		0									
SY											
RATE		.3E-07									
RIVER		100									
M		10									
GRND		105									
QRE		.2E-07									
DELX	50	1				1					
	20	14		9		9	14	21	31	41	
	37	25		17		11	9	13			
DELY	50	1				1					
	10	5		7		10	14	18	27	30	
	31	12									
Group IV	1	0		6		1	1	1.0	24		
	4	4		.05							
	5	4		.05							
	6	4		.05							
	7	4		.05							
	4	11		-10		2					
	6	6		-10							

FIGURE 31.—JCL and data deck to copy some of the data sets on disk, compute for 5 iterations, and store the results on disk.

```

// EXEC FORTGCG
//FORT.SYSIN DD *

```

		---- SAMPLE AQUIFER PROBLEM ----									
Group I	WATE LEAK	10	EVAP REOH	14	SIP	1	CHEC	50	DK1	NUME HEAD	
	CONT	1	1	1	1500	1	.1	FEET	10	10	
Group II		1	1	.003	.01	0	.4E-06				
		1	1	1							
Group III	STRT	100									
	S			1	1						
	PERM			1	1						
	BOTTOM	0									
	SY										
	RATE	.3E-07									
	RIVER	100									
	M	10									
	GRND	105									
	QRE	.2E-07									
DELX				1	1						
DELY						1	1	1.0	24		
Group IV		1	0	6	1	1	1.0	24			
		4	4	.05							
		5	4	.05							
		6	4	.05							
		7	4	.05							
		4	11	-10	2						
	6	6	-10								

```

//
//

```

FIGURE 32.—JCL and data deck to continue the previous run (fig. 31) to a solution.

lem generated using the JCL and problem deck shown in figure 33.

Figures 31 to 33 show that the source cards are being compiled for each run. It is more efficient, of course, to compile the source

deck once and store it as a load module on disk. Subsequent runs can use the load module with considerable reduction in cards read, CPU time, and lines printed.

```
// EXEC FORTGCG
//FORT.SYSIN DD *
Model
source
cards
/*
//GO.SYSIN DD *
```

		---- SAMPLE AQUIFER PROBLEM ----									
Group I	WATE LEAK	EVAP RECH	SIP	CHEC	NUME HEAD						
	10	14	1	50							
	CONT	1	1	1500	1	.1	FEET	10	10		
Group II	1	1	.003	.01	0	.4E-06					
	1	1	1								
STRT	100										
	-1	1									
S										1	
					1	1	1	1	1	1	
Group III	.002	1									
		1	1	1	2	2	2	2	2	2	
		1	1	1	2	2	2	2	2	2	2
PERM		1	1	1	2	2	2	2	2	2	2
		2	2	2	2	2	2	2	2	2	2
		2	2	2	3	3	3	3	3	3	3
		3	3	3	3	3	3	3	3	3	3
	4	4	4	4	4	4	4	4	4	4	4
					3	3	3	3	3	3	3
BOTTOM	0										
SY											
RATE	.3E-07										
RIVER	100										
M	10										
GRND	105										
QRE	.2E-07										
DELX	50	1									
	20	14	9	9	14	21	31	41			
	37	25	17	11	9	13					
DELY	50	1									
	10	5	7	10	14	18	27	30			
	31	12									
Group IV	1	0	6	1	1	1.0	24				
	4	4	.05								
	5	4	.05								
	6	4	.05								
	7	4	.05								
	4	11	-10	2							
	6	6	-10								

FIGURE 33.—JCL and data deck to simulate the sample problem without using disk files.

Program Output using data deck illustrated in figure 33

U. S. G. S.

FINITE-DIFFERENCE MODEL  
FOR  
SIMULATION OF GROUND-WATER FLOW  
JANUARY, 1975

\*\*\*\*\*

---- SAMPLE AQUIFER PROBLEM ----

\*\*\*\*\*

SIMULATION OPTIONS: WATE LEAK EVAP RECH SIP CHEC NUME HEAD

NUMBER OF ROWS = 10  
NUMBER OF COLUMNS = 14  
MAXIMUM PERMITTED NUMBER OF ITERATIONS = 50  
WORDS OF Y VECTOR USED = 3731

ON ALPHAMERIC MAP:  
MULTIPLICATION FACTOR FOR X DIMENSION = 1.000000  
MULTIPLICATION FACTOR FOR Y DIMENSION = 1.000000  
MAP SCALE IN UNITS OF FEET  
NUMBER OF FEET PER INCH = 1500.000  
MULTIPLICATION FACTOR FOR DRAWDOWN = 1.000000  
MULTIPLICATION FACTOR FOR HEAD = 0.9999996E-01  
NUMBER OF PUMPING PERIODS = 1  
TIME STEPS BETWEEN PRINTOUTS = 1  
ERROR CRITERIA FOR CLOSURE = 0.3000000E-02  
STEADY STATE ERROR CRITERIA = 0.9999998E-02  
SPECIFIC STORAGE OF CONFINING BED = 0.0  
EVAPOTRANSPIRATION RATE = 0.4000000E-06  
EFFECTIVE DEPTH OF ET = 10.00000  
MULTIPLICATION FACTOR FOR TRANSMISSIVITY IN X DIRECTION = 1.000000  
IN Y DIRECTION = 1.000000  
STARTING HEAD = 100.0000







GRID SPACING IN PROTOTYPE IN X DIRECTION  
 -----  
 1000. 700. 450. 450. 700. 1050. 1550. 2050. 1850. 1250. 850. 550.  
 450. 650.

GRID SPACING IN PROTOTYPE IN Y DIRECTION  
 -----  
 500. 250. 350. 500. 700. 900. 1350. 1500. 1550. 600.

SOLUTION BY THE STRONGLY IMPLICIT PROCEDURE

BETA= 1.00

10 ITERATION PARAMETERS: 0.0 0.6903903D 00 0.9041418D 00 0.9703214D 00 0.9908112D 00. 0.0  
 0.6903903D 00 0.9041418D 00 0.9703214D 00 0.9908112D 00

PUMPING PERIOD NO. 1: 1.00 DAYS  
 -----

NUMBER OF TIME STEPS= 1  
 DELT IN HOURS = 24,000  
 MULTIPLIER FOR DELT = 1,000

6 WELLS  
 -----  

| I | J  | PUMPING RATE | WELL RADIUS |
|---|----|--------------|-------------|
| 4 | 4  | 0.05         |             |
| 5 | 4  | 0.05         |             |
| 6 | 4  | 0.05         |             |
| 7 | 4  | 0.05         |             |
| 4 | 11 | -10.00       | 2.00        |
| 6 | 6  | -10.00       |             |

-----  
 | TIME STEP NUMBER = 1 |  
 -----

SIZE OF TIME STEP IN SECONDS = 86400.00  
 TOTAL SIMULATION TIME IN SECONDS = 86400.00  
 MINUTES = 1440.00  
 HOURS = 24.00  
 DAYS = 1.00  
 YEARS = 0.00

DURATION OF CURRENT PUMPING PERIOD IN DAYS = 1.00  
 YEARS = 0.00

|                          |            |                                |          |
|--------------------------|------------|--------------------------------|----------|
| CUMULATIVE MASS BALANCE: | L**3       | RATES FOR THIS TIME STEP:      | L**3/T   |
| -----                    | -----      | -----                          | -----    |
| SOURCES:                 |            | STORAGE =                      | 0.0      |
| -----                    |            | RECHARGE =                     | 1.1873   |
| STORAGE =                | 0.0        | CONSTANT FLUX =                | 0.2000   |
| RECHARGE =               | 102586.63  | PUMPING =                      | -20.0000 |
| CONSTANT FLUX =          | 17279.99   | EVAPOTRANSPIRATION =           | -0.2412  |
| CONSTANT HEAD =          | 1418633.00 | CONSTANT HEAD:                 |          |
| LEAKAGE =                | 210060.38  | IN =                           | 16.4194  |
| TOTAL SOURCES =          | 1748559.00 | OUT =                          | 0.0      |
|                          |            | LEAKAGE:                       |          |
| DISCHARGES:              |            | FROM PREVIOUS PUMPING PERIOD = | 0.0      |
| -----                    |            | TOTAL =                        | 2.4313   |
| EVAPOTRANSPIRATION =     | 20837.56   | SUM OF RATES =                 | -0.0032  |
| CONSTANT HEAD =          | 0.0        |                                |          |
| QUANTITY PUMPED =        | 1727998.00 |                                |          |
| LEAKAGE =                | 0.0        |                                |          |
| TOTAL DISCHARGE =        | 1748835.00 |                                |          |
|                          |            |                                |          |
| DISCHARGE-SOURCES =      | 276.00     |                                |          |
| PER CENT DIFFERENCE =    | 0.02       |                                |          |

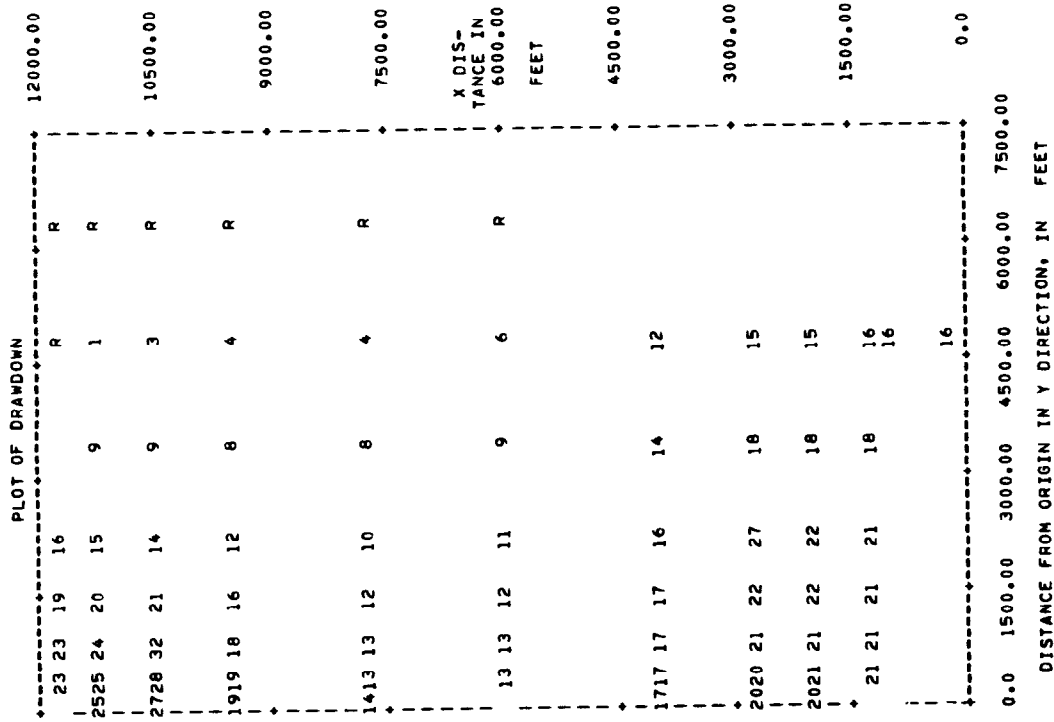
MAXIMUM HEAD CHANGE FOR EACH ITERATION:

|        |        |        |        |        |        |        |        |        |        |
|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|
| 9.5204 | 4.8325 | 3.7815 | 7.4434 | 3.4337 | 2.6980 | 1.3149 | 1.6210 | 1.1354 | 0.8168 |
| 0.4495 | 0.5055 | 0.3512 | 0.3693 | 0.2810 | 0.2107 | 0.0960 | 0.1267 | 0.0765 | 0.0509 |
| 0.0297 | 0.0322 | 0.0225 | 0.0234 | 0.0179 | 0.0133 | 0.0060 | 0.0080 | 0.0048 | 0.0032 |
| 0.0019 |        |        |        |        |        |        |        |        |        |

MAXIMUM CHANGE IN HEAD FOR THIS TIME STEP = 32.067

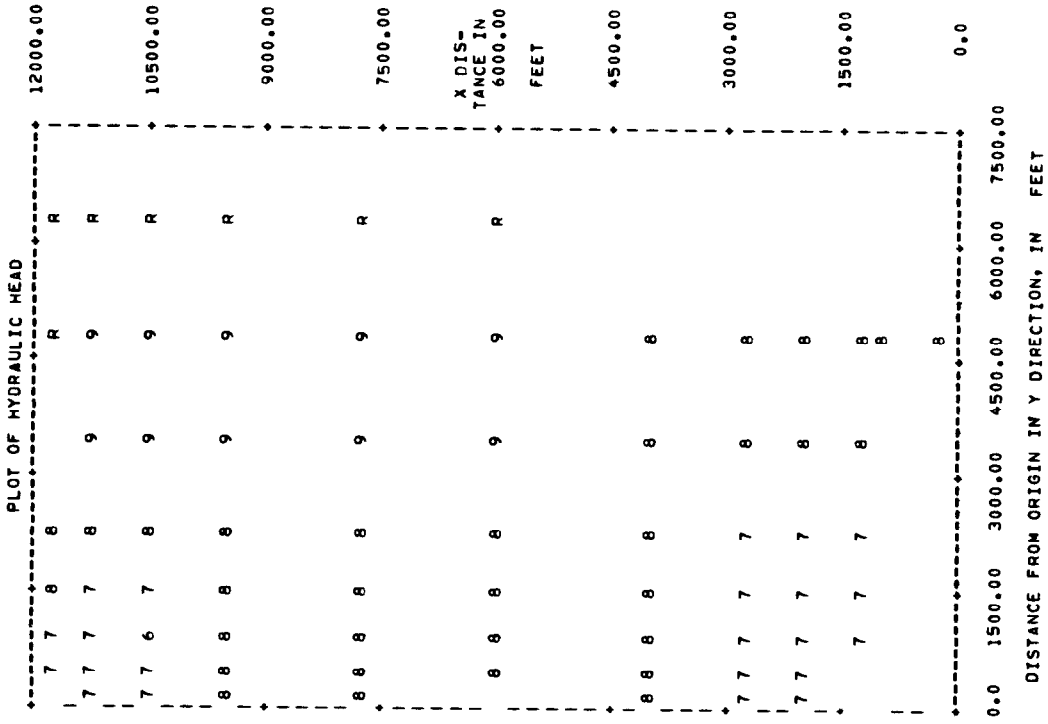
TIME STEP : 1

ITERATIONS: 30



EXPLANATION

- R = CONSTANT HEAD BOUNDARY
- \*\*\* = VALUE EXCEEDED 3 FIGURES
- MULTIPLICATION FACTOR = 1.000



EXPLANATION

R = CONSTANT HEAD BOUNDARY  
 \*\*\* = VALUE EXCEEDED 3 FIGURES  
 MULTIPLICATION FACTOR = 0.100



```

DRAWDOWN
-----
1  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
2  0.0  0.0  0.0  0.0  21.0  20.5  17.3  0.0  14.1  19.8  27.4  25.7  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
3  0.0  0.0  0.0  21.1  21.1  20.6  17.2  13.2  13.9  19.6  28.3  25.4  23.9  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
4  0.0  0.0  0.0  21.2  21.4  21.3  17.2  13.1  13.5  18.9  32.1  24.7  23.1  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
5  0.0  0.0  0.0  21.6  22.0  23.0  17.3  12.7  12.5  16.2  21.4  20.2  19.9  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
6  0.0  0.0  0.0  21.4  22.5  27.9  17.0  11.7  10.8  12.6  14.5  15.1  16.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
7  0.0  0.0  0.0  18.5  18.6  18.4  14.6  9.9  8.6  8.9  9.1  9.1  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
8  0.0  16.0  16.1  16.1  16.0  15.2  12.4  6.6  4.9  4.3  3.4  1.9  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
9  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0
10 0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0  0.0

```

```

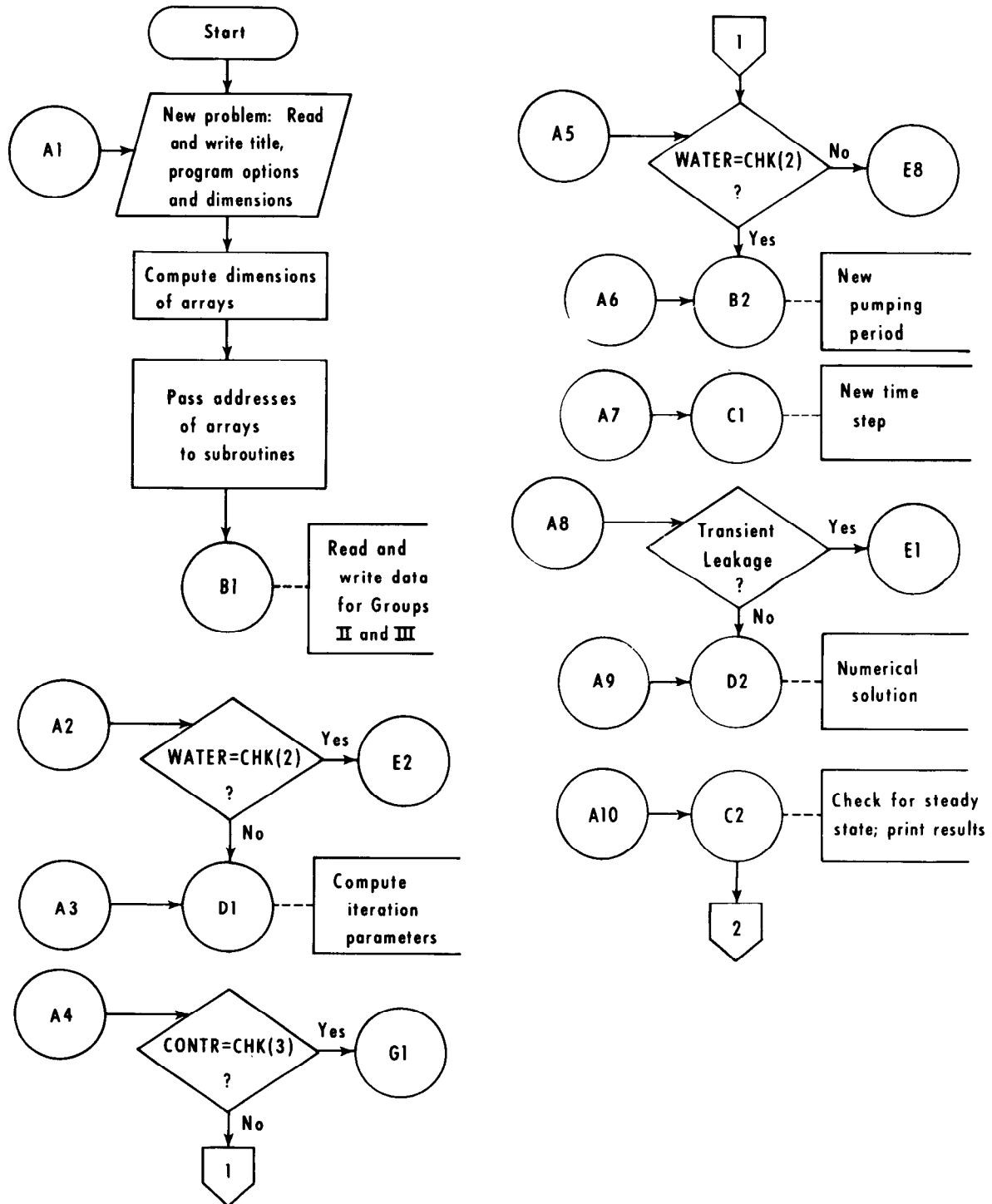
HEAD AND DRAWDOWN IN PUMPING WELLS
-----
I  J  WELL RADIUS  HEAD  DRAWDOWN
4  11  2.00  35.10  64.90

```

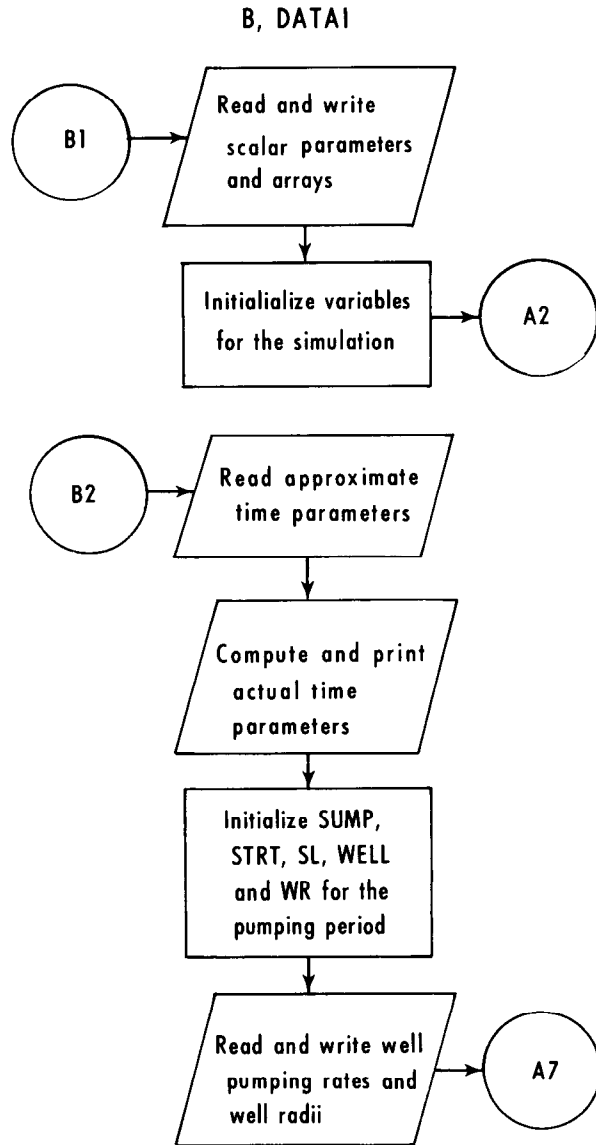
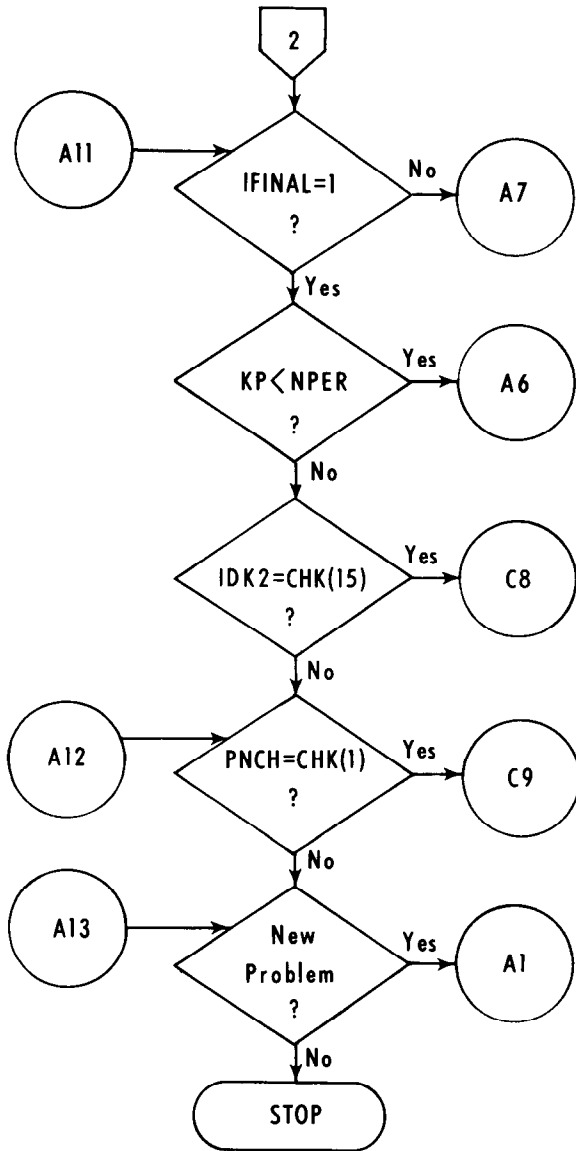
# Attachment V

## Generalized Flow Chart For Aquifer Simulation Model

### A, MAIN PROGRAM



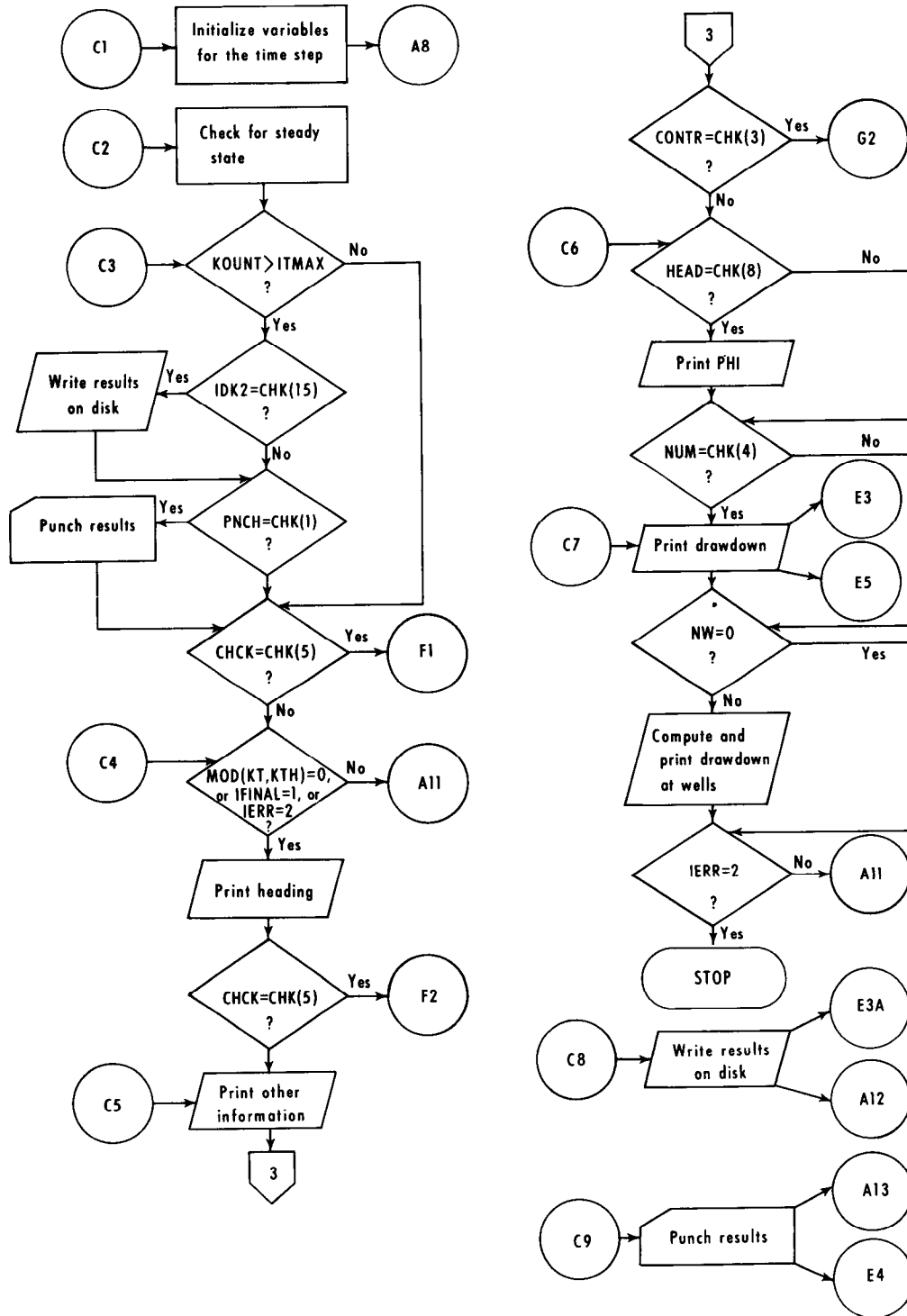
Flow chart—Continued



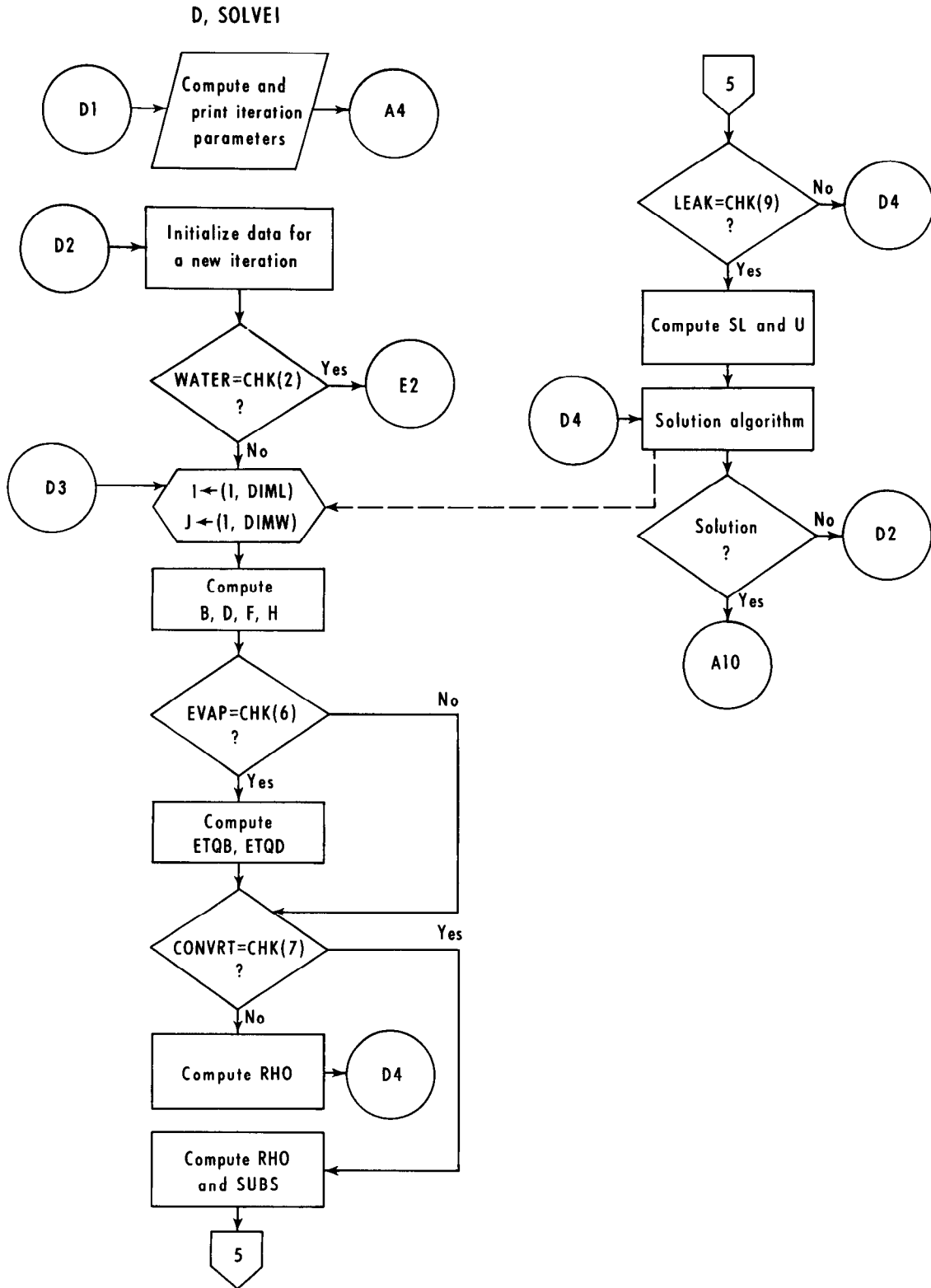


Flow chart—Continued

C, STEP

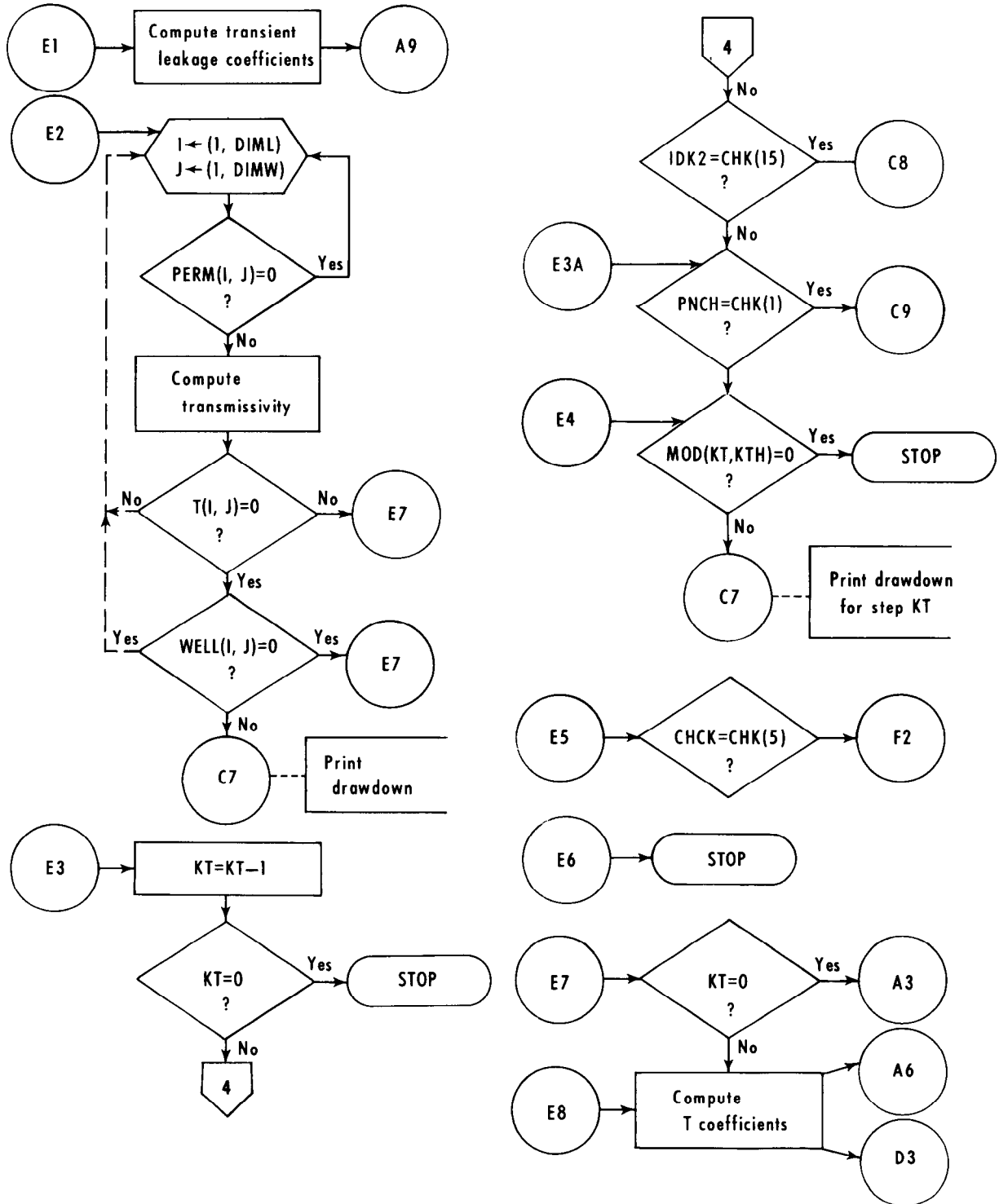


Flow chart—Continued

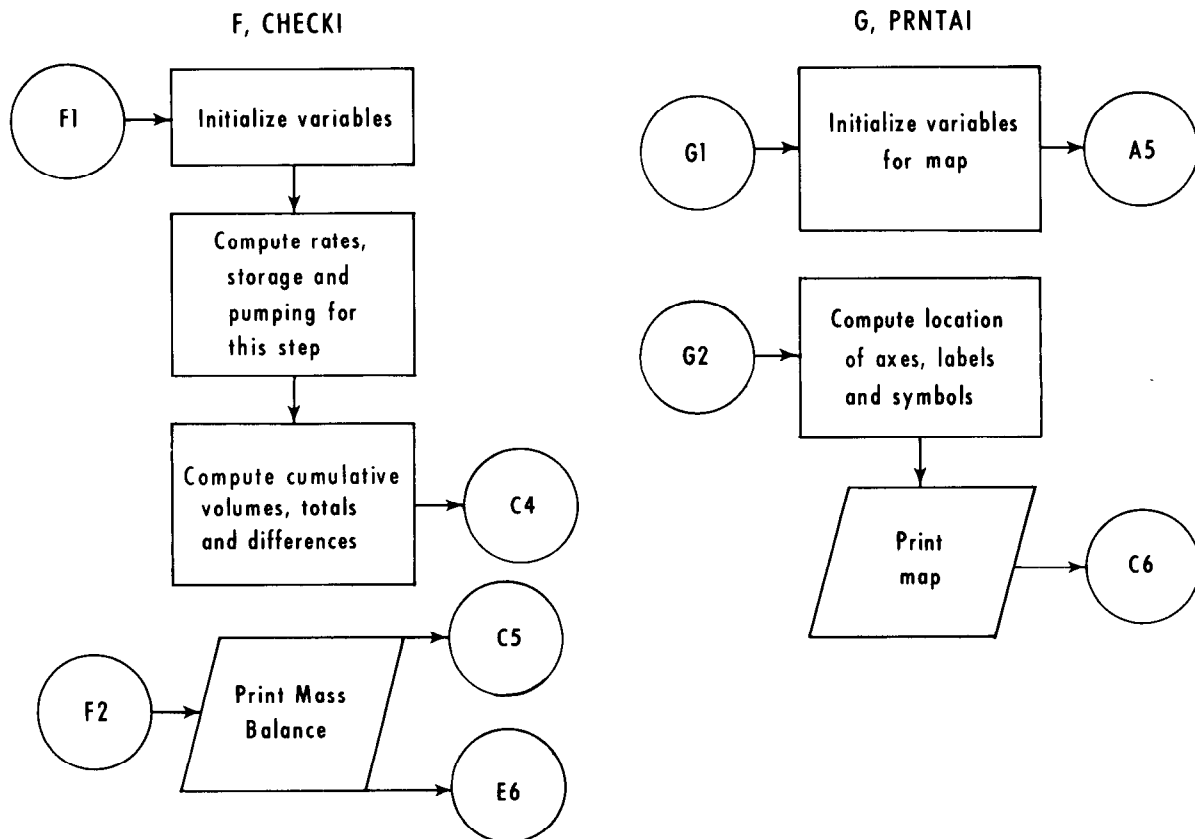


Flow chart—Continued

E, COEF



Flow chart—Continued



## Attachment VI

### Definition Of Program Variables

A IN DATAI, DUMMY ARRAY (DOES NOT USE CORE SPACE) USED TO OBTAIN ADDRESSES OF ARRAY DATA SETS;

ALFA CORRECTION VECTOR FOR ROWS (LSOR);  
PARAMETER IN SIP ALGORITHM;

B  $TC(I-1,J)/DELY(I)$  (1/T);

BE PARAMETER IN THOMAS ALGORITHM;

BOTTOM ELEVATION OF THE BOTTOM OF THE AQUIFER (L);

CDLT MULTIPLYING FACTOR FOR THE TIME STEP;

CHCK CONTAINS CHARACTER STRING FOR MASS BALANCE OPTION;

CHK VECTOR CONTAINING PROBLEM OPTIONS;

CONTR CONTAINS CHARACTER STRING FOR OPTION TO PRINT MARS OF DRAWDOWN AND/OR HEAD;

CONVRT CONTAINS CHARACTER STRING FOR WATER TABLE-ARTESIAN OPTION;

D  $TR(I,J-1)/DELX(J)$  (1/T);

DDN VECTOR THAT CONTAINS DRAWDOWN VALUES (L);

DEL ARRAY USED IN SIP ALGORITHM;

DELT TIME INCREMENT (T);

DELX GRID SPACING IN THE X DIRECTION (L);

DELY GRID SPACING IN THE Y DIRECTION (L);

DIML NUMBER OF ROWS;

DIMW NUMBER OF COLUMNS;

EROR STEADY STATE EROR CRITERION (L);

ERR CLOSURE CRITERION (L);

ETA ARRAY USED IN SIP ALGORITHM;

ETDIST DEPTH AT WHICH EVAPOTRANSPIRATION CEASES BELOW LAND SURFACE (L);

ETQB THAT PART OF ET SOURCE TERM TREATED IMPLICITLY;

ETQD THAT PART OF ET SOURCE TERM TREATED EXPLICITLY;

EVAP CONTAINS CHARACTER STRING FOR EVAPOTRANSPIRATION OPTION;

F  $TR(I,J)/DELX(J)$  (1/T);

FACT SEE EXPLANATION IN GROUP III; ARRAY DATA;

FACTX MULTIPLICATION FACTOR FOR TRANSMISSIVITY IN X DIRECTION;

FACTY MULTIPLICATION FACTOR FOR TRANSMISSIVITY IN Y DIRECTION;

G PARAMETER IN THOMAS ALGORITHM;

H  $TC(I,J)/DELY(I)$  (1/T);

GRND ELEVATION OF LAND SURFACE (L);

HEAD CONTAINS CHARACTER STRING FOR OPTION TO PRINT HEAD VALUES;

HEADNG TITLE FOR SIMULATION;

HMAX MAXIMUM ITERATION PARAMETER (ADI);  
ACCELERATION PARAMETER (LSOR);  
BETA PARAMETER (SIP);

IC INDICATOR USED TO DETERMINE THE TYPE OF ARRAY DATA;

IERR = 0 PUMPING WELLS ARE IN SATURATED PART OF WATER TABLE AQUIFER;  
= 1 PUMPING WELL HAS GONE DRY;

IFINAL = 0 ALL TIME STEPS EXCEPT THE LAST;  
= 1 LAST TIME STEP IN PUMPING PERIOD;

IFMT1,IFMT2,IFMT3 VARIABLE FORMAT ARRAYS PASSED TO DATAI VIA ARRAY ENTRY POINT;

IN IN DATAI, DUMMY ARRAY TO WHICH NAME IS PASSED;

IN01 DIML-1;

IPRN SEE EXPLANATION IN GROUP III; ARRAY DATA;

IRECS,IRECD SEE EXPLANATION IN GROUP III; ARRAY DATA;

IRN RECORD NUMBER USED FOR DISK STORAGE AND RETRIEVAL OF ARRAY DATA;

*Definition of program variables—Continued*

ITMAX      MAXIMUM NUMBER OF ITERATIONS PER TIME STEP;  
 IVAR      SEE EXPLANATION IN GROUP III; ARRAY DATA;  
 ISUM      THE CUMULATIVE WORDS OF STORAGE USED IN THE Y VECTOR;  
 IZ,JZ,ETC. DIMENSIONS OF ARRAYS IN MODEL, COMPUTED IN MAIN PROGRAM;  
 JNO1      DIMW-1;  
 KEEP      HYDRAULIC HEAD AT THE PREVIOUS TIME STEP (L);  
 KKK      ASSOCIATED VARIABLE IN DEFINE FILE, INDICATES NUMBER OF  
           NEXT RECORD;  
 KOUNT     ITERATION COUNTER;  
 KP        NUMBER OF THE PUMPING PERIOD;  
 KPM1     NUMBER OF PREVIOUS PUMPING PERIOD;  
 KT        TIME STEP COUNTER;  
 KTH      NUMBER OF TIME STEPS BETWEEN PRINTOUTS;  
 L        VECTOR CONTAINING INITIAL ADDRESS OF ARRAYS;  
 LEAK     CONTAINS CHARACTER STRING FOR LEAKAGE OPTION;  
 LENGTH   NUMBER OF ITERATION PARAMETERS (SIP,ADI);  
           NUMBER OF ITERATIONS BETWEEN 2-D CORRECTION (LSOR);  
 M        THICKNESS OF CONFINING OR STREAM BED (L);  
 NPER     NUMBER OF PUMPING PERIODS;  
 NUM      CONTAINS CHARACTER STRING FOR OPTION TO PRINT DRAWDOWN;  
 NUMT     NUMBER OF TIME STEPS;  
 NW       NUMBER OF PUMPING WELLS FOR WHICH DRAWDOWN IS TO BE  
           COMPUTED AT A 'REAL' WELL RADIUS;  
 NWEL     NUMBER OF WELLS FOR A PUMPING PERIOD;  
 NWR      LOCATION OF WELLS;  
 PNCH     CONTAINS CHARACTER STRING FOR OPTION TO PUNCH HYDRAULIC  
           HEAD VALUES;  
 P        PRINTER UNIT NUMBER;  
 PARAM    ITERATION PARAMETER;  
 PERM     HYDRAULIC CONDUCTIVITY OF THE AQUIFER (L/T);  
 PHE      HYDRAULIC HEAD AT THE START OF THE ITERATION (L);  
 PHI      HYDRAULIC HEAD (L);  
 PU       PUNCH UNIT NUMBER;  
 QET      MAXIMUM EVAPOTRANSPIRATION RATE (L/T);  
 QRE      RECHARGE RATE (L/T);  
 R        READER UNIT NUMBER;  
 RADIUS   REAL WELL RADIUS (L);  
 RATE     VERTICAL HYDRAULIC CONDUCTIVITY OF THE CONFINING BED  
           OR STREAM BED (L/T);  
 RECH     CONTAINS CHARACTER STRING FOR RECHARGE OPTION;  
 RHO      S/DELTA (1/T);  
 RHOP     VECTOR CONTAINING ITERATION PARAMETERS;  
 RIVER    HYDRAULIC HEAD OF THE STREAM OR IN THE AQUIFER  
           ABOVE OR BELOW THE PUMPED AQUIFER (L);  
 RW       WELL AND RECHARGE SOURCE TERM (L/T);  
 S        STORAGE COEFFICIENT;  
 SIP      CONTAINS CHARACTER STRING FOR SIP OPTION;  
 SL       STEADY PART OF LEAKAGE COEFFICIENT (L/T);  
 SLEAK    INITIAL & TRANSIENT LEAKAGE (L/T);  
 SS       SPECIFIC STORAGE OF CONFINING BED (1/L);  
 STORE    CONTAINS EITHER THE STORAGE COEFFICIENT OR SPECIFIC  
           YIELD DEPENDING ON THE TYPE OF AQUIFER;  
 STRT     HYDRAULIC HEAD AT THE BEGINNING OF THE CURRENT  
           PUMPING PERIOD (L);  
 SUBS     MODIFIES STORAGE TERM IN WATER TABLE-ARTESIAN CONVERSION;  
 SUM      TOTAL ELAPSED TIME IN THE SIMULATION (T);  
 SUMP     TOTAL ELAPSED TIME IN THE PUMPING PERIOD (T);  
 SURI     HYDRAULIC HEAD AT THE START OF THE SIMULATION (L);  
 SY       SPECIFIC YIELD;  
 T        TRANSMISSIVITY (L\*\*2/T);  
 TC       HARMONIC AVERAGE OF T/DELY \* I+1/2.J (L/T);

*Definition of program variables—Continued*

TEMP VECTOR FOR TEMPORARY STORAGE OF HYDRAULIC HEAD (L);  
 TEST = 0 CLOSURE CRITERION SATISFIED;  
       = 1 CLOSURE CRITERION NOT SATISFIED;  
 TEST2 MAXIMUM CHANGE IN HEAD FOR THE TIME STEP (L);  
 TEST3 VECTOR CONTAINING THE SUM OF THE ABSOLUTE VALUES  
       OF HEAD CHANGES FOR EACH ITERATION (L);  
 TL TRANSIENT PART OF LEAKAGE COEFFICIENT (1/T);  
 TMAX NUMBER OF DAYS IN THE PUMPING PERIOD (T);  
 TMIN MINIMUM VALUE OF DIMENSIONLESS TIME FOR THE CURRENT  
       PUMPING PERIOD;  
 TOP ELEVATION OF THE TOP OF THE AQUIFER (L);  
 TR HARMONIC AVERAGE OF  $T/DELX @ I, J+1/2$  (L/T);  
 TT MAXIMUM VALUE OF DIMENSIONLESS TIME FOR THE CURRENT  
       PUMPING PERIOD;  
 U = 0 EXPLICIT TREATMENT OF TRANSIENT LEAKAGE;  
       = 1 IMPLICIT TREATMENT OF TRANSIENT LEAKAGE;  
 U INDICATES DEFINE FILE RECORD LENGTH SPECIFICATION IN WORDS;  
 V ARRAY USED IN SIP ALGORITHM;  
 VF4 VARIABLE FORMAT FOR PRINTING HEAD AND DRAWDOWN;  
 WATER CONTAINS CHARACTER STRING FOR WATER TABLE OPTION;  
 WELL WELL DISCHARGE (L\*\*3/T);  
 WR WELL RADIUS (L);  
 XI ARRAY CONTAINING INCREMENTAL HEAD VALUES IN SIP SOLUTION (L);  
 Y VECTOR CONTAINING ARRAY STORAGE;  
 YDIM LENGTH OF AQUIFER IN Y DIRECTION (L).

## DEFINITION OF VARIABLES IN CHECKI SUBROUTINE

-----  
 CFLUX INFLOW FROM RECHARGE WELLS (L\*\*3/T);  
 CFLUXT CUMULATIVE VOLUME OF WATER FROM RECHARGE WELLS (L\*\*3);  
 CHD1 RATE OF OUTFLOW TO CONSTANT HEAD BOUNDARY (L\*\*3/T);  
 CHD2 RATE OF INFLOW FROM CONSTANT HEAD BOUNDARY (L\*\*3/T);  
 CHDT CUMULATIVE DISCHARGE TO CONSTANT HEAD BOUNDARY (L\*\*3);  
 CHST CUMULATIVE VOLUME OF WATER INFLOW FROM CONSTANT  
       HEAD BOUNDARY (L\*\*3);  
 DIFF ERROR IN MASS BALANCE (L\*\*3);  
 ETFLUX EVAPOTRANSPIRATION RATE (L\*\*3/T);  
 ETFLXT CUMULATIVE DISCHARGE BY ET (L\*\*3);  
 FLUX RATE OF LEAKAGE DUE TO GRADIENTS AT THE START  
       OF THE PUMPING PERIOD (L\*\*3/T);  
 FLUXS NET LEAKAGE RATE (L\*\*3/T);  
 FLXN RATE OF DISCHARGE BY LEAKAGE (L\*\*3/T);  
 FLXNT CUMULATIVE VOLUME OF WATER DISCHARGED BY LEAKAGE (L\*\*3);  
 FLXPT CUMULATIVE VOLUME OF WATER INFLOW FROM LEAKAGE (L\*\*3);  
 PERCENT PERCENT ERROR IN CUMULATIVE MASS BALANCE;  
 PUMP DISCHARGE FROM WELLS (L\*\*3/T);  
 PUMPT CUMULATIVE VOLUME OF WATER DISCHARGED BY PUMPING WELLS (L\*\*3);  
 QREFLX RECHARGE RATE (L\*\*3/T);  
 QRET CUMULATIVE VOLUME OF WATER DERIVED FROM RECHARGE (L\*\*3);  
 STOR RATE OF CHANGE IN STORAGE FOR THE TIME STEP (L\*\*3/T);  
 STORT CUMULATIVE VOLUME OF WATER DERIVED FROM STORAGE (L\*\*3);  
 SUMR SUM OF RECHARGE AND DISCHARGE RATES FOR THE TIME STEP (L\*\*3/T);  
 TOTL1 CUMULATIVE VOLUME OF WATER FROM ALL SOURCES (L\*\*3);  
 TOTL2 CUMULATIVE VOLUME OF WATER DISCHARGED FROM THE SYSTEM (L\*\*3);  
 XNET NET LEAKAGE RATE FOR A CELL (L\*\*3/T).

## DEFINITION OF VARIABLES IN THE PRINTAI SUBROUTINE

-----  
 BLANK CONTAINS BLANK SYMBOLS;  
 DINCH NUMBER OF MAP UNITS PER INCH;  
 DIST LOCATION OF NEXT COLUMN OF NODAL VALUES TO BE PRINTED;

*Definition of variables in the PRNTAI subroutine—Continued*

FACT1 FACTOR FOR ADJUSTING VALUE OF DRAWDOWN PRINTED;  
FACT2 FACTOR FOR ADJUSTING VALUE OF HEAD PRINTED;  
K ADJUSTED VALUE OF DRAWDOWN OR HEAD;  
MESUR NAME OF MAP LENGTH UNIT;  
N INDEX FOR SYMBOLS;  
NA INDICES FOR LOCATING X LABEL;  
NC NUMBER OF BLANKS BEFORE GRAPH;  
N1 NUMBER OF LINES PER INCH;  
N2 NUMBER OF CHARACTERS PER INCH;  
N3 NUMBER OF CHARACTERS PER LINE;  
N4 NUMBER OF LINES IN THE PLOT;  
NB MAXIMUM NUMBER OF CHARACTERS IN Y DIRECTION;  
NXD NUMBER OF INCHES IN THE X DIMENSION OF PLOT;  
NYD NUMBER OF INCHES IN THE Y DIMENSION OF PLOT;  
PRNT CONTAINS THE ARRANGEMENT OF SYMBOLS FOR EACH LINE;  
SPACNG CONTOUR INTERVAL (L);  
SYM VECTOR CONTAINING SYMBOLS USED IN THE PLOT;  
TITLE TITLE FOR PLOT;  
VF1,VF2,VF3 VARIABLE FORMATS FOR CENTERING PLOT;  
XLABEL LABEL FOR X AXIS;  
XN NUMBERS FOR X AXIS;  
XN1 1 INCH/(N1\*2);  
XSCALE MULTIPLICATION FACTOR TO CONVERT MODEL LENGTH UNIT  
TO UNIT USED IN X DIRECTION ON MAPS;  
XSF X SCALE FACTOR;  
YLABEL LABEL FOR Y AXIS;  
YLEN LOCATION OF NEXT VALUE IN THE COLUMN TO BE PRINTED;  
YN NUMBERS FOR Y AXIS;  
YSCALE MULTIPLICATION FACTOR TO CONVERT MODEL LENGTH UNIT  
TO UNIT USED IN Y DIRECTION ON MAPS;  
YSF Y SCALE FACTOR;  
Z LOCATION OF NEXT LINE TO BE PRINTED.



# Attachment VII

## Program Listing

```

C *****MAN 10
C          FINITE-DIFFERENCE MODEL          MAN 20
C          FOR                              MAN 30
C          SIMULATION OF GROUND-WATER FLOW  MAN 40
C          IN TWO DIMENSIONS                MAN 50
C
C          BY P. C. TRESPOTT, G. F. PINDER AND S. P. LARSON  MAN 60
C          U. S. GEOLOGICAL SURVEY         MAN 70
C          SEPTEMBER, 1975                 MAN 80
C *****MAN 90
C *****MAN 100
C MAIN PROGRAM TO DIMENSION DIGITAL MODEL AND CONTROL SEQUENCE  MAN 110
C OF COMPUTATIONS                               MAN 120
C -----MAN 130
C SPECIFICATIONS:                               MAN 140
C REAL *4KEEP,M,HEADNG(32)                     MAN 150
C REAL *8PHI,G,BE,TEMP,Z,YY                   MAN 160
C INTEGER R,P,PU,DIWL,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,MAN 170
C ICONTR,LEAK,RECH,SIP,ADI                     MAN 180
C
C DIMENSION Y(70000), L(37), IFMT1(9), IFMT2(9), IFMT3(9), NAME(99),MAN 200
C 1 YY(1)                                       MAN 210
C EQUIVALENCE (YY(1),Y(1))                     MAN 220
C
C COMMON /SARRAY/ VF4(11),CHK(15)              MAN 230
C COMMON /ARSIZE/ IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IH,IMAX,IMX1  MAN 240
C COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,ERROR,LEMAN 250
C 1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,QET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,MAN 260
C 2NUMS,LSOR,ADI,DELT,SUM,SUMP,SURS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,MAN 270
C 3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIWL,DIMAN 280
C 4MW,JNO1,INO1,R,P,PU,I,J,IDX1,IDX2         MAN 290
C
C DATA IFMT1/4H(1H0,4H,I5,,4H10E1,4H1.3/,4H(1H ,4H,5X,,4H10E1,4H1.3)MAN 320
C 1,4H) /                                       MAN 330
C DATA IFMT2/4H('0',4H,I2,,4H2X,2,4H0F6,,4H1/(5,4HX,20,4HF6.1,4H)) MAN 340
C 1,4H /                                       MAN 350
C DATA IFMT3/4H(1H0,4H,I5,,4H14F9,4H.5/(,4H1H ,.4H5X,1,4H4F9,,4H5)) MAN 360
C 1,4H /                                       MAN 370
C DATA NAME/2*4H ,4H STO,4HRAGE,4H COE,4HFFIC,4HIENT,4*4H ,4H MAN 380
C 1 T,4HRANS,4HMISS,4HIVIT,4HY ,2*4H ,4H A,4HQUIF,4HER H,4HYCMAN 390
C 2RA,4HULIC,4H CON,4HDUCT,4HIVIT,4HY ,4H ,4H A,4HQUIF,4HER B,MAN 400
C 34HASE ,4HELEV,4HATIO,4HN ,3*4H ,4H S,4HPECI,4HFIC,4HYIEL,4MAN 410
C 4HD ,4*4H ,4HAQUI,4HFER ,4HTOP ,4HELEV,4HATIO,4HN ,4H ,4HMAN 420
C 5CONF,4HININ,4HG BE,4HD HY,4HDRAU,4HLIC ,4HCOND,4HUCTI,4HVITY,3*4H MAN 430
C 6 ,4H RIV,4HER H,4HEAD ,4*4H ,4H C,4HONFI,4HNING,4H BED,4H TMAN 440
C 7HI,4HCKNE,4HSS ,2*4H ,4H L,4HAND ,4HSURF,4HACE ,4HELEV,4HATIMAN 450
C 80,4HN ,3*4H ,4H ARE,4HAL R,4HECHA,4HRGE ,4HRATE,2*4H / MAN 460
C
C DEFINE FILE 2(14,2624,U,KKK)                 MAN 470
C *****MAN 480
C *****MAN 490
C *****MAN 500
C ---READ TITLE,PROGRAM OPTIONS AND PROGRAM SIZE--- MAN 510
C 10 READ (R,370) HEADNG                        MAN 520
C WRITE (P,360) HEADNG                        MAN 530
C READ (R,380) WATER,LEAK,CONVRT,EVAP,RECH,NUMS,CHCK,PNCH,IDX1,IDX2,MAN 540
C 1NUM,HEAD                                    MAN 550
C WRITE (P,390) WATER,LEAK,CONVRT,EVAP,RECH,NUMS,CHCK,PNCH,IDX1,IDX2MAN 560
C 1,NUM,HEAD                                    MAN 570

```

## Program listing—Continued

```

      IF (NUMS.EQ.CHK(11).OR.NUMS.EQ.CHK(12).OR.NUMS.EQ.CHK(13)) GO TO 2MAN 580
10      WRITE (P,350) MAN 590
      STOP MAN 600
20      READ (R,320) DIML,DIMW,NW,ITMAX MAN 610
      WRITE (P,340) DIML,DIMW,NW,ITMAX MAN 620
C      MAN 630
C      ---COMPUTE DIMENSIONS FOR ARRAYS--- MAN 640
      IZ=DIML MAN 650
      JZ=DIMW MAN 660
      IH=MAX0(1,NW) MAN 670
      IMAX=MAX0(DIML,DIMW) MAN 680
      ISIZ=DIML*DIMW MAN 690
      ISUM=2*ISIZ+1 MAN 700
      IMX1=ITMAX+1 MAN 710
      L(1)=1 MAN 720
      DO 30 I=2,4 MAN 730
      L(I)=ISUM MAN 740
30      ISUM=ISUM+2*IMAX MAN 750
      DO 40 I=5,16 MAN 760
      L(I)=ISUM MAN 770
40      ISUM=ISUM+ISIZ MAN 780
      IF (WATER.NE.CHK(2)) GO TO 60 MAN 790
      DO 50 I=17,19 MAN 800
      L(I)=ISUM MAN 810
50      ISUM=ISUM+ISIZ MAN 820
      IP=DIML MAN 830
      JP=DIMW MAN 840
      GO TO 80 MAN 850
60      DO 70 I=17,19 MAN 860
      L(I)=ISUM MAN 870
70      ISUM=ISUM+1 MAN 880
      IP=1 MAN 890
      JP=1 MAN 900
80      IF (LEAK.NE.CHK(9)) GO TO 100 MAN 910
      DO 90 I=20,22 MAN 920
      L(I)=ISUM MAN 930
90      ISUM=ISUM+ISIZ MAN 940
      IR=DIML MAN 950
      JR=DIMW MAN 960
      GO TO 120 MAN 970
100     DO 110 I=20,22 MAN 980
      L(I)=ISUM MAN 990
110     ISUM=ISUM+1 MAN1000
      IR=1 MAN1010
      JR=1 MAN1020
120     IF (CONVRT.NE.CHK(7)) GO TO 130 MAN1030
      L(23)=ISUM MAN1040
      ISUM=ISUM+ISIZ MAN1050
      IC=DIML MAN1060
      JC=DIMW MAN1070
      GO TO 140 MAN1080
130     L(23)=ISUM MAN1090
      ISUM=ISUM+1 MAN1100
      IC=1 MAN1110
      JC=1 MAN1120
140     IF (EVAP.NE.CHK(6)) GO TO 150 MAN1130
      L(24)=ISUM MAN1140
      ISUM=ISUM+ISIZ MAN1150
      IL=DIML MAN1160
      JL=DIMW MAN1170
      GO TO 160 MAN1180
      MAN1190

```

## Program listing—Continued

```

150 L(24)=ISUM
    ISUM=ISUM+1
    IL=1
    JL=1
160 IF (NUMS.NE.CHK(11)) GO TO 180
    DO 170 I=25,28
    L(I)=ISUM
170 ISUM=ISUM+ISIZ
    IS=DIML
    JS=DIMW
    GO TO 200
180 DO 190 I=25,28
    L(I)=ISUM
190 ISUM=ISUM+1
    IS=1
    JS=1
200 DO 210 I=29,31
    L(I)=ISUM
210 ISUM=ISUM+DIMW
    DO 220 I=32,33
    L(I)=ISUM
220 ISUM=ISUM+DIML
    L(34)=ISUM
    ISUM=ISUM+IH
    L(35)=ISUM
    ISUM=ISUM+2*IH
    IF (MOD(ISUM,2).EQ.0) ISUM=ISUM+1
    CONTINUE
230 L(36)=ISUM
    ISUM=ISUM+2*IMAX
    L(37)=ISUM
    ISUM=ISUM+IMX1
    WRITE (P,330) ISUM
C
C ---PASS INTIIAL ADDRESSES OF ARRAYS TO SUBROUTINES---
CALL DATA1(Y(L(1)),Y(L(7)),Y(L(8)),Y(L(9)),Y(L(10)),Y(L(11)),Y(L(12)),Y(L(13)),Y(L(14)),Y(L(15)),Y(L(16)),Y(L(17)),Y(L(18)),Y(L(19)),Y(L(20)),Y(L(21)),Y(L(22)),Y(L(23)),Y(L(24)),Y(L(29)),Y(L(32)),Y(L(34)),Y(L(35)))
CALL STEP(Y(L(1)),Y(L(5)),Y(L(7)),Y(L(8)),Y(L(9)),Y(L(14)),Y(L(17)),Y(L(18)),Y(L(23)),Y(L(29)),Y(L(30)),Y(L(32)),Y(L(34)),Y(L(35)),Y(L(37)))
IF (NUMS.EQ.CHK(11)) CALL SOLVE1(Y(L(1)),Y(L(2)),Y(L(3)),Y(L(4)),Y(L(5)),Y(L(6)),Y(L(7)),Y(L(9)),Y(L(12)),Y(L(13)),Y(L(14)),Y(L(15)),Y(L(16)),Y(L(25)),Y(L(26)),Y(L(27)),Y(L(28)),Y(L(29)),Y(L(31)),Y(L(32)),Y(L(33)),Y(L(37)),Y(L(10)),Y(L(11)),Y(L(24)),Y(L(19)),Y(L(20)),Y(L(22)),Y(L(21)))
IF (NUMS.EQ.CHK(12)) CALL SOLVE2(Y(L(1)),Y(L(2)),Y(L(3)),Y(L(4)),Y(L(5)),Y(L(6)),Y(L(7)),Y(L(9)),Y(L(12)),Y(L(13)),Y(L(14)),Y(L(15)),Y(L(16)),Y(L(25)),Y(L(26)),Y(L(27)),Y(L(28)),Y(L(29)),Y(L(31)),Y(L(32)),Y(L(33)),Y(L(37)),Y(L(10)),Y(L(11)),Y(L(24)),Y(L(19)),Y(L(20)),Y(L(22)),Y(L(21)))
IF (NUMS.EQ.CHK(13)) CALL SOLVE3(Y(L(1)),Y(L(2)),Y(L(3)),Y(L(4)),Y(L(5)),Y(L(6)),Y(L(7)),Y(L(9)),Y(L(12)),Y(L(13)),Y(L(14)),Y(L(15)),Y(L(16)),Y(L(25)),Y(L(26)),Y(L(27)),Y(L(28)),Y(L(29)),Y(L(31)),Y(L(32)),Y(L(33)),Y(L(36)),Y(L(37)),Y(L(10)),Y(L(11)),Y(L(24)),Y(L(19)),Y(L(20)),Y(L(22)),Y(L(21)))
CALL COEF(Y(L(1)),Y(L(5)),Y(L(6)),Y(L(7)),Y(L(8)),Y(L(9)),Y(L(10)),Y(L(11)),Y(L(12)),Y(L(14)),Y(L(15)),Y(L(16)),Y(L(17)),Y(L(18)),Y(L(19)),Y(L(20)),Y(L(21)),Y(L(22)),Y(L(23)),Y(L(24)),Y(L(29)),Y(L(32)))

```

```

MAN1200
MAN1210
MAN1220
MAN1230
MAN1240
MAN1250
MAN1260
MAN1270
MAN1280
MAN1290
MAN1300
MAN1310
MAN1320
MAN1330
MAN1340
MAN1350
MAN1360
MAN1370
MAN1380
MAN1390
MAN1400
MAN1410
MAN1420
MAN1430
MAN1440
MAN1450
MAN1460
MAN1470
MAN1480
MAN1490
MAN1500
MAN1510
MAN1520
MAN1530
MAN1540
MAN1550
MAN1560
MAN1570
MAN1580
MAN1590
MAN1600
MAN1610
MAN1620
MAN1630
MAN1640
MAN1650
MAN1660
MAN1670
MAN1680
MAN1690
MAN1700
MAN1710
MAN1720
MAN1730
MAN1740
MAN1750
MAN1760
MAN1770
MAN1780
MAN1790
MAN1800

```

## Program listing—Continued

```

CALL CHECKI(Y(L(1)),Y(L(5)),Y(L(6)),Y(L(7)),Y(L(9)),Y(L(10)),Y(L(11)),Y(L(12)),Y(L(13)),Y(L(14)),Y(L(15)),Y(L(17)),Y(L(18)),Y(L(19)),Y(L(20)),Y(L(21)),Y(L(22)),Y(L(23)),Y(L(24)),Y(L(29)),Y(L(32)))
CALL PRNTAI(Y(L(1)),Y(L(8)),Y(L(9)),Y(L(12)),Y(L(14)),Y(L(29)),Y(L(32)))
C .....
C ---START COMPUTATIONS---
C *****
C ---READ AND WRITE DATA FOR GROUPS II AND III---
C CALL DATIN
CALL ARRAY(Y(L(12)),IFMT3,NAME(1),2)
IF (WATER.EQ.CHK(2)) GO TO 240
CALL ARRAY(Y(L(9)),IFMT3,NAME(10),3)
GO TO 250
240 CALL ARRAY(Y(L(17)),IFMT1,NAME(19),4)
CALL ARRAY(Y(L(18)),IFMT2,NAME(28),5)
CALL ARRAY(Y(L(19)),IFMT3,NAME(37),6)
250 IF (CONVRT.EQ.CHK(7)) CALL ARRAY(Y(L(23)),IFMT2,NAME(46),7)
IF (LEAK.NE.CHK(9)) GO TO 260
CALL ARRAY(Y(L(20)),IFMT1,NAME(55),8)
CALL ARRAY(Y(L(21)),IFMT2,NAME(64),9)
CALL ARRAY(Y(L(22)),IFMT2,NAME(73),10)
260 IF (EVAP.EQ.CHK(6)) CALL ARRAY(Y(L(24)),IFMT2,NAME(82),11)
IF (RECH.EQ.CHK(10)) CALL ARRAY(Y(L(13)),IFMT1,NAME(91),12)
CALL MDAT
C
C ---INITIALIZE TRANSMISSIVITY VALUES IN WATER TABLE PROBLEM---
KT=0
IF (WATER.EQ.CHK(2)) CALL TRANS
C
C ---COMPUTE ITERATION PARAMETERS---
IF (NUMS.EQ.CHK(11)) CALL ITER1
IF (NUMS.EQ.CHK(12)) CALL ITER2
IF (NUMS.EQ.CHK(13)) CALL ITER3
C
C ---INITIALIZE PARAMETERS FOR ALPHAMERIC MAP---
IF (CONTR.EQ.CHK(3)) CALL MAP
C
C ---COMPUTE T COEFFICIENTS FOR ARTESIAN PROBLEM---
IF (WATER.NE.CHK(2)) CALL TCOF
C
C ---READ TIME PARAMETERS AND PUMPING DATA FOR A NEW PUMPING PERIOD---
270 CALL NEWPER
C
KT=0
IFINAL=0
IERR=0
C
C ---START NEW TIME STEP COMPUTATIONS---
280 CALL NEWSTP
C
C ---COMPUTE TRANSIENT PART OF LEAKAGE TERM---
IF (LEAK.EQ.CHK(9).AND.SS.NE.0.) CALL CLAY
C
C ---ENTER APPROPRIATE SOLUTION ROUTINE AND COMPUTE SOLUTION---
IF (NUMS.EQ.CHK(11)) CALL NEWITA
IF (NUMS.EQ.CHK(12)) CALL NEWITB
IF (NUMS.EQ.CHK(13)) CALL NEWITC
C
C ---CHECK FOR STEADY STATE AND PRINT OUTPUT AT DESIGNATED

```

## Program listing—Continued

```

C      TIME STEPS---                                MAN2420
      CALL STEADY                                    MAN2430
C
C      ---LAST TIME STEP IN PUMPING PERIOD ?---    MAN2440
C      IF (IFINAL.NE.1) GO TO 280                  MAN2450
C
C      ---CHECK FOR NEW PUMPING PERIOD---          MAN2460
C      IF (KP.LT.NPER) GO TO 270                  MAN2470
C
C      ---DISK OUTPUT IF DESIRED---               MAN2480
C      IF (IDK2.NE.CHK(15)) GO TO 290            MAN2490
C      CALL DISK                                    MAN2500
C
C      ---PUNCHED OUTPUT IF DESIRED---           MAN2510
C      IF (PNCH.NE.CHK(1)) GO TO 300            MAN2520
290   CALL PUNCH                                    MAN2530
C
C      ---CHECK FOR NEW PROBLEM---                MAN2540
C      300 READ (R,320,END=310) NEXT             MAN2550
C      IF (NEXT.EQ.0) GO TO 10                  MAN2560
310   STOP                                          MAN2570
C
C      .....MAN2580
C      .....MAN2590
C      ---FORMATS---                               MAN2600
C      -----MAN2610
C      -----MAN2620
C      -----MAN2630
C      -----MAN2640
C      -----MAN2650
C      -----MAN2660
C      -----MAN2670
C      -----MAN2680
C      -----MAN2690
320   FORMAT (4I10)                                MAN2700
330   FORMAT ('0',54X,'WORDS OF Y VECTOR USED =',I7)
340   FORMAT ('0',62X,'NUMBER OF ROWS =',I5/60X,'NUMBER OF COLUMNS =',I5MAN2710
      1/9X,'NUMBER OF WELLS FOR WHICH DRAWDOWN IS COMPUTED AT A SPECIFIEDMAN2720
      2 RADIUS =',I5,/,39X,'MAXIMUM PERMITTED NUMBER OF ITERATIONS =',I5)MAN2730
350   FORMAT ('-',36X,'NO EQUATION SOLVING SCHEME SPECIFIED, EXECUTION TMAN2740
      1ERMINATED'/37X,58('*'))                    MAN2750
360   FORMAT ('1',60X,'U. S. G. S.'//55X,'FINITE-DIFFERENCE MODEL'/65X,'MAN2760
      1FOR'/51X,'SIMULATION OF GROUND-WATER FLOW'//60X,'JANUARY, 1975'//1MAN2770
      233('*')/'0',32A4//133('*'))                MAN2780
370   FORMAT (20A4)                                MAN2790
380   FORMAT (16(A4,1X))                            MAN2800
390   FORMAT ('-SIMULATION OPTIONS: ',13(A4,4X))   MAN2810
      END                                          MAN2820-

      SUBROUTINE DATAI(PHI,STRT,SURI,T,TR,TC,S,GRE,WELL,TL,SL,PERM,BOTTCDAT 10
      1M,SY,RATE,RIVER,M,TOP,GRND,DELX,DELY,WR,NWR)  DAT 20
C      -----DAT 30
C      READ AND WRITE INPUT DATA                    DAT 40
C      -----DAT 50
C      -----DAT 60
C      SPECIFICATIONS:                               DAT 70
C      REAL *8PHI,DBLE,XLABEL,YLABEL,TITLE,XN1,MESUR DAT 80
C      REAL *4M                                       DAT 90
C      INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,DAT 100
      1CONTR,LEAK,RECH,SIP,ADI                      DAT 110
C      -----DAT 120
C      DIMENSION PHI(IZ,JZ), STRT(IZ,JZ), SURI(IZ,JZ), T(IZ,JZ), TR(IZ,JZ)DAT 130
      1), TC(IZ,JZ), S(IZ,JZ), GRE(IZ,JZ), WELL(IZ,JZ), TL(IZ,JZ), SL(IZ,DAT 140
      2JZ), PERM(IP,JP), BOTTOM(IP,JP), SY(IP,JP), RATE(IR,JR), RIVER(IR,DAT 150
      3JR), M(IR,JR), TOP(IC,JC), GRND(IL,JL), DELX(JZ), DELY(IZ), WR(IH)DAT 160
      4, NWR(IH,2), A(IZ,JZ), IN(9), IFMT(9)        DAT 170
C      -----DAT 180

```

## Program listing—Continued

```

COMMON /SARRAY/ VF4(11),CHK(15) DAT 190
COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,FROR,LEDAT 200
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,DAT 210
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETGD,FACTX,FACTY,DAT 220
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DIDAT 230
4MW,JNO1,INO1,R,P,PU,I,J,IDK1,IDK2 DAT 240
COMMON /CK/ ETFLXT,STORT,QRET,CHST,CHDT,FLUXT,PUMPT,CFLUXT,FLXNT DAT 250
COMMON /PR/ XLABEL(3),YLABEL(6),TITLE(5),XN1,MESUR,PRNT(122),BLANKDAT 260
1(60),DIGIT(122),VF1(6),VF2(6),VF3(7),XSCALE,DINCH,SYM(17),XN(100),DAT 270
2YN(13),NA(4),N1,N2,N3,YSCALE,FACT1,FACT2 DAT 280
COMMON /ARSIZE/ IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IH,IMAX,IMX1 DAT 290
RETURN DAT 300
C ..... DAT 310
C ***** DAT 320
ENTRY DATAIN DAT 330
C ***** DAT 340
C ..... DAT 350
C ---READ AND WRITE SCALAR PARAMETERS--- DAT 360
READ (R,500) CONTR,XSCALE,YSCALE,DINCH,FACT1,FACT2,MESUR DAT 370
IF (CONTR.EQ.CHK(3)) WRITE (P,610) XSCALE,YSCALE,MESUR,MESUR,DINCHDAT 380
1,FACT1,FACT2 DAT 390
READ (R,490) NPER,KTH,ERR,EROR,SS,GET,ETDIST,LENGTH,HMAX,FACTX,FACDAT 400
ITY DAT 410
IF (ETDIST.LE.0.) ETDIST=1. DAT 420
WRITE (P,520) NPER,KTH,ERR,EROR,SS,GET,ETDIST,FACTX,FACTY DAT 430
C ..... DAT 440
C ---READ CUMULATIVE MASS BALANCE PARAMETERS--- DAT 450
READ (R,600) SUM,SUMP,PUMPT,CFLUXT,QRET,CHST,CHDT,FLUXT,STORT,ETFLDAT 460
1XT,FLXNT DAT 470
IF (IDK1.EQ.CHK(14)) GO TO 20 DAT 480
IF (SUM.EQ.0.0) GO TO 40 DAT 490
WRITE (P,480) SUM DAT 500
C ..... DAT 510
C ..... DAT 520
C ---HEAD DATA TO CONTINUE PREVIOUS COMPUTATIONS READ HERE--- DAT 530
C -----FROM CARDS: DAT 540
DO 10 I=1,DIML DAT 550
READ (R,540) (PHI(I,J),J=1,DIMW) DAT 560
10 WRITE (P,530) I,(PHI(I,J),J=1,DIMW) DAT 570
GO TO 40 DAT 580
C -----READ AND WRITE DATA FROM UNIT 4 ON DISK RATHER THAN CARDS: DAT 590
20 READ (4) PHI,SUM,SUMP,PUMPT,CFLUXT,QRET,CHST,CHDT,FLUXT,STORT,ETFLDAT 600
1XT,FLXNT DAT 610
WRITE (P,480) SUM DAT 620
DO 30 I=1,DIML DAT 630
30 WRITE (P,530) I,(PHI(I,J),J=1,DIMW) DAT 640
REWIND 4 DAT 650
C ..... STRT (STARTING HEAD) ..... DAT 660
40 READ (R,490) FACT,IVAR,IPRN,IRECS,IRECD DAT 670
IF (IRECS.EQ.1) READ (2,1) STRT DAT 680
IF ((IVAR.EQ.1.OR.IRECS.EQ.1).AND.IPRN.NE.1) WRITE (P,470) DAT 690
DO 80 I=1,DIML DAT 700
IF (IVAR.EQ.1) READ (R,540) (STRT(I,J),J=1,DIMW) DAT 710
DO 70 J=1,DIMW DAT 720
IF (IRECS.EQ.1) GO TO 60 DAT 730
IF (IVAR.NE.1) GO TO 50 DAT 740
STRT(I,J)=STRT(I,J)*FACT DAT 750
GO TO 60 DAT 760
50 STRT(I,J)=FACT DAT 770
60 SURI(I,J)=STRT(I,J) DAT 780
T(I,J)=0. DAT 785

```

## Program listing—Continued

```

      TL(I,J)=0.                                DAT 790
      SL(I,J)=0.                                DAT 800
      TR(I,J)=0.                                DAT 810
      TC(I,J)=0.                                DAT 820
      WELL(I,J)=0.0                             DAT 830
      QRE(I,J)=0.                                DAT 840
70  IF (SUM.EQ.0.0.AND.IDK1.NE.CHK(14)) PHI(I,J)=STRT(I,J)  DAT 850
      IF (IVAR.EQ.0.AND.IRECS.EQ.0.OR.IPRN.EQ.1) GO TO 80  DAT 860
      WRITE (P,530) I,(STRT(I,J),J=1,DIMW)             DAT 870
80  CONTINUE                                       DAT 880
      IF (IVAR.NE.1.AND.IRECS.NE.1) WRITE (P,420) FACT  DAT 890
      IF (IRECD.EQ.1) WRITE (2,1) STRT                 DAT 900
      RETURN                                           DAT 910
C
C  ---READ REMAINING ARRAYS FROM CARDS OR DISK (AS SPECIFIED IN THE  DAT 920
C  .OPTIONS) AND WRITE THEM ON DISK IF SPECIFIED IN THE OPTIONS---  DAT 930
C  *****                                         DAT 940
C  ENTRY ARRAY(A,IFMT,IN,IRN)                       DAT 950
C  *****                                         DAT 960
C  READ (R,490) FACT,IVAR,IPRN,IRECS,IRECD          DAT 970
      IK=4*IRECS+2*IVAR+IPRN+1                       DAT 980
      GO TO (90,90,110,110,140,140), IK              DAT1000
90  DO 100 I=1,DIML                                 DAT1010
      DO 100 J=1,DIMW                                 DAT1020
100  A(I,J)=FACT                                     DAT1030
      WRITE (P,430) IN,FACT                           DAT1040
      GO TO 160                                       DAT1050
110  IF (IK.EQ.3) WRITE (P,440) IN                  DAT1060
      DO 130 I=1,DIML                                 DAT1070
      READ (R,510) (A(I,J),J=1,DIMW)                 DAT1080
      DO 120 J=1,DIMW                                 DAT1090
120  A(I,J)=A(I,J)*FACT                             DAT1100
130  IF (IK.EQ.3) WRITE (P,IFMT) I,(A(I,J),J=1,DIMW)  DAT1110
      GO TO 160                                       DAT1120
140  READ (2,IRN) A                                  DAT1130
      IF (IK.EQ.6) GO TO 160                           DAT1140
      WRITE (P,440) IN                                 DAT1150
      DO 150 I=1,DIML                                 DAT1160
150  WRITE (P,IFMT) I,(A(I,J),J=1,DIMW)             DAT1170
160  IF (IRECD.EQ.1) WRITE (2,IRN) A                DAT1180
      RETURN                                           DAT1190
C
C  ---INSERT ZERO VALUES IN THE T OR PERM MATRIX AROUND THE  DAT1200
C  BORDER OF THE MODEL---                           DAT1210
C  *****                                         DAT1220
C  ENTRY MDAAT                                       DAT1230
C  *****                                         DAT1240
C  DO 180 I=1,DIML                                 DAT1250
      DO 180 J=1,DIMW                                 DAT1260
      IF (WATER.EQ.CHK(2)) GO TO 170                 DAT1280
      IF (I.EQ.1.OR.I.EQ.DIML.OR.J.EQ.1.OR.J.EQ.DIMW) T(I,J)=0.  DAT1290
      GO TO 180                                       DAT1300
170  IF (I.EQ.1.OR.I.EQ.DIML.OR.J.EQ.1.OR.J.EQ.DIMW) PERM(I,J)=0.  DAT1310
180  CONTINUE                                       DAT1320
C  ..... DELX,DELY ..... DAT1330
      READ (R,490) FACT,IVAR,IPRN,IRECS,IRECD        DAT1340
      IF (IRECS.EQ.1) GO TO 210                       DAT1350
      IF (IVAR.EQ.1) READ (R,490) DELX                DAT1360
      DO 200 J=1,DIMW                                 DAT1370
      IF (IVAR.NE.1) GO TO 190                       DAT1380
      DELX(J)=DELX(J)*FACT                            DAT1390

```

## Program listing—Continued

```

      GO TO 200
190 DELX(J)=FACT
200 CONTINUE
      GO TO 220
210 READ (2,13) DELX
220 IF (IRECD.EQ.1) WRITE (2,13) DELX
      IF (IVAR.EQ.1.OR.IRECS.EQ.1.AND.IPRN.NE.1) WRITE (P,550) DELX
      IF (IVAR.NE.1.AND.IRECS.NE.1) WRITE (P,450) FACT
      READ (R,490) FACT,IVAR,IPRN,IRECS,IRECD
      IF (IRECS.EQ.1) GO TO 250
      IF (IVAR.EQ.1) READ (R,490) DELY
      DO 240 I=1,DIML
      IF (IVAR.NE.1) GO TO 230
      DELY(I)=DELY(I)*FACT
      GO TO 240
230 DELY(I)=FACT
240 CONTINUE
      GO TO 260
250 READ (2,14) DELY
260 IF (IRECD.EQ.1) WRITE (2,14) DELY
      IF (IVAR.EQ.1.OR.IRECS.EQ.1.AND.IPRN.NE.1) WRITE (P,560) DELY
      IF (IVAR.NE.1.AND.IRECS.NE.1) WRITE (P,460) FACT
C
C ---INITIALIZE VARIABLES---
      JN01=DIMW-1
      IN01=DIML-1
      IF (LEAK.NE.CHK(9).OR.SS.NE.0.) GO TO 280
      DO 270 I=2,IN01
      DO 270 J=2,JN01
      IF (M(I,J).EQ.0.) GO TO 270
      TL(I,J)=RATE(I,J)/M(I,J)
270 CONTINUE
280 ETQB=0.0
      ETQD=0.0
      SUBS=0.0
      U=1.0
      TT=0.0
      IM=MIN0(6*DIMW+4,124)
      IM=(132-IM)/2
      VF4(3)=DIGIT(IM)
      VF4(8)=DIGIT(IM+5)
      WIDTH=0.
      DO 290 J=2,JN01
290 WIDTH=WIDTH+DELX(J)
      YDIM=0.
      DO 300 I=2,IN01
300 YDIM=YDIM+DELY(I)
      RETURN
C .....
C
C ---READ TIME PARAMETERS AND PUMPING DATA FOR A NEW PUMPING PERIOD---
C *****
C ENTRY NEWPER
C *****
C
C READ (R,490) KP,KPM1,NWEL,TMAX,NUMT,CDLT,DELT
C
C ---COMPUTE ACTUAL DELT AND NUMT---
      DT=DELT/24.
      TM=0.0
      DO 310 I=1,NUMT

```



## Program listing—Continued

```

DT=CDLT*DT
TM=TM+DT
IF (TM.GE.TMAX) GO TO 320
310 CONTINUE
GO TO 330
320 DELT=TMAX/TM*DELT
NUMT=I
330 WRITE (P,570) KP,TMAX,NUMT,DELT,CDLT
DELT=DELT*3600.
TMAX=TMAX*86400.
C
C ---INITIALIZE SUMP, STRT, SL, WELL AND WR---
WRITE (P,580) NWEL
IF (KP.GT.KPM1) SUMP=0.
DO 350 I=1,DIML
DO 350 J=1,DIMW
IF (KP.EQ.KPM1) GO TO 340
STRT(I,J)=PHI(I,J)
340 IF (LEAK.NE.CHK(9)) GO TO 350
IF (M(I,J).EQ.0.) GO TO 350
SL(I,J)=RATE(I,J)/M(I,J)*(RIVER(I,J)-STRT(I,J))
350 WELL(I,J)=0.
IF (NW.EQ.0) GO TO 370
DO 360 I=1,NW
360 WR(I)=0.
370 IF (NWEL.EQ.0) GO TO 410
C
C ---READ AND WRITE WELL PUMPING RATES AND WELL RADII---
KW=0
DO 400 II=1,NWEL
READ (R,490) I,J,WELL(I,J),RADIUS
IF (RADIUS.EQ.0.) GO TO 380
KW=KW+I
IF (KW.GT.NW) GO TO 380
NWR(KW,1)=I
NWR(KW,2)=J
WR(KW)=RADIUS
WRITE (P,590) I,J,WELL(I,J),WR(KW)
GO TO 390
380 WRITE (P,590) I,J,WELL(I,J)
390 WELL(I,J)=WELL(I,J)/(DELX(J)*DELY(I))
400 CONTINUE
410 RETURN
C
C .....
C
C FORMATS:
C
C -----
C
C
420 FORMAT ('0',63X,'STARTING HEAD =',G15.7)
430 FORMAT ('0',41X,9A4,'=',G15.7)
440 FORMAT ('1',49X,9A4,'/,65X,'MATRIX',/,50X,36('-'))
450 FORMAT ('0',72X,'DELX =',G15.7)
460 FORMAT ('0',72X,'DELY =',G15.7)
470 FORMAT ('1',60X,'STARTING HEAD MATRIX'/61X,20('-'))
480 FORMAT ('1',40X,' CONTINUATION - HEAD AFTER ',G20.7,' SEC PUMPING
1'/42X,58('-'))
490 FORMAT (8G10.0)
500 FORMAT (A4,6X,5G10.0,A8)
510 FORMAT (20F4.0)

```

DAT2010  
DAT2020  
DAT2030  
DAT2040  
DAT2050  
DAT2060  
DAT2070  
DAT2080  
DAT2090  
DAT2100  
DAT2110  
DAT2120  
DAT2130  
DAT2140  
DAT2150  
DAT2160  
DAT2170  
DAT2180  
DAT2190  
DAT2200  
DAT2210  
DAT2220  
DAT2230  
DAT2240  
DAT2250  
DAT2260  
DAT2270  
DAT2280  
DAT2290  
DAT2300  
DAT2310  
DAT2320  
DAT2330  
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DAT2380  
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DAT2400  
DAT2410  
DAT2420  
DAT2430  
DAT2440  
DAT2450  
DAT2460  
DAT2470  
DAT2480  
DAT2490  
DAT2500  
DAT2510  
DAT2520  
DAT2530  
DAT2540  
DAT2550  
DAT2560  
DAT2570  
DAT2580  
DAT2590  
DAT2600  
DAT2610



## Program listing—Continued

```

C      ---START A NEW TIME STEP---                      STP 340
C      *****                                         STP 350
C      ENTRY NEWSTP                                     STP 360
C      *****                                         STP 370
C      KT=KT+1                                          STP 380
C      KOUNT=0                                          STP 390
C      DO 10 I=1,DIML                                   STP 400
C      DO 10 J=1,DIMW                                   STP 410
10  KEEP(I,J)=PHI(I,J)                                 STP 420
C      DELT=CDLT*DELT                                  STP 430
C      SUM=SUM+DELT                                    STP 440
C      SUMP=SUMP+DELT                                  STP 450
C      DAYSP=SUMP/86400.                                STP 460
C      YRSP=DAYSP/365.                                  STP 470
C      HRS=SUM/3600.                                    STP 480
C      MINS=HRS*60.                                    STP 490
C      DAYS=HRS/24.                                    STP 500
C      YRS=DAYS/365.                                    STP 510
C      RETURN                                          STP 520
C      .....                                         STP 530
C      .....                                         STP 540
C      ---CHECK FOR STEADY STATE---                    STP 550
C      *****                                         STP 560
C      ENTRY STEADY                                     STP 570
C      *****                                         STP 580
C      TEST2=0.                                         STP 590
C      DO 20 I=2,IN01                                   STP 600
C      DO 20 J=2,JN01                                   STP 610
20  TEST2=DMAX1(TEST2,DABS(DBLE(KEEP(I,J))-PHI(I,J))) STP 620
C      IF (TEST2.GE.ERROR) GO TO 30                    STP 630
C      WRITE (P,330) KT                                 STP 640
C      IFINAL=1                                         STP 650
C      GO TO 40                                         STP 660
30  IF (KT.EQ.NUMT) IFINAL=1                           STP 670
C      .....                                         STP 680
C      ---ENTRY FOR TERMINATING COMPUTATIONS IF MAXIMUM ITERATIONS STP 690
C      EXCEEDED---                                     STP 700
C      *****                                         STP 710
C      ENTRY TERM1                                     STP 720
C      *****                                         STP 730
C      40 IF (KT.GT.200) WRITE (P,400)                 STP 740
C      ITTO(KT)=KOUNT                                   STP 750
C      IF (KOUNT.LE.ITMAX) GO TO 80                    STP 760
C      IERR=2                                           STP 770
C      KOUNT=KOUNT-1                                    STP 780
C      ITTO(KT)=KOUNT                                   STP 790
C      IF (KT.EQ.1) GO TO 60                           STP 800
C      .....                                         STP 810
C      ---WRITE ON DISK OR PUNCH CARDS AS SPECIFIED IN THE OPTIONS--- STP 820
C      XXX=SUM-DELT                                     STP 830
C      IF (IDK2.EQ.CHK(15)) WRITE (4) ((KEEP(I,J),YYY,I=1,DIML),J=1,DIMW) STP 840
C      1,XXX,SUMP,PUMPT,CFLUXT,QRET,CHST,CHDT,FLUXT,STORT,ETFLXT,FLXNT STP 850
C      IF (PNCH.NE.CHK(1)) GO TO 80                    STP 860
C      WRITE (PU,360) XXX,SUMP,PUMPT,CFLUXT,QRET,CHST,CHDT,FLUXT,STORT,ET STP 870
C      IFLXT,FLXNT                                     STP 880
C      DO 50 I=1,DIML                                   STP 890
50  WRITE (PU,350) (KEEP(I,J),J=1,DIMW)                STP 900
C      GO TO 80                                         STP 910
60  IF (IDK2.EQ.CHK(15)) WRITE (4) PHI,SUM,SUMP,PUMPT,CFLUXT,QRET,CHST STP 920
C      1,CHDT,FLUXT,STORT,ETFLXT,FLXNT                STP 930
C      IF (PNCH.NE.CHK(1)) GO TO 80                    STP 940

```

## Program listing—Continued

```

WRITE (PU,360) SUM,SUMP,PUMPT,CFLUXT,GRET,CHST,CHDT,FLUXT,STORT,ETSTP 950
1FLXT,FLXNT STP 960
DO 70 I=1,DIML STP 970
70 WRITE (PU,350) (PHI(I,J),J=1,DIMW) STP 980
C STP 990
80 IF (CHCK.EQ.CHK(5)) CALL CHECK STP1000
IF (IERR.EQ.2) GO TO 90 STP1010
C STP1020
C ---PRINT OUTPUT AT DESIGNATED TIME STEPS--- STP1030
IF (MOD(KT,KTH).NE.0.AND.IFINAL.NE.1) RETURN STP1040
90 WRITE (P,340) KT,DELT,SUM,MINS,HRS,DAYS,YRS,DAYSP,YRSP STP1050
IF (CHCK.EQ.CHK(5)) CALL CWRITE STP1060
IF (TT.NE.0.) WRITE (P,320) TMIN,TT STP1070
KOUNT=KOUNT+1 STP1080
WRITE (P,300) (TEST3(J),J=1,KOUNT) STP1090
WRITE (P,290) TEST2 STP1100
I3=1 STP1110
I5=0 STP1120
100 I5=I5+40 STP1130
I4=MIN0(KT,I5) STP1140
WRITE (P,390) (I,I=I3,I4) STP1150
WRITE (P,380) STP1160
WRITE (P,370) (ITTO(I),I=I3,I4) STP1170
WRITE (P,380) STP1180
IF (KT.LE.I5) GO TO 110 STP1190
I3=I3+40 STP1200
GO TO 100 STP1210
C STP1220
C ---PRINT ALPHAMERIC MAPS--- STP1230
110 IF (CONTR.NE.CHK(3)) GO TO 120 STP1240
IF (FACT1.NE.0.) CALL PRNTA(1) STP1250
IF (FACT2.NE.0.) CALL PRNTA(2) STP1260
120 IF (HEAD.NE.CHK(8)) GO TO 140 STP1270
C STP1280
C ---PRINT HEAD MATRIX--- STP1290
WRITE (P,310) STP1300
DO 130 I=1,CIML STP1310
130 WRITE (P,VF4) I,(PHI(I,J),J=1,DIMW) STP1320
140 IF (NUM.NE.CHK(4)) GO TO 170 STP1330
C STP1340
C ---PRINT DRAWDOWN--- STP1350
WRITE (P,280) STP1360
***** STP1370
ENTRY DRDN STP1380
***** STP1390
DO 160 I=1,DIML STP1400
DO 150 J=1,DIMW STP1410
150 DDN(J)=SURI(I,J)-PHI(I,J) STP1420
160 WRITE (P,VF4) I,(DDN(J),J=1,DIMW) STP1430
170 IF (NW.EQ.0.OR.IERR.EQ.1) GO TO 230 STP1440
C STP1450
C ..... STP1460
C STP1470
C ---COMPUTE APPROXIMATE HEAD FOR PUMPING WELLS--- STP1480
WRITE (P,260) STP1490
DO 220 KW=1,NW STP1500
IF (WR(KW).EQ.0.) GO TO 220 STP1510
I=NWR(KW,1) STP1520
J=NWR(KW,2) STP1530
C STP1540
C COMPUTE EFFECTIVE RADIUS OF WELL IN MODEL--- STP1550
RE=(DELT(J)+DELT(I))/9.62

```



## Program listing—Continued

```

310 FORMAT ('1',60X,'HEAD MATRIX'/61X,11('='))          STP2170
320 FORMAT ('0DIMENSIONLESS TIME FOR THIS STEP RANGES FROM',G15.7,'  TSTP2180
10',G15.7)          STP2190
330 FORMAT ('-*****STEADY STATE AT TIME STEP',I4,'*****') STP2200
340 FORMAT (1H1,44X,57('=')/45X,'|',14X,'TIME STEP NUMBER =',I9,14X,'|STP2210
1'/45X,57('=')//50X,29HSIZE OF TIME STEP IN SECONDS=,F14.2//55X,'|TOSTP2220
2TAL SIMULATION TIME IN SECONDS=,F14.2/80X,8HMINUTES=,F14.2/82X,6HSTP2230
3HOURS=,F14.2/83X,5HSDAYS=,F14.2/82X,'YEARS=,F14.2//45X,'|DURATION STP2240
4OF CURRENT PUMPING PERIOD IN DAYS=,F14.2/82X,'YEARS=,F14.2//) STP2250
350 FORMAT (8F10.4)          STP2260
360 FORMAT (4G20.10)          STP2270
370 FORMAT ('0ITERATIONS:',40I3)          STP2280
380 FORMAT (' ',10('='))          STP2290
390 FORMAT ('0TIME STEP ',40I3)          STP2300
400 FORMAT ('0',10('*'),'THE NUMBER OF TIME STEPS EXCEEDS THE DIMENSIOSTP2310
1N OF THE VECTOR ITTO AND MAY CAUSE UNEXPECTED RESULTS IN ADDITIONASTP2320
2L'/'0COMPUTATION. AVOID PROBLEMS BY INCREASING THE DIMENSION OF TSTP2330
3HE VECTOR ITTO IN STEP',10('*'))          STP2340
END          STP2350-

SUBROUTINE SOLVE1(PHI,BE,G,TEMP,KEEP,PHE,STRT,T,S,QRE,WELL,TL,SL,DSIP 10
1EL,ETA,V,XI,DELX,BET,DELY,ALF,TEST3,TR,TC,GRND,SY, TOP,RATE,M,RIVERSIP 20
2)          SIP 30
C -----SIP 40
C SOLUTION BY THE STRONGLY IMPLICIT PROCEDURE          SIP 50
C -----SIP 60
C          SIP 70
C SPECIFICATIONS:          SIP 80
REAL *8PHI,DBLE,RHOP(20),G,BE,TEMP,DABS,W,TEST2,DMAX1,RHO,B,D,F,H,SIP 90
1B1,E,CH,GH,BH,DH,EH,FH,HH,ALFA,BETA,GAMA,RES          SIP 100
REAL *4KEEP,M          SIP 110
INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,SIP 120
1CONTR,LEAK,RECH,SIP,IORDER(21),ADI          SIP 130
C          SIP 140
DIMENSION PHI(1), BE(1), G(1), TEMP(1), KEEP(1), PHE(1), STRT(1), SIP 150
1T(1), S(1), QRE(1), WELL(1), TL(1), SL(1), DEL(1), ETA(1), V(1), XSIP 160
2I(1), DELX(1), BET(1), DELY(1), ALF(1), TEST3(1), TR(1), TC(1), GRSIP 170
3ND(1), SY(1), TOP(1), RATE(1), M(1), RIVER(1)          SIP 180
C          SIP 190
COMMON /SARRAY/ VF4(11),CHK(15)          SIP 200
COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,ERROR,LESIP 210
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,SIP 220
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,SIP 230
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DISIP 240
4MW,JN01,IN01,R,P,PU,I,J,IDX1,IDX2          SIP 250
RETURN          SIP 260
C .....SIP 270
C          SIP 280
C ---COMPUTE AND PRINT ITERATION PARAMETERS---          SIP 290
C *****SIP 300
C ENTRY ITER1          SIP 310
C *****SIP 320
C ---INITIALIZE ORDER OF ITERATION PARAMETERS (OR REPLACE WITH A          SIP 330
C READ STATEMENT)---          SIP 340
DATA IORDER/1,2,3,4,5,1,2,3,4,5,11*1/          SIP 350
I2=IN01-1          SIP 360
J2=JN01-1          SIP 370
L2=LENGTH/2          SIP 380
PL2=L2-1.          SIP 390
W=0.          SIP 400

```

## Program listing—Continued

```

          PI=0.                                SIP 410
C
C      ---COMPUTE AVERAGE MAXIMUM PARAMETER FOR PROBLEM--- SIP 420
          DO 10 I=2,IN01                       SIP 430
          DO 10 J=2,JN01                       SIP 440
          N=I+DIML*(J-1)                      SIP 450
          IF (T(N).EQ.0.) GO TO 10             SIP 460
          PI=PI+1.                             SIP 470
          DX=DELX(J)/WIDTH                    SIP 480
          DY=DELY(I)/YDIM                     SIP 490
          W=W+1.-AMIN1(2.*DX*DX/(1.+FACTY*DX*DX/(FACTX*DY*DY)),2.*DY*DY/(1.+SIP 510
          1FACTX*DY*DY/(FACTY*DX*DX)))        SIP 520
10 CONTINUE                                  SIP 530
          W=W/PI                               SIP 540
C
C      ---COMPUTE PARAMETERS IN GEOMETRIC SEQUENCE--- SIP 550
          PJ=-1.                               SIP 560
          DO 20 I=1,L2                         SIP 570
          PJ=PJ+1.                             SIP 580
          P=PL2**PJ                            SIP 590
20 TEMP(I)=1.-(1.-W)**(PJ/PL2)              SIP 600
C
C      ---ORDER SEQUENCE OF PARAMETERS--- SIP 610
          DO 30 J=1,LENGTH                     SIP 620
          RHOP(J)=TEMP(IORDER(J))             SIP 630
          WRITE (P,370) HMAX                   SIP 640
          WRITE (P,380) LENGTH,(RHOP(J),J=1,LENGTH) SIP 650
          RETURN                               SIP 660
          .....                               SIP 670
          .....                               SIP 680
C
C      ---INITIALIZE DATA FOR A NEW ITERATION--- SIP 690
          40 KOUNT=KOUNT+1                     SIP 700
          IF (KOUNT.LE.ITMAX) GO TO 50        SIP 710
          WRITE (P,360)                        SIP 720
          CALL TERM1                           SIP 730
          50 IF (MOD(KOUNT,LENGTH)) 60,60,70 SIP 740
          *****                             SIP 750
          ENTRY NEWITA                          SIP 760
          *****                             SIP 770
C
          60 NTH=0                             SIP 780
          70 NTH=NTH+1                         SIP 790
          W=RHOP(NTH)                         SIP 800
          TEST3(KOUNT+1)=0.                   SIP 810
          TEST=0.                              SIP 820
          N=DIML*DIMW                         SIP 830
          DO 80 I=1,N                          SIP 840
          PHE(I)=PHI(I)                       SIP 850
          DEL(I)=0.                           SIP 860
          ETA(I)=0.                            SIP 870
          V(I)=0.                              SIP 880
          80 XI(I)=0.                          SIP 890
          BIGI=0.0                            SIP 900
C
C      ---COMPUTE TRANSMISSIVITY AND T COEFFICIENTS IN WATER TABLE SIP 910
C      OR WATER TABLE-ARTESIAN SIMUATION--- SIP 920
          IF (WATER.NE.CHK(2)) GO TO 90      SIP 930
          CALL TRANS                           SIP 940
C
C      ---CHOOSE SIP NORMAL OR REVERSE ALGORITHM--- SIP 950
          90 IF (MOD(KOUNT,2)) 100,230,100  SIP 960
          .....                               SIP 970
          .....                               SIP 980
C
C      ---ORDER EQUATIONS WITH ROW 1 FIRST - 3X3 EXAMPLE: SIP 990
          .....                               SIP1000
          .....                               SIP1010

```

## Program listing—Continued

```

C           1 2 3                               SIP1020
C           4 5 6                               SIP1030
C           7 8 9                               SIP1040
C           .....SIP1050
100 DO 210 I=2,IN01                             SIP1060
    DO 210 J=2,JN01                             SIP1070
    N=I+DIML*(J-1)                             SIP1080
    NL=N-DIML                                   SIP1090
    NR=N+DIML                                   SIP1100
    NA=N-1                                       SIP1110
    NB=N+1                                       SIP1120
C                                           SIP1130
C    ---SKIP COMPUTATIONS IF NODE IS OUTSIDE AQUIFER BOUNDARY--- SIP1140
C    IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 210    SIP1150
C                                           SIP1160
C    ---COMPUTE COEFFICIENTS---                SIP1170
C    D=TR(NL)/DELX(J)                          SIP1180
C    F=TR(N)/DELX(J)                          SIP1190
C    B=TC(NA)/DELY(I)                         SIP1200
C    H=TC(N)/DELY(I)                         SIP1210
C    IF (EVAP.NE.CHK(6)) GO TO 120            SIP1220
C                                           SIP1230
C    ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE--- SIP1240
C    ETQB=0.                                   SIP1250
C    ETQD=0.0                                  SIP1260
C    IF (PHE(N).LE.GRND(N)=ETDIST) GO TO 120  SIP1270
C    IF (PHE(N).GT.GRND(N)) GO TO 110        SIP1280
C    ETQB=QET/ETDIST                          SIP1290
C    ETQD=ETQB*(ETDIST-GRND(N))              SIP1300
C    GO TO 120                                 SIP1310
110 ETQD=QET                                   SIP1320
C                                           SIP1330
C    ---COMPUTE STORAGE TERM---                SIP1340
C    120 IF (CONVRT.EQ.CHK(7)) GO TO 130      SIP1350
C    RHO=S(N)/DELT                             SIP1360
C    IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT     SIP1370
C    GO TO 200                                 SIP1380
C                                           SIP1390
C    ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM--- SIP1400
C    130 SUBS=0.0                              SIP1410
C    IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 170 SIP1420
C    IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 160 SIP1430
C    IF (KEEP(N)-PHE(N)) 140,150,150        SIP1440
C    140 SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N)) SIP1450
C    GO TO 170                                 SIP1460
C    150 SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N)) SIP1470
C    160 RHO=SY(N)/DELT                       SIP1480
C    GO TO 180                                 SIP1490
C    170 RHO=S(N)/DELT                       SIP1500
C    180 IF (LEAK.NE.CHK(9)) GO TO 200      SIP1510
C                                           SIP1520
C    ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION--- SIP1530
C    IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 200 SIP1540
C    HED1=AMAX1(STRT(N),TOP(N))              SIP1550
C    U=1.                                       SIP1560
C    HED2=0.                                   SIP1570
C    IF (PHE(N).GE.TOP(N)) GO TO 190        SIP1580
C    HED2=TOP(N)                              SIP1590
C    U=0.                                       SIP1600
C    190 SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N)) SIP1610
C    200 CONTINUE                             SIP1620

```



## Program listing—Continued

```

C                                     SIP1630
C   ---SIP 'NORMAL' ALGORITHM---      SIP1640
C   ---FORWARD SUBSTITUTE, COMPUTING INTERMEDIATE VECTOR V--- SIP1650
E=-B-D-F-H-RHO-TL(N)*U-ETQB          SIP1660
CH=DEL(NA)*B/(1.+W*DEL(NA))          SIP1670
GH=ETA(NL)*D/(1.+W*ETA(NL))          SIP1680
BH=B-W*CH                             SIP1690
DH=D-W*GH                             SIP1700
EH=E+W*CH+W*GH                         SIP1710
FH=F-W*CH                             SIP1720
HH=H-W*GH                             SIP1730
ALFA=BH                               SIP1740
BETA=DH                               SIP1750
GAMA=EH-ALFA*ETA(NA)-BETA*DEL(NL)     SIP1760
DEL(N)=FH/GAMA                         SIP1770
ETA(N)=HH/GAMA                         SIP1780
RES=-D*PHI(NL)-F*PHI(NR)-H*PHI(NB)-B*PHI(NA)-E*PHI(N)-RHO*KEEP(N)-SIP1790
1SL(N)=QRE(N)-WELL(N)+ETQD-SUBS-TL(N)*STRT(N) SIP1800
V(N)=(HMAX*RES-ALFA*V(NA)-BETA*V(NL))/GAMA SIP1810
210 CONTINUE                           SIP1820
C                                     SIP1830
C   ---BACK SUBSTITUTE FOR VECTOR XI--- SIP1840
DO 220 I=1,I2                          SIP1850
I3=DIML-I                               SIP1860
DO 220 J=1,J2                          SIP1870
J3=DIMW-J                               SIP1880
N=I3+DIML*(J3-1)                       SIP1890
IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 220 SIP1900
XI(N)=V(N)-DEL(N)*XI(N+DIML)-ETA(N)*XI(N+1) SIP1910
C                                     SIP1920
C   ---COMPARE MAGNITUDE OF CHANGE WITH CLOSURE CRITERION-- SIP1930
TCHK=ABS(XI(N))                         SIP1940
IF (TCHK.GT.BIGI) BIGI=TCHK             SIP1950
PHI(N)=PHI(N)+XI(N)                   SIP1960
220 CONTINUE                           SIP1970
IF (BIGI.GT.ERR) TEST=1.                SIP1980
TEST3(KOUNT+1)=BIGI                   SIP1990
IF (TEST.EQ.1.) GO TO 40               SIP2000
RETURN                                 SIP2010
C                                     SIP2020
C   .....SIP2030
C   ---ORDER EQUATIONS WITH THE LAST ROW FIRST - 3X3 EXAMPLE: SIP2040
C           7 8 9                      SIP2050
C           4 5 6                      SIP2060
C           1 2 3                      SIP2070
C   .....SIP2080
230 DO 340 II=1,I2                     SIP2090
I=DIML-II                               SIP2100
DO 340 J=2,JN01                        SIP2110
N=I+DIML*(J-1)                         SIP2120
NL=N-DIML                              SIP2130
NR=N+DIML                              SIP2140
NA=N-1                                  SIP2150
NB=N+1                                  SIP2160
C                                     SIP2170
C   ---SKIP COMPUTATIONS IF NODE IS OUTSIDE AQUIFER BOUNDARY--- SIP2180
IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 340 SIP2190
C                                     SIP2200
C   ---COMPUTE COEFFICIENTS---          SIP2210
D=TR(NL)/DELX(J)                       SIP2220
F=TR(N)/DELX(J)                        SIP2230

```

## Program listing—Continued

```

B=TC(NA)/DELY(I)
H=TC(N)/DELY(I)
IF (EVAP.NE.CHK(6)) GO TO 250
C
C ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---
ETQB=0.
ETQD=0.0
IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 250
IF (PHE(N).GT.GRND(N)) GO TO 240
ETQB=QET/ETDIST
ETQD=ETQB*(ETDIST-GRND(N))
GO TO 250
240 ETQD=QET
C
C ---COMPUTE STORAGE TERM---
250 IF (CONVRT.EQ.CHK(7)) GO TO 260
RHO=S(N)/DELT
IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT
GO TO 330
C
C ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM---
260 SUBS=0.0
IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 300
IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 290
IF (KEEP(N)-PHE(N)) 270,280,280
270 SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))
GO TO 300
280 SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))
290 RHO=SY(N)/DELT
GO TO 310
300 RHO=S(N)/DELT
310 IF (LEAK.NE.CHK(9)) GO TO 330
C
C ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION---
IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 330
HED1=AMAX1(STRT(N),TOP(N))
U=1.
HED2=0.
IF (PHE(N).GE.TOP(N)) GO TO 320
HED2=TOP(N)
U=0.
320 SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N))
330 CONTINUE
C
C ---SIP 'REVERSE' ALGORITHM---
C ---FORWARD SUBSTITUTE, COMPUTING INTERMEDIATE VECTOR V---
E=-B-D-F-H-RHO-TL(N)*U-ETQB
CH=DEL(NB)*H/(1.+W*DEL(NB))
GH=ETA(NL)*D/(1.+W*ETA(NL))
BH=H-W*CH
DH=D-W*GH
EH=E+W*CH+W*GH
FH=F-W*CH
HH=B-W*GH
ALFA=BH
BETA=DH
GAMA=EH-ALFA*ETA(NB)-BETA*DEL(NL)
DEL(N)=FH/GAMA
ETA(N)=HH/GAMA
RES=-D*PHI(NL)-F*PHI(NR)-H*PHI(NB)-B*PHI(NA)-E*PHI(N)-RHO*KEEP(N)-SIP2830
ISL(N)-QRE(N)-WELL(N)+ETQD-SUBS-TL(N)*STRT(N)
SIP2240
SIP2250
SIP2260
SIP2270
SIP2280
SIP2290
SIP2300
SIP2310
SIP2320
SIP2330
SIP2340
SIP2350
SIP2360
SIP2370
SIP2380
SIP2390
SIP2400
SIP2410
SIP2420
SIP2430
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SIP2450
SIP2460
SIP2470
SIP2480
SIP2490
SIP2500
SIP2510
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SIP2690
SIP2700
SIP2710
SIP2720
SIP2730
SIP2740
SIP2750
SIP2760
SIP2770
SIP2780
SIP2790
SIP2800
SIP2810
SIP2820
SIP2830
SIP2840

```

## Program listing—Continued

```

      V(N)=(HMAX*RES-ALFA*V(NB)-BETA*V(NL))/GAMA
340 CONTINUE
C
C      ---BACK SUBSTITUTE FOR VECTOR XI---
      DO 350 I3=2,IN01
      DO 350 J=1,J2
      J3=DIMW-J
      N=I3+DIML*(J3-1)
      IF (T(N).EQ.0.,OR.S(N).LT.0.) GO TO 350
      XI(N)=V(N)-DEL(N)*XI(N+DIML)-ETA(N)*XI(N-1)
C
C      ---COMPARE MAGNITUDE OF CHANGE WITH CLOSURE CRITERION--
      TCHK=ABS(XI(N))
      IF (TCHK.GT.BIGI) BIGI=TCHK
      PHI(N)=PHI(N)+XI(N)
350 CONTINUE
      IF (BIGI.GT.ERR) TEST=1.
      TEST3(KOUNT+1)=BIGI
      IF (TEST.EQ.1.) GO TO 40
      RETURN
C
C      .....
C
C      ---FORMATS---
C
C      -----
C
360 FORMAT ('0EXCEEDED PERMITTED NUMBER OF ITERATIONS'/' ',39('*'))
370 FORMAT ('-',44X,'SOLUTION BY THE STRONGLY IMPLICIT PROCEDURE'/45X,
143('*'),'//',61X,'BETA=',F5.2)
380 FORMAT (1H0,15,22H ITERATION PARAMETERS:,6D15.7/(/28X,6D15.7/))
      END

      SUBROUTINE SOLVE2(PHI,BE,G,TEMP,KEEP,PHE,STRT,T,S,QRE,WELL,TL,SL,DSOR 10
1EL,ETA,V,XI,DELX,BETA,DELY,ALFA,TEST3,TR,TC,GRND,SY, TOP,RATE,M,RIVSOR 20
2ER)
      -----SOR 30
C      SOLUTION BY LINE SUCCESSIVE OVERRELAXATION
C      -----SOR 40
C      -----SOR 50
C      -----SOR 60
C      -----SOR 70
C      -----SOR 80
C      SPECIFICATIONS:
      REAL *8PHI,DBLE,RHOP(20),G,BE,TEMP,IMK,DABS,W,PARAM,TEST2,DMAX1,R2SOR 90
1,A,C,B1,E,Q,RHO,B,D,F,H
      REAL *4KEEP,M
      INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,SOR 120
1CONTR,LEAK,RECH,SIP,ADI
C      -----SOR 130
C      -----SOR 140
      DIMENSION PHI(1), BE(1), G(1), TEMP(1), KEEP(1), PHE(1), STRT(1), SOR 150
1T(1), S(1), QRE(1), WELL(1), TL(1), SL(1), DEL(1), ETA(1), V(1), XSOR 160
2I(1), DELX(1), BETA(1), DELY(1), ALFA(1), TEST3(1), TR(1), TC(1), SOR 170
3GRND(1), SY(1), TOP(1), RATE(1), M(1), RIVER(1)
C      -----SOR 180
C      -----SOR 190
      COMMON /SARRAY/ VF4(11),CHK(15)
      -----SOR 200
      COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,ERROR,LESOR 210
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,SOR 220
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,SOR 230
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DISOR 240
4MW,JN01,IN01,R,P,PU,I,J,IDX1,IDX2
      RETURN
      -----SOR 250
      -----SOR 260

```

## Program listing—Continued

```

C .....SOR 270
C .....SOR 280
C ---WRITE ACCELERATION PARAMETER---SOR 290
C *****SOR 300
C ENTRY ITER2SOR 310
C *****SOR 320
C WRITE (P,490)SOR 330
C WRITE (P,500) HMAX,LENGTHSOR 340
C RETURNSOR 350
C .....SOR 360
C .....SOR 370
C ---INITIALIZE DATA FOR A NEW ITERATION---SOR 380
10 KOUNT=KOUNT+1SOR 390
IF (KOUNT.LE.ITMAX) GO TO 20SOR 400
WRITE (P,510)SOR 410
CALL TERM1SOR 420
C *****SOR 430
C ENTRY NEWITBSOR 440
C *****SOR 450
20 TEST3(KOUNT+1)=0.SOR 460
TEST=0.SOR 470
N=DIML*DIMWSOR 480
DO 30 I=1,NSOR 490
30 PHE(I)=PHI(I)SOR 500
BIGI=0.0SOR 510
C .....SOR 520
C ---COMPUTE TRANSMISSIVITY AND T COEFFICIENTS IN WATER TABLESOR 530
C OR WATER TABLE=ARTESIAN SIMUATION---SOR 540
C IF (WATER.NE.CHK(2)) GO TO 40SOR 550
C CALL TRANSOR 560
C .....SOR 570
C .....SOR 580
C ---SOLUTION BY LSOR---SOR 590
C -----SOR 600
40 NO3=DIMW-2SOR 610
TEMP(DIMW)=0.0SOR 620
DO 170 I=2,IN01SOR 630
DO 150 J=2,JN01SOR 640
N=I+DIML*(J-1)SOR 650
NA=N-1SOR 660
NB=N+1SOR 670
NL=N-DIMLSOR 680
NR=N+DIMLSOR 690
BE(J)=0.0SOR 700
G(J)=0.0SOR 710
C .....SOR 720
C ---SKIP COMPUTATIONS IF NODE IS OUTSIDE AQUIFER BOUNDARY---SOR 730
C IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 150SOR 740
C .....SOR 750
C ---COMPUTE COEFFICIENTS---SOR 760
C D=TR(N-DIML)/DELX(J)SOR 770
C F=TR(N)/DELX(J)SOR 780
C B=TC(N-1)/DELY(I)SOR 790
C H=TC(N)/DELY(I)SOR 800
C IF (EVAP.NE.CHK(6)) GO TO 60SOR 810
C .....SOR 820
C ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---SOR 830
C ETQB=0.SOR 840
C ETQD=0.0SOR 850
C IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 60SOR 860
C IF (PHE(N).GT.GRND(N)) GO TO 50SOR 870

```

## Program listing—Continued

|     |  |         |
|-----|--|---------|
|     | ETQB=QET/ETDIST  | SOR 880 |
|     | ETQD=ETQB*(ETDIST-GRND(N))   | SOR 890 |
|     | GO TO 60   | SOR 900 |
| 50  | ETQD=QET   | SOR 910 |
| C   |  | SOR 920 |
| C   | ---COMPUTE STORAGE TERM---   | SOR 930 |
| 60  | IF (CONVRT.EQ.CHK(7)) GO TO 70                                     | SOR 940 |
|     | RHO=S(N)/DELT  | SOR 950 |
|     | IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT                                | SOR 960 |
|     | GO TO 140  | SOR 970 |
| C   |  | SOR 980 |
| C   | ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM---           | SOR 990 |
| 70  | SUBS=0.0   | SOR1000 |
|     | IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 110              | SOR1010 |
|     | IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 100              | SOR1020 |
|     | IF (KEEP(N)-PHE(N)) 80,90,90                                       | SOR1030 |
| 80  | SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))                            | SOR1040 |
|     | GO TO 110  | SOR1050 |
| 90  | SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))                            | SOR1060 |
| 100 | RHO=SY(N)/DELT   | SOR1070 |
|     | GO TO 120  | SOR1080 |
| 110 | RHO=S(N)/DELT  | SOR1090 |
| 120 | IF (LEAK.NE.CHK(9)) GO TO 140                                      | SOR1100 |
| C   |  | SOR1110 |
| C   | ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION---           | SOR1120 |
|     | IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 140                         | SOR1130 |
|     | HED1=AMAX1(STRT(N),TOP(N))   | SOR1140 |
|     | U=1.   | SOR1150 |
|     | HED2=0.  | SOR1160 |
|     | IF (PHE(N).GE.TOP(N)) GO TO 130                                    | SOR1170 |
|     | HED2=TOP(N)  | SOR1180 |
|     | U=0.   | SOR1190 |
| 130 | SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N))       | SOR1200 |
| 140 | CONTINUE   | SOR1210 |
| C   |  | SOR1220 |
| C   | ---FORWARD SUBSTITUTE, COMPUTING INTERMEDIATE VECTOR G---          | SOR1230 |
|     | E=-D-F-B-H-RHO-TL(N)*U-ETQB  | SOR1240 |
|     | W=E-D*BE(J-1)  | SOR1250 |
|     | BE(J)=F/W  | SOR1260 |
|     | Q=-B*PHI(NA)-H*PHI(NB)-RHO*KEEP(N)-SL(N)-QRE(N)-WELL(N)+ETQD-SUBS- | SOR1270 |
|     | 1TL(N)*STRT(N)-D*PHI(NL)-F*PHI(NR)-E*PHI(N)                        | SOR1280 |
|     | G(J)=(Q-D*G(J-1))/W  | SOR1290 |
| 150 | CONTINUE   | SOR1300 |
| C   |  | SOR1310 |
| C   | ---BACK SUBSTITUTE FOR TEMP---                                     | SOR1320 |
|     | DO 160 KNO4=1,NO3  | SOR1330 |
|     | NO4=DIMW-KNO4  | SOR1340 |
|     | TEMP(NO4)=G(NO4)-BE(NO4)*TEMP(NO4+1)                               | SOR1350 |
| 160 | CONTINUE   | SOR1360 |
| C   |  | SOR1370 |
| C   | ---EXTRAPOLATED VALUE OF PHI---                                    | SOR1380 |
|     | DO 170 J=2,JNO1  | SOR1390 |
|     | N=I+DIML*(J-1)   | SOR1400 |
|     | PHI(N)=PHI(N)+HMAX*TEMP(J)   | SOR1410 |
| C   |  | SOR1420 |
| C   | ---COMPARE DIFFERENCE WITH CLOSURE CRITERION--                     | SOR1430 |
|     | TCHK=DABS(TEMP(J))   | SOR1440 |
|     | IF (TCHK.GT.BIGI) BIGI=TCHK  | SOR1450 |
| 170 | CONTINUE   | SOR1460 |
|     | IF (BIGI.GT.ERR) TEST=1.   | SOR1470 |
|     | TEST3(KOUNT+1)=BIGI  | SOR1480 |

## Program listing—Continued

```

IF (KOUNT.EQ.0) GO TO 10
IF (TEST.EQ.0.) RETURN
C
C ---TEST FOR TWO DIMENSIONAL CORRECTION---
IF (MOD(KOUNT,LENGTH).NE.0) GO TO 10
GO TO 200
180 DO 190 I=2,IN01
DO 190 J=2,JN01
N=I+DIML*(J-1)
IF (T(N).EQ.0.) GO TO 190
PHI(N)=PHI(N)+ALFA(I)+BETA(J)
190 CONTINUE
GO TO 10
C
C .....
C
C ---TWO DIMENSIONAL CORRECTION TO LSOR---
C -----
C
C ---COMPUTE ALFA CORRECTION FOR ROWS---
200 DO 210 I=1,DIML
ALFA(I)=0.
BE(I)=0.0
210 G(I)=0.0
DO 330 I=2,IN01
A=0.
B2=0.
C=0.
Q=0.
C
C ---SUMMATION OF COEFFICIENTS FOR EACH ROW---
DO 320 J=2,JN01
N=I+DIML*(J-1)
NA=N-1
NB=N+1
NL=N-DIML
NR=N+DIML
IF (S(N).LT.0.) GO TO 330
IF (T(N).EQ.0.) GO TO 320
C
C ---COMPUTE COEFFICIENTS---
D=TR(N-DIML)/DELX(J)
F=TR(N)/DELX(J)
B=TC(N-1)/DELY(I)
H=TC(N)/DELY(I)
IF (EVAP.NE.CHK(6)) GO TO 230
C
C ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---
ETQB=0.
ETQD=0.0
IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 230
IF (PHE(N).GT.GRND(N)) GO TO 220
ETQB=QET/ETDIST
ETQD=ETQB*(ETDIST-GRND(N))
GO TO 230
220 ETQD=QET
C
C ---COMPUTE STORAGE TERM---
230 IF (CONVRT.EQ.CHK(7)) GO TO 240
RHO=S(N)/DELT
IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT
GO TO 310

```

SOR1490  
SOR1500  
SOR1510  
SOR1520  
SOR1530  
SOR1540  
SOR1550  
SOR1560  
SOR1570  
SOR1580  
SOR1590  
SOR1600  
SOR1610  
SOR1620  
SOR1630  
SOR1640  
SOR1650  
SOR1660  
SOR1670  
SOR1680  
SOR1690  
SOR1700  
SOR1710  
SOR1720  
SOR1730  
SOR1740  
SOR1750  
SOR1760  
SOR1770  
SOR1780  
SOR1790  
SOR1800  
SOR1810  
SOR1820  
SOR1830  
SOR1840  
SOR1850  
SOR1860  
SOR1870  
SOR1880  
SOR1890  
SOR1900  
SOR1910  
SOR1920  
SOR1930  
SOR1940  
SOR1950  
SOR1960  
SOR1970  
SOR1980  
SOR1990  
SOR2000  
SOR2010  
SOR2020  
SOR2030  
SOR2040  
SOR2050  
SOR2060  
SOR2070  
SOR2080  
SOR2090

## Program listing—Continued

```

240 SUBS=0.0
IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 280
IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 270
IF (KEEP(N)-PHE(N)) 250,260,260
250 SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))
GO TO 280
260 SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))
270 RHO=SY(N)/DELT
GO TO 290
280 RHO=S(N)/DELT
290 IF (LEAK.NE.CHK(9)) GO TO 310
C
C ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION---
IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 310
HED1=AMAX1(STRT(N),TOP(N))
U=1.
HED2=0.
IF (PHE(N).GE.TOP(N)) GO TO 300
HED2=TOP(N)
U=0.
300 SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N))
310 CONTINUE
C
A=A-B
B1=B+H+RHO+TL(N)*U+ETQB
B2=B2+B1
C=C-H
Q=Q+(D*PHI(NL)+F*PHI(NR)+B*PHI(NA)+H*PHI(NB)+RHO*KEEP(N)+SL(N)+QRE
1(N)+WELL(N)-ETGD+SUBS+TL(N)*STRT(N)-(D+F+B1)*PHI(N))
320 CONTINUE
C
C ---COMPUTATION OF INTERMEDIATE VECTOR G---
W=B2-A*BE(I-1)
BE(I)=C/W
G(I)=(Q-A*G(I-1))/W
330 CONTINUE
C
C ---BACK SUBSTITUTE FOR ALFA---
NO3=DIML-2
DO 340 KNO4=1,NO3
NO4=DIML-KNO4
340 ALFA(NO4)=G(NO4)-BE(NO4)*ALFA(NO4+1)
*****
C
C ---COMPUTE BETA CORRECTION FOR COLUMNS---
DO 350 J=1,DIMW
BETA(J)=0.
BE(J)=0.0
350 G(J)=0.0
DO 470 J=2,JNO1
A=0.
B2=0.
C=0.
Q=0.
C
C ---SUMMATION OF COEFFICIENTS FOR EACH COLUMN---
DO 460 I=2,INO1
N=I+DIML*(J-1)
NA=N-1
NB=N+1
NL=N-DIML

```

```

SOR2100
SOR2110
SOR2120
SOR2130
SOR2140
SOR2150
SOR2160
SOR2170
SOR2180
SOR2190
SOR2200
SOR2210
SOR2220
SOR2230
SOR2240
SOR2250
SOR2260
SOR2270
SOR2280
SOR2290
SOR2300
SOR2310
SOR2320
SOR2330
SOR2340
SOR2350
SOR2360
SOR2370
SOR2380
SOR2390
SOR2400
SOR2410
SOR2420
SOR2430
SOR2440
SOR2450
SOR2460
SOR2470
SOR2480
SOR2490
SOR2500
SOR2510
SOR2520
SOR2530
SOR2540
SOR2550
SOR2560
SOR2570
SOR2580
SOR2590
SOR2600
SOR2610
SOR2620
SOR2630
SOR2640
SOR2650
SOR2660
SOR2670
SOR2680
SOR2690
SOR2700

```

## Program listing—Continued

```

NR=N*DIML
IF (S(N).LT.0.) GO TO 470
IF (T(N).EQ.0.) GO TO 460
D=TR(N-DIML)/DELX(J)
F=TR(N)/DELX(J)
B=TC(N-1)/DELY(I)
H=TC(N)/DELY(I)
IF (EVAP.NE.CHK(6)) GO TO 370
C
C ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---
ETQB=0.
ETQD=0.0
IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 370
IF (PHE(N).GT.GRND(N)) GO TO 360
ETQB=QET/ETDIST
ETQD=ETQB*(ETDIST-GRND(N))
GO TO 370
360 ETQD=QET
C
C ---COMPUTE STORAGE TERM---
370 IF (CONVRT.EQ.CHK(7)) GO TO 380
RHO=S(N)/DELT
IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT
GO TO 450
C
C ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM---
380 SUBS=0.0
IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 420
IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 410
IF (KEEP(N)-PHE(N)) 390,400,400
390 SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))
GO TO 420
400 SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))
410 RHO=SY(N)/DELT
GO TO 430
420 RHO=S(N)/DELT
430 IF (LEAK.NE.CHK(9)) GO TO 450
C
C ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION---
IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 450
HED1=AMAX1(STRT(N),TOP(N))
U=1.
HED2=0.
IF (PHE(N).GE.TOP(N)) GO TO 440
HED2=TOP(N)
U=0.
440 SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)*TL(N)*(HED1-HED2-STRT(N))
450 CONTINUE
C
A=A-D
B1=D+F+RHO*TL(N)*U+ETQB
B2=B2+B1
C=C-F
Q=Q+(D*PHI(NL)+F*PHI(NR)+B*PHI(NA)+H*PHI(NB)+RHO*KEEP(N)+SL(N)+QRE
1(N)+WELL(N)-ETQD+SUBS+TL(N)*STRT(N)-(B+H+B1)*PHI(N))
460 CONTINUE
C
C ---COMPUTATION OF INTERMEDIATE VECTOR G---
W=B2-A*BE(J-1)
BE(J)=C/W
G(J)=(Q-A*G(J-1))/W

```

```

SOR2710
SOR2720
SOR2730
SOR2740
SOR2750
SOR2760
SOR2770
SOR2780
SOR2790
SOR2800
SOR2810
SOR2820
SOR2830
SOR2840
SOR2850
SOR2860
SOR2870
SOR2880
SOR2890
SOR2900
SOR2910
SOR2920
SOR2930
SOR2940
SOR2950
SOR2960
SOR2970
SOR2980
SOR2990
SOR3000
SOR3010
SOR3020
SOR3030
SOR3040
SOR3050
SOR3060
SOR3070
SOR3080
SOR3090
SOR3100
SOR3110
SOR3120
SOR3130
SOR3140
SOR3150
SOR3160
SOR3170
SOR3180
SOR3190
SOR3200
SOR3210
SOR3220
SOR3230
SOR3240
SOR3250
SOR3260
SOR3270
SOR3280
SOR3290
SOR3300
SOR3310

```



## Program listing—Continued

```

470 CONTINUE                                SOR3320
C                                             SOR3330
C ---BACK SUBSTITUTE FOR BETA---            SOR3340
  NO3=DIMW-2                                SOR3350
  DO 480 KNO4=1,NO3                          SOR3360
  NO4=DIMW-KNO4                              SOR3370
480 BETA(NO4)=G(NO4)-BE(NO4)*BETA(NO4+1)    SOR3380
  GO TO 180                                  SOR3390
C .....                                SOR3400
C .....                                SOR3410
C ---FORMATS---                             SOR3420
C .....                                SOR3430
C .....                                SOR3440
C .....                                SOR3450
C .....                                SOR3460
490 FORMAT ('-',45X,'SOLUTION BY LINE SUCCESSIVE OVERRELAXATION',/46X,4SOR3470
  12(' ',' '))                              SOR3480
500 FORMAT ('-',26X,'ACCELERATION PARAMETER =',F6.3,' TWO DIMENSIONALSOR3490
  1 CORRECTION EVERY',I5,' ITERATIONS')     SOR3500
510 FORMAT ('0EXCEEDED PERMITTED NUMBER OF ITERATIONS'/ ' ',39(' ',' ')) SOR3510
  END                                         SOR3520-
  SUBROUTINE SOLVE3(PHI,BE,G,TEMP,KEEP,PHE,STRT,T,S,QRE,WELL,TL,SL,DADI 10
  1EL,ETA,V,XI,DELX,BETA,DELY,ALFA,XII,TEST3,TR,TC,GRND,SY, TOP,RATE,MADI 20
  2,RIVER)                                  ADI 30
C -----AD I 40
C SOLUTION BY THE ALTERNATING DIRECTION IMPLICIT PROCEDURE ADI 50
C -----AD I 60
C -----AD I 70
C SPECIFICATIONS: ADI 80
  REAL *8PHI,DBLE,RHOP(20),G,BE,TEMP,IMK,DABS,W,PARAM,TEST2,DMAX1,DTADI 90
  1ERMS,B1,E,Q,B,D,F,H,RHO,XII ADI 100
  REAL *4KEEP,M ADI 110
  INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,ADI 120
  1CONTR,LEAK,RECH,SIP,ADI ADI 130
C -----AD I 140
  DIMENSION PHI(1), BE(1), G(1), TEMP(1), KEEP(1), PHE(1), STRT(1), ADI 150
  1T(1), S(1), QRE(1), WELL(1), TL(1), SL(1), DEL(1), ETA(1), V(1), XADI 160
  2I(1), DELX(1), BETA(1), DELY(1), ALFA(1), XII(1), TEST3(1), TR(1),ADI 170
  3 TC(1), GRND(1), SY(1), TOP(1), RATE(1), M(1), RIVER(1) ADI 180
C -----AD I 190
  COMMON /SARRAY/ VF4(11),CHK(15) ADI 200
  COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,EROR,LEADI 210
  1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,COLT,HMAX,YDIM,WIDTH,ADI 220
  2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,ADI 230
  3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DIADI 240
  4MW,JNO1,INO1,R,P,PU,I,J,IDX1,IDX2 ADI 250
  RETURN ADI 260
C .....AD I 270
C .....AD I 280
C ---COMPUTE AND PRINT ITERATION PARAMETERS--- ADI 290
C ***** ADI 300
C ENTRY ITER3 ADI 310
C ***** ADI 320
  HMIN=2. ADI 330
  IN4=DIMW-2 ADI 340
  IN5=DIML-2 ADI 350
  XVAL=3.1415**2/(2.*IN4*IN4) ADI 360
  YVAL=3.1415**2/(2.*IN5*IN5) ADI 370
  DO 10 I=2,INO1 ADI 380

```

## Program listing—Continued

```

DO 10 J=2,JN01                                ADI 390
N=I+DIML*(J-1)                                ADI 400
IF (T(N).EQ.0.) GO TO 10                      ADI 410
XPART=XVAL*(1/(1+DELX(J)**2*FACTY/DELY(I)**2*FACTX)) ADI 420
YPART=YVAL*(1/(1+DELY(I)**2*FACTX/DELX(J)**2*FACTY)) ADI 430
HMIN=AMIN1(HMIN,XPART,YPART)                  ADI 440
10 CONTINUE                                    ADI 450
ALPHA=EXP(ALOG(HMAX/HMIN)/(LENGTH-1))         ADI 460
RHOP(1)=HMIN                                  ADI 470
DO 20 NTIME=2,LENGTH                           ADI 480
20 RHOP(NTIME)=RHOP(NTIME-1)*ALPHA            ADI 490
WRITE (P,400)                                  ADI 500
WRITE (P,410) LENGTH,(RHOP(J),J=1,LENGTH)     ADI 510
RETURN                                          ADI 520
C .....ADJ 530
C .....ADJ 540
C ---INITIALIZE DATA FOR A NEW ITERATION---   ADI 550
30 KOUNT=KOUNT+1                               ADI 560
IF (KOUNT.LE.ITMAX) GO TO 40                   ADI 570
WRITE (P,390)                                  ADI 580
CALL TERM1                                     ADI 590
40 IF (MOD(KOUNT,LENGTH)) 50,50,60            ADI 600
*****ADJ 610
C ENTRY NEWITC                                  ADI 620
*****ADJ 630
C .....ADJ 640
50 NTH=0                                       ADI 650
60 NTH=NTH+1                                   ADI 660
PARAM=RHOP(NTH)                               ADI 670
TEST3(KOUNT+1)=0.                             ADI 680
TEST=0.                                        ADI 690
N=DIML*DIMW                                   ADI 700
DO 70 I=1,N                                   ADI 710
70 PHE(I)=PHI(I)                              ADI 720
BIGI=0.0                                       ADI 730
C .....ADJ 740
C ---COMPUTE TRANSMISSIVITY AND T COEFFICIENTS IN WATER TABLE ADI 750
C OR WATER TABLE-ARTESIAN SIMUATION---       ADI 760
IF (WATER.NE.CHK(2)) GO TO 80                 ADI 770
CALL TRANS                                     ADI 780
C .....ADJ 790
C .....ADJ 800
C ---SOLUTION BY ADI---                         ADI 810
C -----ADJ 820
C ---COMPUTE IMPLICITLY ALONG ROWS---          ADI 830
80 N03=DIMW-2                                  ADI 840
DO 90 J=1,DIMW                                ADI 850
N=1+DIML*(J-1)                                ADI 860
90 TEMP(J)=PHI(N)                              ADI 870
DO 230 I=2,DIML                               ADI 880
DO 200 J=2,JN01                               ADI 890
N=I+DIML*(J-1)                                ADI 900
NA=N-1                                         ADI 910
NB=N+1                                         ADI 920
NL=N-DIML                                     ADI 930
NR=N+DIML                                     ADI 940
BE(J)=0.0                                     ADI 950
G(J)=0.0                                       ADI 960
C .....ADJ 970
C ---SKIP COMPUTATIONS IF NODE IS OUTSIDE AQUIFER BOUNDARY--- ADI 980
IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 200      ADI 990
C .....ADJ 990

```

## Program listing—Continued

```

C      ---COMPUTE COEFFICIENTS---
      D=TR(N-DIML)/DELX(J)
      F=TR(N)/DELX(J)
      B=TC(N-1)/DELY(I)
      H=TC(N)/DELY(I)
      IF (EVAP.NE.CHK(6)) GO TO 110
C
C      ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---
      ETQB=0.
      ETQD=0.0
      IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 110
      IF (PHE(N).GT.GRND(N)) GO TO 100
      ETQB=QET/ETDIST
      ETQD=ETQB*(ETDIST-GRND(N))
      GO TO 110
100  ETQD=QET
C
C      ---COMPUTE STORAGE TERM---
110  IF (CONVRT.EQ.CHK(7)) GO TO 120
      RHO=S(N)/DELT
      IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT
      GO TO 190
C
C      ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM---
120  SUBS=0.0
      IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 160
      IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 150
      IF (KEEP(N)=PHE(N)) 130,140,140
130  SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))
      GO TO 160
140  SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))
150  RHO=SY(N)/DELT
      GO TO 170
160  RHO=S(N)/DELT
170  IF (LEAK.NE.CHK(9)) GO TO 190
C
C      ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION---
      IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 190
      HED1=AMAX1(STRT(N),TOP(N))
      U=1.
      HED2=0.
      IF (PHE(N).GE.TOP(N)) GO TO 180
      HED2=TOP(N)
      U=0.
180  SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N))
190  CONTINUE
C
C      ---CALCULATE VALUES FOR PARAMETERS USED IN THOMAS ALGORITHM
      AND FORWARD SUBSTITUTE TO COMPUTE INTERMEDIATE VECTOR G---
      IMK=(B+D+F+H)*PARAM
      E=-D-F-RHO-IMK-TL(N)*U-ETQB
      W=E-D*BE(J-1)
      BE(J)=F/W
      Q=-B*PHI(NA)+(B+H-IMK-E)*PHI(N)-H*PHI(NB)-RHO*KEEP(N)-SL(N)-QRE(N)
1  -WELL(N)+ETQD-SUBS-TL(N)*STRT(N)-D*PHI(NL)-F*PHI(NR)
      G(J)=(Q-D*G(J-1))/W
200  CONTINUE
C
C      ---BACK SUBSTITUTE FOR HEAD VALUES AND PLACE THEM IN TEMP---
      XII(DIMW)=0.00
      DO 220 KNO4=1,N03

```

```

ADI1000
ADI1010
ADI1020
ADI1030
ADI1040
ADI1050
ADI1060
ADI1070
ADI1080
ADI1090
ADI1100
ADI1110
ADI1120
ADI1130
ADI1140
ADI1150
ADI1160
ADI1170
ADI1180
ADI1190
ADI1200
ADI1210
ADI1220
ADI1230
ADI1240
ADI1250
ADI1260
ADI1270
ADI1280
ADI1290
ADI1300
ADI1310
ADI1320
ADI1330
ADI1340
ADI1350
ADI1360
ADI1370
ADI1380
ADI1390
ADI1400
ADI1410
ADI1420
ADI1430
ADI1440
ADI1450
ADI1460
ADI1470
ADI1480
ADI1490
ADI1500
ADI1510
ADI1520
ADI1530
ADI1540
ADI1550
ADI1560
ADI1570
ADI1580
ADI1590
ADI1600

```

## Program listing—Continued

```

      NO4=DIMW-KNO4
      N=I+DIML*(NO4-1)
C
C      ---FIRST PLACE TEMP VALUES IN PHI(N-1)---
      PHI(N-1)=TEMP(NO4)
      IF (T(N).NE.0..AND.S(N).GE.0.) GO TO 210
      XII(NO4)=0.00
      GO TO 220
210  XII(NO4)=G(NO4)-BE(NO4)*XII(NO4+1)
220  TEMP(NO4)=PHI(N)+XII(NO4)
230  CONTINUE
C
C      .....
C
C      ---COMPUTE IMPLICITLY ALONG COLUMNS---
      NO3=DIML-2
      DO 240 I=1,DIML
240  TEMP(I)=PHI(I)
      DO 380 J=2,DIMW
      DO 350 I=2,IN01
      N=I+DIML*(J-1)
      NA=N-1
      NB=N+1
      NL=N-DIML
      NR=N+DIML
      BE(I)=0.0
      G(I)=0.0
C
C      ---SKIP COMPUTATIONS IF NODE IS OUTSIDE AQUIFER BOUNDARY---
      IF (T(N).EQ.0..OR.S(N).LT.0.) GO TO 350
C
C      ---COMPUTE COEFFICIENTS---
      D=TR(N-DIML)/DELX(J)
      F=TR(N)/DELX(J)
      B=TC(N-1)/DELY(I)
      H=TC(N)/DELY(I)
      IF (EVAP.NE.CHK(6)) GO TO 260
C
C      ---COMPUTE EXPLICIT AND IMPLICIT PARTS OF ET RATE---
      ETQB=0.
      ETQD=0.0
      IF (PHE(N).LE.GRND(N)-ETDIST) GO TO 260
      IF (PHE(N).GT.GRND(N)) GO TO 250
      ETQB=GET/ETDIST
      ETQD=ETQB*(ETDIST-GRND(N))
      GO TO 260
250  ETQD=GET
C
C      ---COMPUTE STORAGE TERM---
260  IF (CONVRT.EQ.CHK(7)) GO TO 270
      RHO=S(N)/DELT
      IF (WATER.EQ.CHK(2)) RHO=SY(N)/DELT
      GO TO 340
C
C      ---COMPUTE STORAGE COEFFICIENT FOR CONVERSION PROBLEM---
270  SUBS=0.0
      IF (KEEP(N).GE.TOP(N).AND.PHE(N).GE.TOP(N)) GO TO 310
      IF (KEEP(N).LT.TOP(N).AND.PHE(N).LT.TOP(N)) GO TO 300
      IF (KEEP(N)-PHE(N)) 280,290,290
280  SUBS=(SY(N)-S(N))/DELT*(KEEP(N)-TOP(N))
      GO TO 310
290  SUBS=(S(N)-SY(N))/DELT*(KEEP(N)-TOP(N))

```

```

      ADI1610
      ADI1620
      ADI1630
      ADI1640
      ADI1650
      ADI1660
      ADI1670
      ADI1680
      ADI1690
      ADI1700
      ADI1710
      ADI1720
      ADI1730
      ADI1740
      ADI1750
      ADI1760
      ADI1770
      ADI1780
      ADI1790
      ADI1800
      ADI1810
      ADI1820
      ADI1830
      ADI1840
      ADI1850
      ADI1860
      ADI1870
      ADI1880
      ADI1890
      ADI1900
      ADI1910
      ADI1920
      ADI1930
      ADI1940
      ADI1950
      ADI1960
      ADI1970
      ADI1980
      ADI1990
      ADI2000
      ADI2010
      ADI2020
      ADI2030
      ADI2040
      ADI2050
      ADI2060
      ADI2070
      ADI2080
      ADI2090
      ADI2100
      ADI2110
      ADI2120
      ADI2130
      ADI2140
      ADI2150
      ADI2160
      ADI2170
      ADI2180
      ADI2190
      ADI2200
      ADI2210

```

## Program listing—Continued

```

300 RHO=SY(N)/DELT                                ADI2220
    GO TO 320                                      ADI2230
310 RHO=S(N)/DELT                                  ADI2240
320 IF (LEAK.NE.CHK(9)) GO TO 340                  ADI2250
C                                                    ADI2260
C    ---COMPUTE NET LEAKAGE TERM FOR CONVERSION SIMULATION--- ADI2270
    IF (RATE(N).EQ.0..OR.M(N).EQ.0.) GO TO 340    ADI2280
    HED1=AMAX1(STRT(N),TOP(N))                    ADI2290
    U=1.                                           ADI2300
    HED2=0.                                         ADI2310
    IF (PHE(N).GE.TOP(N)) GO TO 330               ADI2320
    HED2=TOP(N)                                    ADI2330
    U=0.                                           ADI2340
330 SL(N)=RATE(N)/M(N)*(RIVER(N)-HED1)+TL(N)*(HED1-HED2-STRT(N)) ADI2350
340 CONTINUE                                       ADI2360
C                                                    ADI2370
C    ---CALCULATE VALUES FOR PARAMETERS USED IN THOMAS ALGORITHM ADI2380
C    AND FORWARD SUBSTITUTE TO COMPUTE INTERMEDIATE VECTOR G--- ADI2390
C    IMK=(B+D+F+H)*PARAM                           ADI2400
    E=-B-H-RHO-IMK-TL(N)*U-ETQB                   ADI2410
    W=E-B*BE(I-1)                                  ADI2420
    BE(I)=H/W                                       ADI2430
    Q=-D*PHI(NL)+(D+F-IMK-E)*PHI(N)-F*PHI(NR)-RHO*KEEP(N)-SL(N)-GRE(N) ADI2440
    I=WELL(N)+ETQD-SUBS-TL(N)*STRT(N)-B*PHI(NA)-H*PHI(NB) ADI2450
    G(I)=(Q-B*G(I-1))/W                             ADI2460
350 CONTINUE                                       ADI2470
C                                                    ADI2480
C    ---BACK SUBSTITUTE FOR HEAD VALUES AND PLACE THEM IN TEMP--- ADI2490
C    XII(DIML)=0.D0                                 ADI2500
    DO 370 KNO4=1,N03                              ADI2510
    N04=DIML-KNO4                                  ADI2520
    N=N04+DIML*(J-1)                               ADI2530
C                                                    ADI2540
C    ---FIRST PLACE TEMP VALUES IN PHI(N-DIML)---- ADI2550
C    PHI(N-DIML)=TEMP(N04)                         ADI2560
    IF (T(N).NE.0..AND.S(N).GE.0.) GO TO 360      ADI2570
    XII(N04)=0.D0                                  ADI2580
    TEMP(N04)=PHI(N)                              ADI2590
    GO TO 370                                       ADI2600
360 XII(N04)=G(N04)-BE(N04)*XII(N04+1)           ADI2610
    TEMP(N04)=PHI(N)+XII(N04)                     ADI2620
C                                                    ADI2630
C    ---COMPARE CHANGE IN HEAD WITH CLOSURE CRITERION-- ADI2640
C    TCHK=ABS(SNGL(TEMP(N04))-PHE(N))              ADI2650
    IF (TCHK.GT.BIGI) BIGI=TCHK                    ADI2660
370 CONTINUE                                       ADI2670
380 CONTINUE                                       ADI2680
    IF (BIGI.GT.ERR) TEST=1.                       ADI2690
    TEST3(KOUNT+1)=BIGI                            ADI2700
    IF (TEST.EQ.1.) GO TO 30                       ADI2710
    RETURN                                          ADI2720
C    .....ADI2730
C                                                    ADI2740
C    ---FORMATS---                                  ADI2750
C                                                    ADI2760
C    -----ADI2770
C                                                    ADI2780
C                                                    ADI2790
390 FORMAT ('0EXCEEDED PERMITTED NUMBER OF ITERATIONS/' ' ',39('*')) ADI2800
400 FORMAT ('-' ,38X,'SOLUTION BY THE ALTERNATING DIRECTION IMPLICIT PRADI2810
    IOCEDURE'/39X,56('_'))                          ADI2820

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## Program listing—Continued

```

410 FORMAT (///1H0,I5,22H ITERATION PARAMETERS:,8D12.3//28X,10D12.3) ADI2830
      END ADI2840-

      SUBROUTINE COEF(PHI,KEEP,PHE,STRT,SURI,T,TR,TC,S,WELL,TL,SL,PERM,BCOF 10
10TTOM,SY,RATE,RIVER,M,TOP,GRND,DELX,DELY) COF 20
      -----COF 30
      C COMPUTE COEFFICIENTS COF 40
      C -----COF 50
      C COF 60
      C SPECIFICATIONS: COF 70
      REAL *8PHI,DBLE,RHO,B,D,F,H COF 80
      REAL *4KEEP,M COF 90
      INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,COF 100
1CCNTR,LEAK,RECH,SIP,ADI COF 110
      C COF 120
      DIMENSION PHI(1), KEEP(1), PHE(1), STRT(1), SURI(1), T(1), TR(1), COF 130
1TC(1), S(1), WELL(1), TL(1), SL(1), PERM(1), BOTTOM(1), SY(1), RATCOF 140
2E(1), RIVER(1), M(1), TOP(1), GRND(1), DELX(1), DELY(1) COF 150
      C COF 160
      COMMON /SARRAY/ VF4(11),CHK(15) COF 170
      COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,EROR,LECOF 180
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,QET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,COF 190
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SURS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,COF 200
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DICOF 210
4MW,JNO1,INO1,R,P,PU,I,J,IDK1,IDK2 COF 220
      C COF 230
      DATA PIE/3.141593/ COF 235
      RETURN COF 240
      C .....COF 250
      C COF 260
      C ---COMPUTE COEFFICIENTS FOR TRANSIENT PART OF LEAKAGE TERM--- COF 270
      C ***** COF 280
      C ENTRY CLAY COF 290
      C ***** COF 300
      TMIN=1.E30 COF 310
      TT=0.0 COF 320
      PRATE=0. COF 330
      DO 50 I=1,DIML COF 340
      DO 50 J=1,DIMW COF 350
      N=I+DIML*(J-1) COF 360
      C COF 370
      C ---SKIP COMPUTATIONS IF T, RATE OR M = 0, OR IF CONSTANT COF 380
      C HEAD BOUNDARY--- COF 390
      C IF (RATE(N).LE.0..OR.T(N).EQ.0..OR.M(N).EQ.0..OR.S(N).LT.0.) GO TO COF 400
1 50 COF 410
      C COF 420
      C ---IF VALUE FOR TL(N ) WILL EQUAL VALUE FOR PREVIOUS NODE, COF 430
      C SKIP PART OF COMPUTATIONS--- COF 440
      IF (RATE(N)*M(N).EQ.PRATE) GO TO 40 COF 450
      DIMT=RATE(N)*SUMP/(M(N)*M(N)*SS*3) COF 460
      IF (DIMT.GT.TT) TT=DIMT COF 470
      IF (DIMT.LT.TMIN) TMIN=DIMT COF 480
      PPT=PIE*PIE*DIMT COF 490
      C COF 500
      C ---RECOMPUTE PPT IF DIMT WITHIN RANGE FOR SHORT TIME COMPUTATION---COF 510
      IF (DIMT.LT.1.0E-03) PPT=1.0/DIMT COF 520
      CC=(2.3-PPT)/(2.*PPT) COF 530
      C COF 540
      C ---COMPUTE SUM OF EXPONENTIALS--- COF 550
      SUMN=0.0 COF 560

```

## Program listing—Continued

```

DO 20 K=1,200                                COF 570
POWER=K*K*PPT                                COF 580
IF (POWER.LE.150.) GO TO 10                   COF 590
POWER=150                                     COF 600
10 PEX=EXP(-POWER)                            COF 610
SUMN=SUMN+PEX                                 COF 620
IF (PEX.GT.0.00009) GO TO 20                 COF 630
IF (K.GT.CC) GO TO 30                         COF 640
20 CONTINUE                                   COF 650
C                                              COF 660
C ---COMPUTE DENOMINATER DEPENDING ON VALUE OF DIMT--- COF 670
30 DENOM=1.0                                  COF 680
IF (DIMT.LT.1.0E-03) DENOM=SQRT(PIE*DIMT)    COF 690
C                                              COF 700
C ---HEAD VALUES ARE NOT INCLUDED IN COMPUTATION OF Q FACTOR SINCE COF 710
C LEAKAGE IS CONSIDERED IMPLICITLY---        COF 720
40 Q1=RATE(N)/(M(N)*DENOM)                   COF 730
TL(N)=Q1+2.*Q1*SUMN                          COF 740
PRATE=RATE(N)*M(N)                            COF 750
50 CONTINUE                                   COF 760
TMIN=TMIN*3.0                                 COF 770
TT=TT*3.0                                     COF 780
RETURN                                         COF 790
C ..... COF 800
C ..... COF 810
C ---COMPUTE TRANSMISSIVITY IN WT OR WT-ARTESIAN CONVERSION PROBLEM--- COF 820
C ***** COF 830
C ENTRY TRANS COF 840
C ***** COF 850
DO 60 I=1,DIML                                COF 860
DO 60 J=1,DIMW                                COF 870
N=I+DIML*(J-1)                                COF 880
IF (PERM(N).EQ.0.) GO TO 60                   COF 890
HED=PHI(N)                                    COF 900
IF (CONVRT.EQ.CHK(7)) HED=AMINI(SNGL(PHI(N)),TOP(N)) COF 910
T(N)=PERM(N)*(HED-BOTTOM(N))                 COF 920
IF (T(N).GT.0.) GO TO 60                     COF 930
IF (WELL(N).LT.0.) GO TO 70                  COF 940
C                                              COF 950
C ---THE FOLLOWING STATEMENTS APPLY WHEN NODES (EXCEPT WELL NODES) COF 960
C GO DRY--- COF 970
PERM(N)=0. COF 980
T(N)=0.0 COF 990
TR(N-DIML)=0. COF1000
TR(N)=0. COF1010
TC(N-1)=0. COF1020
TC(N)=0. COF1030
PHI(N)=SURI(N) COF1040
WRITE (P,150) I,J COF1050
60 CONTINUE COF1060
IF (KT.EQ.0) RETURN COF1070
GO TO 90 COF1080
C COF1090
C ---START PROGRAM TERMINATION WHEN A WELL GOES DRY--- COF1100
70 WRITE (P,120) I,J COF1110
WRITE (P,130) COF1120
IERR=1 COF1130
CALL ORDN COF1140
DO 80 I=2,INO1 COF1150
DO 80 J=2,JNO1 COF1160
N=I+DIML*(J-1) COF1170

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## Program listing—Continued

```

80 PHI(N)=KEEP(N)                                COF1180
SUM=SUM-DELT                                     COF1190
SUMP=SUMP-DELT                                   COF1200
KT=KT-1                                          COF1210
IF (KT.EQ.0) STOP                                COF1220
IF (IDK2.EQ.CHK(15)) CALL DISK                   COF1230
IF (PNCH.EQ.CHK(1)) CALL PUNCH                   COF1240
IF (MOD(KT,KTH).EQ.0) STOP                       COF1250
WRITE (P,140) KT,SUM                             COF1260
CALL DRDN                                        COF1270
IF (CHCK.EQ.CHK(5)) CALL CWRITE                  COF1280
STOP                                             COF1290
C                                               COF1300
C   ---COMPUTE T COEFFICIENTS---                  COF1310
C   *****                                       COF1320
C   ENTRY TCOF                                     COF1330
C   *****                                       COF1340
90 DO 110 I=1,IN01                               COF1350
DO 110 J=1,JN01                                   COF1360
N=I+DIML*(J-1)                                   COF1370
NR=N+DIML                                        COF1380
NB=N+1                                           COF1390
IF (T(N).EQ.0.) GO TO 110                        COF1400
IF (T(NR).EQ.0.) GO TO 100                       COF1410
TR(N)=(2.*T(NR)*T(N))/(T(N)*DELX(J)+T(NR)*DELX(J))*FACTX COF1420
100 IF (T(NB).EQ.0.) GO TO 110                   COF1430
TC(N)=(2.*T(NB)*T(N))/(T(N)*DELY(I)+T(NB)*DELY(I))*FACTY COF1440
110 CONTINUE                                     COF1450
RETURN                                           COF1460
C                                               COF1470
C   ---FORMATS---                                COF1480
C                                               COF1490
C   -----COF1500
C                                               COF1510
C                                               COF1520
120 FORMAT ('-*****WELL',I3,',',I3,' GOES DRY*****') COF1530
130 FORMAT ('1',50X,'DRAWDOWN WHEN WELL WENT DRY') COF1540
140 FORMAT ('1',32X,'DRAWDOWN FOR TIME STEP',I3,',',SIMULATION TIME =' COF1550
11PE15.7,' SECONDS') COF1560
150 FORMAT ('-',20(' '), ' NODE ',I4,',',I4,' GOES DRY ',20(' ')) COF1570
END                                              COF1580-

SUBROUTINE CHECKI(PHI,KEEP,PHE,STRT,T,TR,TC,S,QRE,WELL,TL,PERM,BOTCHK 10
ITOM,SY,RATE,RIVER,M,TOP,GRND,DELX,DELY) CHK 20
C -----CHK 30
C COMPUTE A MASS BALANCE CHK 40
C -----CHK 50
C CHK 60
C SPECIFICATIONS: CHK 70
REAL *8PHI,DBLE CHK 80
REAL *4KEEP,M CHK 90
INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CHK 100
1CONTR,LEAK,RECH,SIP,ADI CHK 110
C CHK 120
DIMENSION PHI(IZ,JZ), KEEP(IZ,JZ), PHE(IZ,JZ), STRT(IZ,JZ), T(IZ,JCHK 130
IZ), TR(IZ,JZ), TC(IZ,JZ), S(IZ,JZ), QRE(IZ,JZ), WELL(IZ,JZ), TL(IZCHK 140
2,JZ), PERM(IZ,JZ), BOTTOM(IP,JP), SY(IP,JP), RATE(IR,JR), RIVER(IRCHK 150
3,JR), M(IR,JR), TOP(IC,JC), GRND(IL,JL), DELX(JZ), DELY(IZ) CHK 160
C CHK 170
COMMON /SARRAY/ VF4(11),CHK(15) CHK 180

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## Program listing—Continued

```

COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,EROR,LECHK 190
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,QET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,CHK 200
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,CHK 210
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DICLK 220
4MW,JN01,IN01,R,P,PU,I,J,IDK1,IDK2 CHK 230
COMMON /CK/ ETFLXT,STORT,QRET,CHST,CHDT,FLUXT,PUMPT,CFLUXT,FLXNT CHK 240
COMMON /ARSize/ IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IH,IMAX,IMX1 CHK 250
RETURN CHK 260
C .....CHK 270
C *****CHK 280
ENTRY CHECK CHK 290
C *****CHK 300
C ---INITIALIZE VARIABLES--- CHK 310
PUMP=0. CHK 320
STOR=0. CHK 330
FLUXS=0.0 CHK 340
CHD1=0.0 CHK 350
CHD2=0.0 CHK 360
QREFLX=0. CHK 370
CFLUX=0. CHK 380
FLUX=0. CHK 390
ETFLUX=0. CHK 400
FLXN=0.0 CHK 410
C .....CHK 420
C .....CHK 430
C ---COMPUTE RATES,STORAGE AND PUMPAGE FOR THIS STEP--- CHK 440
DO 240 I=2,DIML CHK 450
DO 240 J=2,DIMW CHK 460
IF (T(I,J).EQ.0.) GO TO 240 CHK 470
AREA=DELX(J)*DELY(I) CHK 480
IF (S(I,J).GE.0.) GO TO 120 CHK 490
C .....CHK 500
C ---COMPUTE FLOW RATES TO AND FROM CONSTANT HEAD BOUNDARIES--- CHK 510
IF (S(I,J-1).LT.0..OR.T(I,J-1).EQ.0.) GO TO 30 CHK 520
X=(STRT(I,J)-PHI(I,J-1))*TR(I,J-1)*DELY(I) CHK 530
IF (X) 10,30,20 CHK 540
10 CHD1=CHD1+X CHK 550
GO TO 30 CHK 560
20 CHD2=CHD2+X CHK 570
30 IF (S(I,J+1).LT.0..OR.T(I,J+1).EQ.0.) GO TO 60 CHK 580
X=(STRT(I,J)-PHI(I,J+1))*TR(I,J)*DELY(I) CHK 590
IF (X) 40,60,50 CHK 600
40 CHD1=CHD1+X CHK 610
GO TO 60 CHK 620
50 CHD2=CHD2+X CHK 630
60 IF (S(I-1,J).LT.0..OR.T(I-1,J).EQ.0.) GO TO 90 CHK 640
X=(STRT(I,J)-PHI(I-1,J))*TC(I-1,J)*DELX(J) CHK 650
IF (X) 70,90,80 CHK 660
70 CHD1=CHD1+X CHK 670
GO TO 90 CHK 680
80 CHD2=CHD2+X CHK 690
90 IF (S(I+1,J).LT.0..OR.T(I+1,J).EQ.0.) GO TO 240 CHK 700
X=(STRT(I,J)-PHI(I+1,J))*TC(I,J)*DELX(J) CHK 710
IF (X) 100,240,110 CHK 720
100 CHD1=CHD1+X CHK 730
GO TO 240 CHK 740
110 CHD2=CHD2+X CHK 750
GO TO 240 CHK 760
C .....CHK 770
C ---RECHARGE AND WELLS--- CHK 780
120 QREFLX=QREFLX+QRE(I,J)*AREA CHK 790

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## Program listing—Continued

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      IF (WELL(I,J)) 130,150,140                                CHK 800
130  PUMP=PUMP+WELL(I,J)*AREA                                  CHK 810
      GO TO 150                                                CHK 820
140  CFLUX=CFLUX+WELL(I,J)*AREA                                CHK 830
150  IF (EVAP.NE.CHK(6)) GO TO 190                             CHK 840
C
C    ---COMPUTE ET RATE---                                     CHK 850
      IF (PHI(I,J).GE.GRND(I,J)-ETDIST) GO TO 160             CHK 860
      ETQ=0.0                                                  CHK 870
      GO TO 180                                                CHK 880
160  IF (PHI(I,J).LE.GRND(I,J)) GO TO 170                     CHK 890
      ETQ=GET                                                  CHK 900
      GO TO 180                                                CHK 910
170  ETQ=GET/ETDIST*(PHI(I,J)+ETDIST-GRND(I,J))              CHK 920
180  ETFLUX=ETFLUX-ETQ*AREA                                    CHK 930
C
C    ---COMPUTE VOLUME FROM STORAGE---                         CHK 940
190  STORE=S(I,J)                                             CHK 950
      IF (WATER.EQ.CHK(2)) STORE=SY(I,J)                       CHK 960
      IF (CONVRT.NE.CHK(7)) GO TO 230                          CHK 970
      X=KEEP(I,J)-PHI(I,J)                                     CHK 980
      IF (X) 200,210,210                                       CHK 990
200  HED1=PHI(I,J)                                            CHK1000
      HED2=KEEP(I,J)                                           CHK1010
      X=ABS(X)                                                  CHK1020
      GO TO 220                                                CHK1030
210  HED1=KEEP(I,J)                                           CHK1040
      HED2=PHI(I,J)                                           CHK1050
220  STORE=S(I,J)                                             CHK1060
      IF (HED1-TOP(I,J).LE.0.) STORE=SY(I,J)                   CHK1070
      IF ((HED1-TOP(I,J))*(HED2-TOP(I,J)).LT.0.0) STORE=(HED1-TOP(I,J))/CHK1080
      1X*S(I,J)+(TOP(I,J)-HED2)/X*SY(I,J)                       CHK1090
230  STOR=STOR+STORE*(KEEP(I,J)-PHI(I,J))*AREA                CHK1100
C
C    ---COMPUTE LEAKAGE RATE---                                CHK1110
      IF (LEAK.NE.CHK(9)) GO TO 240                             CHK1120
      IF (M(I,J).EQ.0.) GO TO 240                               CHK1130
      HED1=STRT(I,J)                                           CHK1140
      IF (CONVRT.EQ.CHK(7)) HED1=AMAX1(STRT(I,J),TOP(I,J))    CHK1150
      HED2=PHI(I,J)                                             CHK1160
      IF (CONVRT.EQ.CHK(7)) HED2=AMAX1(SNGL(PHI(I,J)),TOP(I,J))  CHK1170
      XX=RATE(I,J)*(RIVER(I,J)-HED1)*AREA/M(I,J)              CHK1180
      YY=TL(I,J)*(HED1-HED2)*AREA                               CHK1190
      FLUX=FLUX+XX                                             CHK1200
      XNET=XX+YY                                               CHK1210
      FLUXS=FLUXS+XNET                                         CHK1220
      IF (XNET.LT.0.) FLXN=FLXN-XNET                            CHK1230
240  CONTINUE                                                 CHK1240
C
C    .....                                                  CHK1250
C
C    ---COMPUTE CUMULATIVE VOLUMES, TOTALS, AND DIFFERENCES---
      STORT=STORT+STOR                                         CHK1260
      STOR=STOR/DELT                                           CHK1270
      ETFLXT=ETFLXT-ETFLUX*DELT                                CHK1280
      FLUXT=FLUXT+FLUXS*DELT                                    CHK1290
      FLXNT=FLXNT+FLXN*DELT                                    CHK1300
      FLXPT=FLXPT+FLXNT                                         CHK1310
      QRET=QRET+QREFLX*DELT                                     CHK1320
      CHDT=CHDT-CHD1*DELT                                       CHK1330
      CHST=CHST+CHD2*DELT                                       CHK1340
      PUMPT=PUMPT-PUMP*DELT                                     CHK1350
      CHK1360
      CHK1370
      CHK1380
      CHK1390
      CHK1400

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## Program listing—Continued

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CFLUXT=CFLUXT+CFLUX*DELT                                CHK1410
TOTL1=STORT+QRET+CFLUXT+CHST+FLXPT                      CHK1420
TOTL2=CHDT+PUMPT+ETFLXT+FLXNT                          CHK1430
SUMR=QREFLX+CFLUX+CHD2+CHD1+PUMP+ETFLUX+FLUXS+STOR     CHK1440
DIFF=TOTL2-TOTL1                                       CHK1450
PERCNT=0.0                                              CHK1460
IF (TOTL2.EQ.0.) GO TO 250                              CHK1470
PERCNT=DIFF/TOTL2*100.                                  CHK1480
250 RETURN                                              CHK1490
C .....                                                CHK1500
C .....                                                CHK1510
C ---PRINT RESULTS---                                  CHK1520
C *****                                              CHK1530
C ENTRY CWRITE                                         CHK1540
C *****                                              CHK1550
C .....                                                CHK1560
WRITE (P,260) STOR,QREFLX,STORT,CFLUX,QRET,PUMP,CFLUXT,ETFLUX,CHST,CHK1570
1,FLXPT,CHD2,TOTL1,CHD1,FLUX,FLUXS,ETFLXT,CHDT,SUMR,PUMPT,FLXNT,TOTL2,CHK1580
2L2,DIFF,PERCNT                                        CHK1590
RETURN                                                 CHK1600
C .....                                                CHK1610
C ---FORMATS---                                        CHK1620
C .....                                                CHK1630
C .....                                                CHK1640
C .....                                                CHK1650
C .....                                                CHK1660
260 FORMAT ('0',10X,'CUMULATIVE MASS BALANCE:',16X,'L**3',23X,'RATES FCHK1670
1OR THIS TIME STEP:',16X,'L**3/T'/11X,24(' '),43X,25(' ')//20X,'SOUCHK1680
2RCES:',69X,'STORAGE =',F20.4/20X,8(' '),68X,'RECHARGE =',F20.4/27XCHK1690
3,'STORAGE =',F20.2,35X,'CONSTANT FLUX =',F20.4/26X,'RECHARGE =',F2CHK1700
40.2,41X,'PUMPING =',F20.4/21X,'CONSTANT FLUX =',F20.2,30X,'EVAPOTRCHK1710
5ANSPIRATION =',F20.4/21X,'CONSTANT HEAD =',F20.2,34X,'CONSTANT HEACHK1720
6D:',27X,'LEAKAGE =',F20.2,46X,'IN =',F20.4/21X,'TOTAL SOURCES =',FCHK1730
720.2,45X,'OUT =',F20.4/96X,'LEAKAGE:',20X,'DISCHARGES:',45X,'FROM CHK1740
8PREVIOUS PUMPING PERIOD =',F20.4/20X,11(' '),68X,'TOTAL =',F20.4/1CHK1750
96X,'EVAPOTRANSPIRATION =',F20.2/21X,'CONSTANT HEAD =',F20.2,36X,'SCHK1760
SUM OF RATES =',F20.4/19X'QUANTITY PUMPED =',F20.2/27X,'LEAKAGE =',CHK1770
$F20.2/19X,'TOTAL DISCHARGE =',F20.2//17X,'DISCHARGE-SOURCES =',F20CHK1780
$.2/15X,'PER CENT DIFFERENCE =',F20.2//)              CHK1790
END                                                    CHK1800-

SUBROUTINE PRNTAI(PHI,SURI,T,S,WELL,DELX,DELY)          PRN 10
-----PRN 20
PRINT MAPS OF DRAWDOWN AND HYDRAULIC HEAD             PRN 30
-----PRN 40
C .....PRN 50
C SPECIFICATIONS:                                     PRN 60
REAL *8PHI,Z,XLABEL,YLABEL,TITLE,XN1,MESUR           PRN 70
REAL *4K                                              PRN 80
INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,PRN 90
1CONTR,LEAK,RECH,SIP,ADI                             PRN 100
C .....PRN 110
C DIMENSION PHI(IZ,JZ),SURI(IZ,JZ),S(IZ,JZ),WELL(IZ,JZ),DELX(JZ)PRN 120
1,DELY(IZ),T(IZ,JZ)                                  PRN 130
C .....PRN 140
C COMMON /SARRAY/ VF4(11),CHK(15)                    PRN 150
COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,EROR,LEPRN 160
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,PRN 170
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETGD,FACTX,FACTY,PRN 180
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DIPRN 190
4MW,JN01,INO1,R,P,PU,I,J,IDX1,IDX2                  PRN 200

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## Program listing—Continued

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COMMON /PR/ XLABEL(3),YLABEL(6),TITLE(5),XN1,MESUR,PRNT(122),BLANKPRN 210
1(60),DIGIT(122),VF1(6),VF2(6),VF3(7),XSCALE,DINCH,SYM(17),XN(100),PRN 220
2YN(13),NA(4),N1,N2,N3,YSCALE,FACT1,FACT2 PRN 230
COMMON /ARSize/ IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IH,IMAX,IMX1 PRN 240
RETURN PRN 250
C ..... PRN 260
C PRN 270
C ---INITIALIZE VARIABLES FOR PLOT--- PRN 280
C ***** PRN 290
C ENTRY MAP PRN 300
C ***** PRN 310
10 XSF=DINCH*XSCALE PRN 320
   YSF=DINCH*YSCALE PRN 330
   NYD=YDIM/YSF PRN 340
   IF (NYD*YSF.LE.YDIM-DELY(INO1)/2,) NYD=NYD+1 PRN 350
   IF (NYD.LE.12) GO TO 20 PRN 360
   DINCH=YDIM/(12.*YSCALE) PRN 370
   WRITE (P,310) DINCH PRN 380
   IF (YSCALE.LT.1.0) WRITE (P,320) PRN 390
   GO TO 10 PRN 400
20 NXD=WIDTH/XSF PRN 410
   IF (NXD*XSF.LE.WIDTH-DELX(JNO1)/2,) NXD=NXD+1 PRN 420
   N4=NXD*N1+1 PRN 430
   N5=NXD+1 PRN 440
   N6=NYD+1 PRN 450
   N8=N2*NYD+1 PRN 460
   NA(1)=N4/2-1 PRN 470
   NA(2)=N4/2 PRN 480
   NA(3)=N4/2+3 PRN 490
   NC=(N3-N8-10)/2 PRN 500
   ND=NC+N8 PRN 510
   NE=MAX0(N5,N6) PRN 520
   VF1(3)=DIGIT(ND) PRN 530
   VF2(3)=DIGIT(ND) PRN 540
   VF3(3)=DIGIT(NC) PRN 550
   XLABEL(3)=MESUR PRN 560
   YLABEL(6)=MESUR PRN 570
   DO 40 I=1,NE PRN 580
   NNX=N5-I PRN 590
   NNY=I-1 PRN 600
   IF (NNY.GE.N6) GO TO 30 PRN 610
   YN(I)=YSF*NNY/YSCALE PRN 620
30 IF (NNX.LT.0) GO TO 40 PRN 630
   XN(I)=XSF*NNX/YSCALE PRN 640
40 CONTINUE PRN 650
   RETURN PRN 660
C ..... PRN 670
C PRN 680
C ***** PRN 690
C ENTRY PRNTA(NG) PRN 700
C ***** PRN 710
C ---VARIABLES INITIALIZED EACH TIME A PLOT IS REQUESTED--- PRN 720
C DIST=WIDTH-DELX(JNO1)/2. PRN 730
   JJ=JNO1 PRN 740
   LL=1 PRN 750
   Z=NXD*XSF PRN 760
   IF (NG.EQ.1) WRITE (P,280) (TITLE(I),I=1,2) PRN 770
   IF (NG.EQ.2) WRITE (P,280) (TITLE(I),I=3,5) PRN 780
   DO 270 I=1,N4 PRN 790
C PRN 800
C ---LOCATE X AXES--- PRN 810

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## Program listing—Continued

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IF (I.EQ.1.OR.I.EQ.N4) GO TO 50 PRN 820
PRNT(1)=SYM(12) PRN 830
PRNT(N8)=SYM(12) PRN 840
IF ((I-1)/N1*N1.NE.I-1) GO TO 70 PRN 850
PRNT(1)=SYM(14) PRN 860
PRNT(N8)=SYM(14) PRN 870
GO TO 70 PRN 880
C PRN 890
C ---LOCATE Y AXES--- PRN 900
50 DO 60 J=1,N8 PRN 910
IF ((J-1)/N2*N2.EQ.J-1) PRNT(J)=SYM(14) PRN 920
60 IF ((J-1)/N2*N2.NE.J-1) PRNT(J)=SYM(13) PRN 930
C PRN 940
C ---COMPUTE LOCATION OF NODES AND DETERMINE APPROPRIATE SYMBOL--- PRN 950
70 IF (DIST.LT.0..OR.DIST.LT.Z-XN1*XSF) GO TO 220 PRN 960
YLEN=DELY(2)/2. PRN 970
DO 200 L=2,IN01 PRN 980
J=YLEN*N2/YSF+1.5 PRN 990
IF (T(L,JJ).EQ.0.) GO TO 140 PRN1000
IF (S(L,JJ).LT.0.) GO TO 190 PRN1010
INDX3=0 PRN1020
GO TO (80,90), NG PRN1030
80 K=(SURI(L,JJ)-PHI(L,JJ))*FACT1 PRN1040
C -TO CYCLE SYMBOLS FOR DRAWDOWN, REMOVE C FROM COL. 1 OF NEXT CARD-PRN1050
C K=AMOD(K,10.) PRN1060
GO TO 100 PRN1070
90 K=PHI(L,JJ)*FACT2 PRN1080
100 IF (K) 110,140,120 PRN1090
110 IF (J-2.GT.0) PRNT(J-2)=SYM(13) PRN1100
N=-K PRN1110
IF (N.LT.100) GO TO 130 PRN1120
GO TO 170 PRN1130
120 N=K PRN1140
IF (N.LT.100) GO TO 130 PRN1150
IF (N.GT.999) GO TO 170 PRN1160
INDX3=N/100 PRN1170
IF (J-2.GT.0) PRNT(J-2)=SYM(INDX3) PRN1180
N=N-INDX3*100 PRN1190
130 INDX1=MOD(N,10) PRN1200
IF (INDX1.EQ.0) INDX1=10 PRN1210
C -TO CYCLE SYMBOLS FOR DRAWDOWN, REMOVE C FROM COL. 1 OF NEXT CARD-PRN1220
C IF (NG.EQ.1) GO TO 150 PRN1230
INDX2=N/10 PRN1240
IF (INDX2.GT.0) GO TO 160 PRN1250
INDX2=10 PRN1260
IF (INDX3.EQ.0) INDX2=15 PRN1270
GO TO 160 PRN1280
140 INDX1=15 PRN1290
150 INDX2=15 PRN1300
160 IF (J-1.GT.0) PRNT(J-1)=SYM(INDX2) PRN1310
PRNT(J)=SYM(INDX1) PRN1320
GO TO 200 PRN1330
170 DO 180 II=1,3 PRN1340
JI=J-3+II PRN1350
180 IF (JI.GT.0) PRNT(JI)=SYM(11) PRN1360
190 IF (S(L,JJ).LT.0.) PRNT(J)=SYM(16) PRN1370
200 YLEN=YLEN+(DELY(L)+DELY(L+1))/2. PRN1380
210 DIST=DIST-(DELX(JJ)+DELX(JJ-1))/2. PRN1390
JJ=JJ-1 PRN1400
IF (JJ.EQ.0) GO TO 220 PRN1410
IF (DIST.GT.Z-XN1*XSF) GO TO 210 PRN1420

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## Program listing—Continued

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220 CONTINUE
C
C   ---PRINT AXES, LABELS, AND SYMBOLS---
C   IF (I-NA(LL).EQ.0) GO TO 240
C   IF ((I-1)/N1*N1-(I-1)) 250,230,250
230 WRITE (P,VF1) (BLANK(J),J=1,NC), (PRNT(J),J=1,N8), XN(1+(I-1)/6)
GO TO 260
240 WRITE (P,VF2) (BLANK(J),J=1,NC), (PRNT(J),J=1,N8), XLABEL(LL)
LL=LL+1
GO TO 260
250 WRITE (P,VF2) (BLANK(J),J=1,NC), (PRNT(J),J=1,N8)
C
C   ---COMPUTE NEW VALUE FOR Z AND INITIALIZE PRNT---
260 Z=Z-2.*XN1*XSF
DO 270 J=1,N8
270 PRNT(J)=SYM(15)
C
C   ---NUMBER AND LABEL Y AXIS AND PRINT LEGEND---
WRITE (P,VF3) (BLANK(J),J=1,NC), (YN(I),I=1,N6)
WRITE (P,300) (YLABEL(I),I=1,6)
IF (NG.EQ.1) WRITE (P,290) FACT1
IF (NG.EQ.2) WRITE (P,290) FACT2
RETURN
C
C   ---FORMATS---
C
C   -----
280 FORMAT ('1',53X,4A8//)
290 FORMAT ('0EXPLANATION'/' ' ,11(' - ')// ' R = CONSTANT HEAD BOUNDARY',/PRN1730
1' *** = VALUE EXCEEDED 3 FIGURES'/' MULTIPLICATION FACTOR =',F8.3)PRN1740
300 FORMAT ('0',39X,6A8)
310 FORMAT ('0',25X,10('**'),' TO FIT MAP WITHIN 12 INCHES, DINCH REVISPRN1760
1ED TO',G15.7,1X,10('**'))
320 FORMAT ('0',45X,'NOTE: GENERALLY SCALE SHOULD BE > OR = 1.0')
END
C
C
BLOCK DATA
-----
C
REAL *8XLABEL,YLABEL,TITLE,XN1,MESUR,RHO,B,D,F,H
INTEGER R,P,PU,DIML,DIMW,CHK,WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,
1CONTR,LEAK,RECH,SIP,ADI
C
COMMON /DPARAM/ RHO,B,D,F,H
COMMON /SARRAY/ VF4(11),CHK(15)
COMMON /SPARAM/ WATER,CONVRT,EVAP,CHCK,PNCH,NUM,HEAD,CONTR,EROR,LEBLD
1AK,RECH,SIP,U,SS,TT,TMIN,ETDIST,GET,ERR,TMAX,CDLT,HMAX,YDIM,WIDTH,BLD
2NUMS,LSOR,ADI,DELT,SUM,SUMP,SUBS,STORE,TEST,ETQB,ETQD,FACTX,FACTY,BLD
3IERR,KOUNT,IFINAL,NUMT,KT,KP,NPER,KTH,ITMAX,LENGTH,NWEL,NW,DIML,DI
4MW,JN01,IN01,R,P,PU,I,J,DK1,DK2
COMMON /PR/ XLABEL(3),YLABEL(6),TITLE(5),XN1,MESUR,PRNT(122),BLANKBLD
1(60),DIGIT(122),VF1(6),VF2(6),VF3(7),XSCALE,DINCH,SYM(17),XN(100),BLD
2YN(13),NA(4),N1,N2,N3,YSCALE,FACT1,FACT2
COMMON /ARSIZE/ IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IH,IMAX,IMX1
C
*****
C
DATA IZ,JZ,IP,JP,IR,JR,IC,JC,IL,JL,IS,JS,IMAX/13*20/,IH/1/
DATA CHK/'PUNC', 'WATE', 'CONT', 'NUME', 'CHEC', 'EVAP', 'CONV', 'HEAD', 'BLD

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## Program listing—Continued

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1LEAK','RECH','SIP ','LSOR','ADI','DK1 ','DK2 ' /,R,P,PU/5,6,7/,B,D;BLD 220
2F,H/4*0.D0/ BLD 230
  DATA SYM/'1','2','3','4','5','6','7','8','9','10','11','12','13','14','15','16','17','18','19','20','21','22','23','24','25','26','27','28','29','30','31','32','33','34','35','36','37','38','39','40','41','42','43','44','45','46','47','48','49','50','51','52','53','54','55','56','57','58','59','60','61','62','63','64','65','66','67','68','69','70','71','72','73','74','75','76','77','78','79','80','81','82','83','84','85','86','87','88','89','90','91','92','93','94','95','96','97','98','99','100','101','102','103','104','105','106','107','108','109','110','111','112','113','114','115','116','117','118','119','120','121','122' / BLD 240
1 ','R','W' / BLD 250
  DATA PRNT/122* ' /,N1,N2,N3,XN1/6,10,133,.833333333D-1/,BLANK/60* BLD 260
1 ' /,NA(4)/1000 / BLD 270
  DATA XLABEL/' X DIS- ',TANCE IN', ' MILES ' /,YLABEL/'DISTANCE', ' BLD 280
1FROM OR','IGIN IN ',Y DIRECT',ION, IN ',MILES ' /,TITLE/'PLOT BLD 290
2OF ',DRAWDCWN',PLOT OF ',HYDRAULI',C HEAD' / BLD 300
  DATA DIGIT/'1','2','3','4','5','6','7','8','9','10','11','12','13' BLD 310
1,'14','15','16','17','18','19','20','21','22','23','24','25','26' BLD 320
2,'27','28','29','30','31','32','33','34','35','36','37','38','39','40' BLD 330
340','41','42','43','44','45','46','47','48','49','50','51','52','53' BLD 340
43','54','55','56','57','58','59','60','61','62','63','64','65','66' BLD 350
51','67','68','69','70','71','72','73','74','75','76','77','78','79' BLD 360
61','80','81','82','83','84','85','86','87','88','89','90','91','92' BLD 370
7,'93','94','95','96','97','98','99','100','101','102','103','104' BLD 380
8,'105','106','107','108','109','110','111','112','113','114','115' BLD 390
9,'116','117','118','119','120','121','122' / BLD 400
  DATA VF1/'(1H ',',',', ' ',A1,F',10.2',')' / BLD 410
  DATA VF2/'(1H ',',',', ' ',A1,1',X,A8',')' / BLD 420
  DATA VF3/'(1H0',',',', ' ',A1,F',3.1',',12F1',0.2)' / BLD 430
  DATA VF4/'(1H0',',',', ' ',X,I2',',2X',',20F6',.1/(', ' ',X,2' BLD 440
10',F6.1',')' / BLD 450
C ***** BLD 460
  END BLD 470-

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