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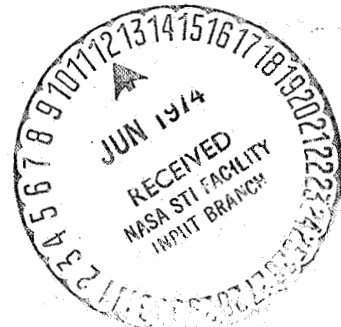
ANALYZER

Engineering-Program Manual

T. ISHIMOTO and L.C. FINK

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Systems Improved Numerical Differencing Analyzer
Engineering-Program Manual

T. Ishimoto and L. C. Fink

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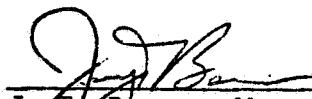
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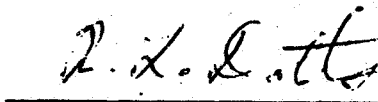
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PREFACE

The present SINDA computer program has evolved from the CINDA-3G program, which in turn evolved from the CINDA program, etc. With each major program revision an updated user manual was generated, but a more in-depth presentation of programming considerations and the theoretical development of the numerous subroutines were not generated. This SINDA program manual represents a preliminary effort to fill some of the existing void by describing the program structure, by identifying the major functions of each processor routine with a functional flow chart, and by a more in-depth mathematical description of the numerical solution subroutines. It is not the intent of this engineering-program manual, however, to provide sufficient detailed information for a user to make modifications and/or additions to the existing subprograms.

1. NOMENCLATURE AND MNEMONICS

1.1 Nomenclature

a_{ij}	= $k(A/\ell)_{ij}$, conduction coefficient between nodes i and j.
A	= array
A	= area
$(A/\ell)_{ij}$	= effective ratio of cross-sectional area to distance between nodes i and j.
b_{ij}	= radiation factor between nodes i and j (composed of radiation interchange factor and area)
C_i	= capacity of ith node
\bar{C}_i	= $C_i/\Delta t$, capacity of ith node divided by time-step
DD	= $1 - DN$ (allows certain fraction of "old" temperature to be included as part of temperature change for current time-step, refer to Section 6.2.5.1)
DN	\equiv DAMPA (user control constant, refer to Sections 6.2.5.1 and 6.2.3.2)
F, F1, F2	= multiplying factors, either user constants or literals, refer to Tables 6.2-1, 6.2-2 and 6.2-3).
G_{ij}	= $a_{ij} + \sigma b_{ij} (T_i^2 + T_j^2)(T_i + T_j)$
G_k	= a_{ij} or σb_{ij} , conduction or radiation coefficient.
k	= thermal conductivity
L	= a literal multiplying factor
N	= number of variable temperature nodes (NNA + NND)
NNA	= number of arithmetic-nodes
NND	= number of diffusion nodes
R	= resistance
p	= total number of nodes
q_i	= impressed heat load into the ith node
t	= time
Δt	= time-step
t_m	= $(TIME\phi + TIMEN)/2.0$, mean time
T	= temperature ($^{\circ}F$ or $^{\circ}R$)
T_m	= $(T_i + T_j)/2.0$, mean temperature ($^{\circ}F$ or $^{\circ}R$)

x, y = coordinate
 α = (k/C) , thermal diffusivity
 α = $\sum_{j=1}^P G_{ij}/C_i$, refer to equation 6.3-7
 β = factor that ranges from 0 to 1/2 (refer to equation 6.2-7)
 β' = 2β (used in subroutine CNVARB)
 \mathcal{F}_{ij} = radiation interchange factor including inter-reflections between nodes i and j .
 σ = Stefan-Boltzmann constant ($.1714 \times 10^{-8}$ Btu/hr $^{\circ}R^4 ft^2$)
 τ_n = $\frac{\Delta t_n}{\Delta t_{n-1} + \Delta t_n}$, weighting factor (equation 6.3-13)

$(A^i:t_m)$ Interpolated value of array A using t_m as the independent variable
 $(A^i:T_i)$ " " " " " " T_i " " " "
 $(A^i:T_j)$ " " " " " " T_j " " " "
 $(A^i:T_m)$ " " " " " " T_m " " " "
 $(A^P:T_i)$ Polynomial " " " " " " T_i " " " "
 $(A^P:T_j)$ " " " " " " T_j " " " "
 $(A^P:T_m)$ " " " " " " T_m " " " "
 $(A^b:T_i, t_m)$ Interpolated value of the bivariate array A using T_i and t_m as independent variables.
 $(A^b:T_m, t_m)$ Interpolated value of the bivariate array A using T_m and t_m as independent variables.

Subscripts

i = i th node
 j = j th node
 ij = between nodes i and j
 i, n = updating of i th temperature, source, etc. at time-step n .
 i, k = updating of i th temperature, etc. at k th iteration.
 ij, n = updating of coefficient between nodes i and j at time-step n
 ij, k = updating of coefficient between nodes i and j at k th iteration
 m = mean

1.2 Mnemonics

1.2.1 Control constants (refer to Sections 6.2.3.1 and 6.2.3.2)

ARLXCA = allowable arithmetic node relaxation temperature change
ARLXCC = calculated maximum arithmetic node relaxation temperature change
ATMPCA = allowable arithmetic node temperature change
ATMPCC = calculated maximum arithmetic node temperature change
BACKUP = back switch
BALENG = specified system energy balance
CSGFAC = time-step factor
CSGMAX = maximum value of $C_i/\Sigma G_{ij}$
CSGMIN = minimum value of $C_i/\Sigma G_{ij}$
CSGRAL = allowable range between CSGMIN and CSGMAX
DAMPA = arithmetic node damping factor
DAMPD = diffusion node damping factor
DRLXCA = allowable diffusion node relaxation temperature change
DRLXCC = calculated diffusion node relaxation temperature change
DTIMEH = allowable maximum time-step
DTIMEI = specified time-step for implicit solutions
DTIMEL = allowable minimum time-step
DTIMEU = contains computed time-step
DTMPCA = allowable diffusion node temperature change
DTMPCC = calculated maximum diffusion node temperature change
ENGBAL = calculated system energy balance
ITEST = contains dummy integer constant
JTEST = contains dummy integer constant
KTEST = contains dummy integer constant
LAXFAC = number of iterations for linearized lumped parameter system,
CINDSM only.
LINECT = line counter location for program output
LØØPCT = contains number of iterations performed
NØCØPY = contains no copy switch for matrix users
NLØØP = number of specified iteration loops
ØPEITR = output each iteration switch
PAGECT = page counter location for program output

RTEST = contains dummy floating point constants
STEST = contains dummy floating point constants
TIMEM = $(\text{TIME}\emptyset + \text{TIMEN})/2.0$, mean time for computational interval
TIMEN = TIMEN + DTIMEU, new time at the end of computational interval
TIMEND = problem stop time
TIME \emptyset = old time at the start of the computational interval
TTEST = contains dummy floating point constants
UTEST = contains dummy floating point constants
VTEST = contains dummy floating point constants

1.2.2 Numerical Solution Routines (refer to Sections 6.3 - 6.5)

CINDSS = steady state routine, refer to Section 6.5.1
CINDSL = steady state routine, refer to Section 6.5.2
CINDSM = steady state routine, refer to Section 6.5.3
CNBACK = implicit routine, refer to Section 6.4.1
CNDUFR = explicit routine, refer to Section 6.3.4
CNEXPN = explicit routine, refer to Section 6.3.3
CNFAST = explicit routine, refer to Section 6.3.2
CNFRDL = explicit routine, refer to Section 6.3.1
CNFWBK = implicit routine, refer to Section 6.4.2
CNFWRD = explicit routine, refer to Section 6.3.1
CNQUIK = explicit routine, refer to Section 6.3.5
CNVARB = implicit routine, refer to Section 6.4.3

1.2.3 Options (used in Tables 6.2-1 - 6.2-3)

BIV = Bivariate Interpolation Variable
DIT = Double Interpolation with Time as variable
DIV = Double Interpolation Variable
DPV = Double Polynomial Variable
DTV = Double interpolation with Time and Temperature as Variables
SIT = Single Interpolation with Time as variable
SIV = Single Interpolation Variable
SPV = Single Polynomial Variable

1.2.4 Routines and Subroutines of Preprocessor

- SINDA = routine that specifies overlay of preprocessor to system allocator.
- PREPRØ = main routine for preprocessor; initializes counters and FØRTRAN logical units; sets length of dynamic storage array and controls major logic.
- ALPINT = subroutine that accepts an integer in alphanumeric format and converts it to integer format; determines relative number of this actual number and converts it back to alphanumeric format.
- BLKCRD = subroutine that formats the five generated FØRTRAN routines (SINDA, EXECTN, VARBL1, VARBL2, and ØUTCAL) in 507 word blocks.
- CØDERD = Subroutine that reads and checks the block header cards for the data blocks.
- CØNVRT = subroutine that converts Hollerith data to integer data.
- DATARD = subroutine that scans the data block card images under an A format and determines appropriate format to reread the card images.
- ERRMES = subroutine that prints most of the error messages generated within the data blocks.
- FINDRM = subroutine that moves the data in the dynamic storage array either up or down by 100 words.
- GENLNK = subroutine that generates the driver (FØRTRAN routine named SINDA) for the user's program.
- GENUK = subroutine that generates user constants.
- INCØRE = subroutine that reads data into the dynamic storage array for the parameter-runs option.
- MXTØFN = subroutine that processes data for the "m" option (converts card images from mixed FØRTRAN/SINDA notation to FØRTRAN notation.
- NØDEDA = subroutine that processes data for node and conductor data blocks.
- PCS2 = subroutine that packs the FØRTRAN addresses for the array and constants locations required by the second pseudo-compute sequence.
- PRESUB = subroutine that reads and checks the block header cards for the operations blocks and generates the non-executable FØRTRAN cards for each of the operations blocks via a call to BLKCRD.

PSEUDØ = subroutine that forms the first and second pseudo-compute sequences.

QDATA = subroutine that checks and processes all data input in the source data block.

RELACT = subroutine that finds the relative node numbers from the actual node number; computes the FORTRAN address for arrays and user constants from the actual number.

SEARCH = subroutine that retains a relative number for nodes, conductors, user constants, and arrays, given the actual number.

SETFMT = subroutine that processes the card for the "new format" option; that is, it sets up the format for data cards as specified by the cards with an "N" in column one.

SINDA4 = subroutine that reads and processes the user input cards from the operations blocks.

SKIP = subroutine that is used when a problem is RECALLED; it positions the tape to the proper problem as specified on the first card of the data deck.

SPLIT = subroutine that reads the data from the RECALL tape and splits the RECALL information onto the proper data "tape" and the dictionary "tape."

SQUEEZ = subroutine that compresses the specified data groups in the dynamic storage array.

STFFB = subroutine that fills out a card image in array KBLK with Hollerith blanks.

TYPCHK = subroutine that checks the input from the data blocks for the correct type (integer, floating point, or alphanumeric).

WRDTA = subroutine that writes the program data "tape" in the format required by INPUTT or INPUTG.

WRTPMT = subroutine that writes the required data for parameter runs on the parameter "runs" "tape" and the dictionary "tape."

WRTBLK = subroutine that writes the 507 word blocks contained in array KBLK on the program FORTRAN "tape."

1.2.5 Others

SINDA = Systems Improved Numerical Differencing Analyzer

LPCS = Long Pseudo-Compute Sequence

SPCS = Short Pseudo-Compute Sequence

LPCS2 = Second Long Pseudo-Compute Sequence

PCS1 = Pseudo-Compute Sequence One

PCS2 = Pseudo-Compute Sequence Two

TSUM = elapsed time from last printout

TPRINT = time of last printout.

2. BACKGROUND ON SINDA

The original CINDA^{*1} (Chrysler Improved Numerical Differencing Analyzer) computer program was developed by the Thermodynamics Section of the Aerospace Physics Branch of Chrysler Corporation Space Division at NASA Michoud Assembly Facility and was coded in FORTRAN-II and FAP for the IBM-7094 computers. CINDA was the product of an intensive analytical, engineering and programming effort that surveyed numerous thermal analyzer-type programs and studied several in-depth. The foundation for CINDA was the storage and addressing of information required only for the network solution and the systems features which allowed the re-utilization of core storage area and brought into core only those instructions necessary for the solution of a particular problem. A systems compiler computer program that automatically optimized the utilization of computer core space was developed. This meant the generation of an integrated operation of relative addressing, packing features, peripheral tape storage units and overlay features.

CINDA evolved into CINDA-3G² which was developed by the same group that generated CINDA with a major portion of the work done under contract NASA/MSC NAS9-7043. CINDA-3G represented (essentially) a complete rework of CINDA in order to take advantage of the improved systems software and machine speeds of the 3rd generation computers. CINDA was unsuitable for standard operation on third generation computers since it was virtually a self-contained program having its own Update, Monitor and Compiler. On the other hand, CINDA-3G consisted of a preprocessor (written in FORTRAN) which accepted the user input data and the block data input. The user input data was converted into advanced FORTRAN language subroutines and block data input was passed onto the system FORTRAN Compiler. This required a double pass on data where previously only one was required, but the increased speed and improved software of the third generation machines more than compensated for the double pass.

SINDA³ (Systems Improved Numerical Differencing Analyzer) was developed by the Heat Transfer and Thermodynamics Department of TRW Systems Group, Redondo Beach. Most of the improvements and subroutine additions to CINDA-3G was done as part of the NASA/MSC contract NAS 9-8389,

* Superscript numbers refer to the literature cited in the Reference Section.

entitled, "Development of Digital Computer program for Thermal Network Correction." Programming and systems integration were directed to the UNIVAC-1108 computer.

SINDA relied quite heavily on CINDA-3G and data deck compatibility was rigorously followed; as a result, CINDA-3G data decks should, in the main, be directly operational on the SINDA program although some differences exist. For example, properties are updated before VARIABLES 1 call in CINDA-3G, whereas the properties are updated within the numerical solution routines after VARIABLES 1 call in SINDA. The primary differences between SINDA and CINDA-3G are: (1) elimination wherever possible of assembly language coding; (2) increased mnemonic options to aid the program user in data input; (3) inclusion of a second pseudo computer sequence for evaluation of nonlinear network elements; and (4) additional subroutines such as STEP (sensitivity analysis) and KALØBS-KALFIL (Kalman filtering). Most of the changes and additions to CINDA-3G were required in order to integrate the thermal network correction subroutine package into the existing CINDA-3G program.

During the development of SINDA a number of useful improvements became apparent. As a result, modifications to SINDA were made a part of the NASA/MSC contract NASA 9-10435 entitled, "Development of an Advanced SINDA Thermal Analyzer System." These changes that include a variable input format, simplified parameter runs, and generated user constants were made by the same group that developed SINDA. These improvements are reported in an updated SINDA users manual.⁴

3. GENERAL SINDA PROGRAM DESCRIPTION

3.1 SINDA Operating System

SINDA is more like an operating system rather than applications program. SINDA is programmed as a preprocessor in order to accommodate the desired operations relative to overlay features, data packing, dynamic storage allocation and subroutine library file, but yet be written in FØRTRAN. This preprocessor operates in an integral fashion with a library of numerous and varied subroutines,^{3, 4} which may be called in any desired sequence but yet operate in an integrated manner. The preprocessor reads the input data, assigns relative numbers, packs this information, forms a pseudo-compute sequence(s) (which will be described briefly in a later paragraph of this section and is described in more detail in Section 4, called Preprocessor), and writes the operations blocks on a peripheral unit as FØRTRAN source language with all of the data values dimensioned exactly in labeled common. In turn, controls are shifted to the system FØRTRAN compiler which compiles the constructed subroutines and enters execution. The FØRTRAN allocator has access to the SINDA subroutine library and loads only those subroutines called by the problem being processed.

As a result of this type of systems operation SINDA is extremely dependent upon the systems software. However, once the program is operational on a particular computer, the user-prepared problem data deck can be confined to the control cards and deck set-up requirements at a particular installation.

It should be recognized that the use of a preprocessor provides a computer with a large capability and considerable flexibility, but because of the numerous options that are generally offered, user instructions are more difficult than other thermal analyzer-type programs which have less flexibility.

3.2 Use of Lumped-Parameter Concept

Use of SINDA is based on a lumped parameter representation of a physical system. This means that SINDA does not solve a set of partial differential equations that represents a distributive system, but rather SINDA numerically solves a set of ordinary (and in general) nonlinear

differential equations that represent a lumped parameter system. The procedure for the formulation and the numerical solutions of the lumped parameter equations are reported extensively in literature and basic considerations are presented in Section 5. For the discussion to follow on the pseudo compute sequence it is convenient to indicate a general set of ordinary linear differential heat balance equations,

$$\frac{dT_i}{dt} = \frac{1}{C_i} \left[q_i + \sum_{j=1}^P a_{ij} (T_j - T_i) \right] \quad (3.2-1)$$

$i = 1, 2, \dots, N$ (number of variable temperatures)

$T_j = \text{constant}, N < j \leq p$

where, C_i = the i th nodal capacity

q_i = the heat load into node i (impressed)

a_{ij} = the conduction coefficient between nodes i and j [$= k \left(\frac{A}{l} \right)_{ij}$]

t = time

Suppose an implicit numerical method as discussed in Section 5.2.2 of this manual is chosen; the implicit finite difference form becomes after letting,

$$dT_i/dt \approx (T_{i,n+1} - T_{i,n})/\Delta t, \quad T_j = T_{j,n+1} \quad \text{and} \quad T_i = T_{i,n+1},$$

$$\frac{C_i}{\Delta t} (T_{i,n+1} - T_{i,n}) = q_i + \sum_{j=1}^P a_{ij} (T_{j,n+1} - T_{i,n+1}) \quad (3.2-2)$$

where, $T_{i,n}$ = temperature at time point t_n

$T_{i,n+1}$ = temperature at time point $t_{n+1} = t_n + \Delta t$

Δt = time-step

Rearrangement of equation (3.2-2) yields,

$$\left(\bar{C}_i + \sum_{\substack{j=1 \\ j \neq i}}^P a_{ij} \right) T_{i,n+1} - \sum_{\substack{j=1 \\ j \neq i}}^N a_{ij} T_{j,n+1} = q_i + \bar{C}_i T_{i,n} + \sum_{j=N+1}^P a_{ij} T_{j,n} \quad (3.2-3)$$

$i = 1, 2, \dots, N$

$T_{j,n} = \text{constant}, N < j \leq p$

where, $\bar{C}_i = C_i/\Delta t$, average capacity of node i over Δt time-step

3.3 Pseudo-Compute Sequence (PCS)

A pseudo-compute sequence as generated by the SINDA preprocessor

is a list of numbers that indicates the position of required data values in various arrays such as conductance, temperature and capacitance. This meaning will become clearer by formulating equation (3.2-3) in a matrix form. The matrix formulation is straightforward since temperatures at time-step $n+1$ are the unknowns and terms on the right side of equation (3.2-3) represent the forcing function. Let us expand equation (3.2-3) to show this,

$$\begin{aligned}
 (\bar{C}_1 + \sum_{j=1}^P a_{1j}) T_{1,n+1} - a_{12} T_{2,n+1}, \dots, -a_{1N} T_{N,n+1} &= q_1 + \bar{C}_1 T_{1,n} + \sum_{j=N+1}^P a_{1j} T_{j,n} \\
 -a_{21} T_{1,n+1} + (\bar{C}_2 + \sum_{j=1}^P a_{2j}) T_{2,n+1}, \dots, -a_{2N} T_{N,n+1} &= q_2 + \bar{C}_2 T_{2,n} + \sum_{j=N+1}^P a_{2j} T_{j,n} \\
 &\vdots \\
 -a_{N1} T_{1,n+1} - a_{N2} T_{2,n+1}, \dots, (\bar{C}_N + \sum_{j=1}^P a_{Nj}) T_{N,n+1} &= q_N + \bar{C}_N T_{N,n} + \sum_{j=N+1}^P a_{Nj} T_{j,n}
 \end{aligned}$$

Thus the matrix form of equation (3.2-3) becomes,

$$[\beta] \{T'\} = \{Q\} \quad (3.3-1)$$

where,

$$\beta = \begin{bmatrix} \sum_{j=2}^P (\bar{C}_1 + a_{1j}), & -a_{12} & \dots, & -a_{1N} \\ -a_{21} & , \sum_{j=1}^P (\bar{C}_2 + a_{2j}), & \dots, & -a_{2N} \\ \vdots & \sum_{j \neq 2}^P & \vdots & \vdots \\ -a_{N1} & , & -a_{N2} & \dots, \sum_{j=1}^P (\bar{C}_N + a_{Nj}) \end{bmatrix} \quad (3.3-2)$$

$$T' = \begin{pmatrix} T_{1,n+1} \\ T_{2,n+1} \\ \vdots \\ T_{N,n+1} \end{pmatrix} \quad (3.3-3) \quad ; \quad \{Q\} = \begin{pmatrix} q_1 + \bar{C}_1 T_{1,n} + \sum_{j=N+1}^P a_{1j} T_{j,n} \\ q_2 + \bar{C}_2 T_{2,n} + \sum_{j=N+1}^P a_{2j} T_{j,n} \\ \vdots \\ q_N + \bar{C}_N T_{N,n} + \sum_{j=N+1}^P a_{Nj} T_{j,n} \end{pmatrix} \quad (3.3-4)$$

The matrix represented by equation (3.3-2) appears to be a full matrix (very small number of elements that are zero), but in reality most

of the off-diagonal elements are zero. Thus, if equation (3.2-3) was to be solved by a matrix inversion technique, all elements including zeros must be stored. Since the number of elements varies as N^2 (N is the number of nodes), the required number of data locations would vary as N^2 and the computer time required for matrix inversion would be proportional to N^3 .

The explicit and iterative implicit numerical methods (refer to Section 5) of solving equation (3.2-1) lend themselves for optimizing the data storage area required and for reducing the solution time. If the conductors are numbered and related to the appropriate adjoining nodes as indicated in Table (3.2-1), retention of adjoining node number for each conductor provides a means of identifying element position in the coefficient matrix. This can be seen by considering the one-dimensional heat conduction example pictured in Figure (3.3-1).

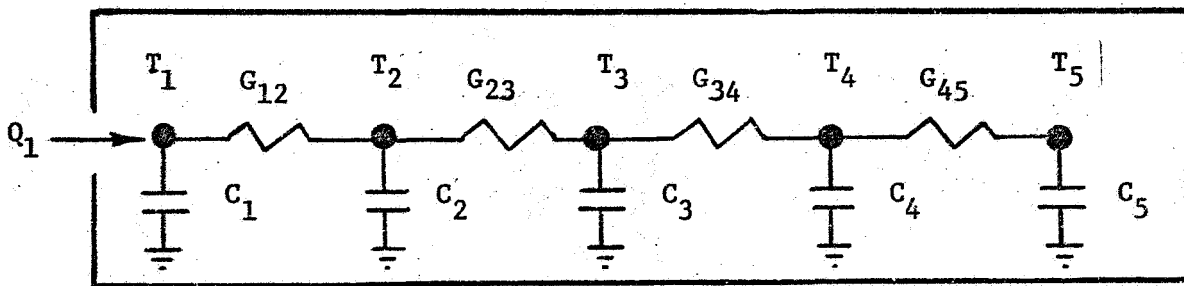


Figure 3.3-1. Thermal Circuit for a One-Dimensional System

The set of equations associated with the problem of Figure (3.3-1) may be readily expressed as,

$$\begin{bmatrix} (\bar{C}_1 + G_1) & -G_1 & 0 & 0 & 0 \\ -G_1 & (\bar{C}_2 + G_1 + G_2) & -G_2 & 0 & 0 \\ 0 & -G_2 & (\bar{C}_3 + G_2 + G_3) & -G_3 & 0 \\ 0 & 0 & -G_3 & (\bar{C}_4 + G_3 + G_4) & -G_4 \\ 0 & 0 & 0 & -G_4 & (\bar{C}_5 + G_4) \end{bmatrix} \begin{pmatrix} T_1 \\ T_2 \\ T_3 \\ T_4 \\ T_5 \end{pmatrix} = \begin{pmatrix} Q_1 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} \quad (3.3-5)$$

By comparing the element position of equation (3.3-5) with the tabular identification in Table (3.2-1) it is seen that elements with zero values need not be stored. The main diagonal term is never zero and is a composite of capacitance and off-diagonal conductors.

Table (3.2-1) Tabular Identification of Conductor and Adjoining Node Numbers

Conductance Number	ith Node	Adjoining Node Number	Comment
G#	N#	N#	
1	1	2	G1 is conductor #1 between nodes 1 and 2.
1	2	1	G1 is conductor #1 between nodes 2 and 1.
2	2	3	G2 is conductor #2 between nodes 2 and 3.
.	.	.	
.	.	.	
4	5	4	G4 is conductor #4 between nodes 5 and 4.

It is of interest to note that the use of a pseudo-compute sequence is only one of a number of ways to store data efficiently. For example, TRW TAP⁵ does not employ a pseudo-compute sequence because of other user requirements. However, from a data storage standpoint, it appears that the use of a pseudo-compute sequence utilizes computer core most efficiently.

More than one pseudo-compute sequence is formed by SINDA. Both a so-called long (LPCS) and a so-called short (SPCS) pseudo-compute sequence as used in CINDA-3G² are formed and in addition a second long pseudo-compute (LPCS2) required for thermal network correction is also formed in SINDA. A detailed discussion of these pseudo-compute sequences will be presented in Section 4.6, but is of interest here to indicate the characteristics of these "sequences."

3.3.1 Long Pseudo-Compute Sequence (LPCS)

A long pseudo-compute sequence identifies the position and value of all off-diagonal elements of the coefficient matrix. This is done by operating on adjoining node numbers which have been assigned relative node numbers by the preprocessor. Since nodal temperatures are calculated sequentially in ascending numerical order, the conductor and adjoining node number are searched until node one is found with the conductor number and the other adjoining node number stored in a single core location. In addition, several indicators are stored in this single core location. These

indicators are : (1) var C (indicates the input of a capacitor as a variable); (2) var G (indicates the input of a conductor as a variable); (3) rad (indicates the input of a radiation conductance); (4) Q (indicates the input of a source in the source data block); (5) one-way (indicates the input of a one-way conductor); and (6) last G (indicates the last conductor to a particular node). Order of indicator storage is indicated in Table (3.3-2).

Search is continued until all node-one's have been located and characteristics processed. The procedure is repeated for all node-two's and so forth sequentially until all nodes have been processed. The important consideration of a LPCS is the encounter of each conductor of the coefficient matrix twice. Formation of a pseudo-compute sequence for the example shown in Table (3.3-1) is given in Table (3.3-2). A pseudo-compute sequence starts with node one and advances the node number by one each time a last conductor indicator (last G) is passed. The conductor and node numbers identify the position of the conductor value in an array of conductor values and the position of the temperature, capacitor and source values in arrays of temperature, capacitor and source values respectively.

A long pseudo-compute sequence is well-suited for "successive point" iteration (refer to Section 5.2.2 for a discussion of this) of the implicit finite difference equations because all elements of the coefficient matrix are identified. Thus, when a row of the coefficient matrix is processed and a new value of temperature obtained, the new temperature can then be used in the calculation procedure of succeeding rows.

3.3.2 Short Pseudo-Compute Sequence (SPCS)

The short pseudo-compute sequence identifies each conductor only once and since the coefficient matrix (equation 3.3-1) is symmetrical, all sparsity and off-diagonal elements of the coefficient matrix are accounted for. The node being processed and the adjoining node number reveal temperature- and source-value locations. The short pseudo-compute sequence for the example in Table (3.3-1) is formed in Table (3.3-3). By placing a minus sign on the initially encountered other-adjoining nodes, these nodes are not recognized on a second encounter. A short pseudo-compute sequence

Table (3.3-1) Example of Conductor Connections

Conductor No.	Adjoining Node Numbers	
	G#	N#
1	1	2
2	1	3
3	1	4
4	2	3
5	2	4
6	3	4

Table (3.3-2) Long Pseudo-Compute Sequence (LPCS)
for the Example of Table (3.3-1)

Node No. Searched	Last G	var C	var G	rad	Q	G#	One- way	Node # Stored
1						1		2
1						2		3
1	1					3		4
2						1		1
2						4		3
2	1					5		4
3						2		1
3						4		2
3	1					6		4
4						5		2
4	1					6		3

Table (3.3-3) Short Pseudo-Compute Sequence (SPCS)
for the Example of Table (3.3-1)

Node No. Searched	Last G	var G	var G	rad	Q	G#	One- way	Node # Stored
1						1		2
1						2		3
1	1					3		4
2						4		3
2	1					5		4
3	1					6		4
4	1					0		0

is well-suited for explicit numerical solutions methods which calculate the energy flow through the conductor, add it to the source location of the node being processed and subtract it from the source location for the adjoining node. The SPCS can be used for implicit methods of solution but the "block" iterative procedure (refer to Section 5.2.2 for a discussion of this) must be used since succeeding rows of conductor and adjoining node numbers do not contain the necessary element information.

3.3.3 Second Long Pseudo-Compute Sequence (LPCS2)

The second long pseudo-compute sequence (LPCS2) as a user input option flags a non-linear conductor between two diffusion nodes twice; LPCS flags the non-linear conductor only one. LPCS2 is required for the thermal network correction of a sparse network by the use of subroutine KAFIL (refer to Reference 3 or 6).

3.3.4 Pseudo-Compute Sequence One (PCS 1) and Pseudo-Compute Sequence Two (PCS 2)

PCS 1 and PCS 2 are not user options but are fixed internally. The contents of PCS 1 and PCS 2 are governed by the user input of LPCS, SPCS or LPCS2). PCS 1 contains two relative addresses (conductor and adjoining node locations), two non-linear type indicators, and an impressed source indicator. Indicators are keyed through a simple counter to a second pseudo-compute sequence (PCS 2) which contains integer addresses or relative constant and array starting locations necessary for evaluation of temperature varying coefficients and time varying coefficients for sources. When the input data contain literal values in SIV type calls, the preprocessor stores the values as extended user constants and supplies the relative constant address to the second pseudo-compute sequence. Detailed discussion on PCS 1 and PCS 2 is presented in Section 4.6.

3.4 Data Logistics

3.4.1 Relative Numbers

Both the long and short pseudo-compute sequences require the storage of only the finite values in the coefficient matrix, thereby taking advantage of matrix sparsity. If the short pseudo-compute sequence is used, the advantage of symmetry is accounted for. Conductors with the same constant value may share the same conductor number and value. The storage efficiency of the pseudo-compute sequences requires the sequential numbering of the nodes and the conductors. Since the numbering of thermal math-models is arbitrary and not sequential, the SINDA program assigns relative numbers (starting from one, sequential and ascending) to the actual numbers of the incoming node data, conductor data, constants data and array data in the order received. Thus, numbers not used in the actual numbering system are neither identified nor required.

3.4.2 Storage Requirements and Dynamic Storage Allocation

All numerical solution subroutines require three locations for each diffusion node data (temperature, capacitance and source), two locations for each arithmetic node data (temperature and source), one location for each boundary data (temperature) and one location for each conductor value. In addition intermediate data storage ranging from zero to three locations per node may be required for the storage of temperatures and temperature differences; acceleration of convergence (refer to Section 6.2.7) used in the implicit and steady state routines (except CINDSS) requires three locations. Storage requirements for conductances depends upon the problem. For example, each internal diffusion and arithmetic node of a three-dimensional conduction system with rectangular-nodalization will be connected with only three being unique; thus, each diffusion node (or arithmetic node) in a three-dimensional conduction system requires from six to nine storage locations for data values (temperature, capacitance, source, three conductors and up to three intermediate locations). Now each of the conductors for the short pseudo-compute sequence requires a single core location that contains two integer values (conductor and adjoining node numbers) and six indicators (refer to Section 3.3.1 for description). Each of the conductors between variable temperatures for the

long pseudo-compute sequence requires two core locations since the conductors are used twice during the computational process. This means that each internal node of a three-dimensional conduction system will require six data addressing locations for the long pseudo-compute sequence and, on the average, three data addressing locations for the short pseudo-compute sequence.

Thus for a three-dimensional conduction system (no radiation), the number of required core locations per node can vary from nine (temperature, capacitance, source, three unique conductors and three data addressing locations) to fifteen (temperature, capacitance, source, six conductors and six data addressing locations) exclusive of the second pseudo-compute sequence which is required for variable coefficients, capacitance and sources.

The user must allocate an array of data locations which is to be used for intermediate data storage and initialize the array start and length indicators. Each subroutine that requires intermediate storage area has access to this array and the start and length indicators. During a subroutine execution a check on the sufficiency of space is made and start and length indicator are updated. If a subroutine calls upon another subroutine that requires intermediate storage, the called subroutine repeats the check and update procedure. Whenever any subroutine terminates its operation, the start and length indicators are returned to their entry values. This process is termed "Dynamic Storage Allocation" and allows subroutines to share a common working area.

3.5 Order of Computation

A network data deck consists of four data blocks (node, conductor, constants, and array), one optional data block (source) and four operations blocks which are preprocessed by the preprocessor and passed on to the system FORTRAN compiler. Non-network problems require no node or conductor data blocks. The operations blocks are named EXECUTION, VARIABLES 1, VARIABLES 2 and OUTPUT CALLS: the SINDA preprocessor constructs these blocks into individual subroutines with the entry names EXECUTN, VARBL1, VARBL2 and OUTCAL, respectively. After a successful FORTRAN compilation, control is passed to the EXECUTN subroutine. This means that the order of computation depends on the sequence of subroutine calls placed in the EXECUTION block

by the program user. No other operations blocks are performed unless called upon by the user either directly by name or indirectly through a subroutine call. The numerical solution subroutines described in Section 6 internally call upon VARBL1, VARBL2 and ØUTCAL; The internal order of computation for these routines is similar with the primary difference being the numerical solution method. A general flow diagram of the numerical solution routines, as well as a detailed description of each is presented in Section 6.

4. PREPROCESSOR

4.1 General Description

The SINDA preprocessor reads and analyzes the user input deck and from this information constructs a program tailored to the user's requirements.

The rationale for a preprocessor is flexibility and speed. Flexibility is achieved by providing the user with a library of routines to solve problems, manipulate data, and print selected values. In addition, the user may insert non-SINDA routines into the constructed program. Speed (defined here as minimal execution time) is achieved by structuring the data in an efficient manner.

The SINDA preprocessor consists of thirty routines with seven overlay links. All of the routines are written in FØRTRAN except for one assembly language routine which writes a "tape" in a format acceptable to the FØRTRAN compiler. These routines provide the user with a number of major options in the type of problem to be solved and the form of the data to be used. Henceforth these major options are designated as "major logic" of the preprocessor. See Figure 4.1-1 for a flow chart of the major logic of the preprocessor and its interface with the user program.

The major logic consists of the five following options: (1) NASA MSC EDIT feature; (2) RECALL option; (3) generation of a THERMAL problem; (4) generation of a GENERAL problem; (5) and PARAMETER RUNS option. The primary features of each item of the major logic is discussed below.

- (1) EDIT feature: The first card of the deck is checked for the user request of the EDIT feature. If the EDIT feature is requested the input "tape" is changed from the system input "tape" to the EDIT "tape" and control is transferred to subroutine EDIT for processing. On return a branch is made to the THERMAL or GENERAL section as specified by the data on the EDIT "tape." If the EDIT feature has not been requested, the check for RECALL is made.
- (2) RECALL option: The first card of the deck is checked for user request of the RECALL option. If the RECALL option

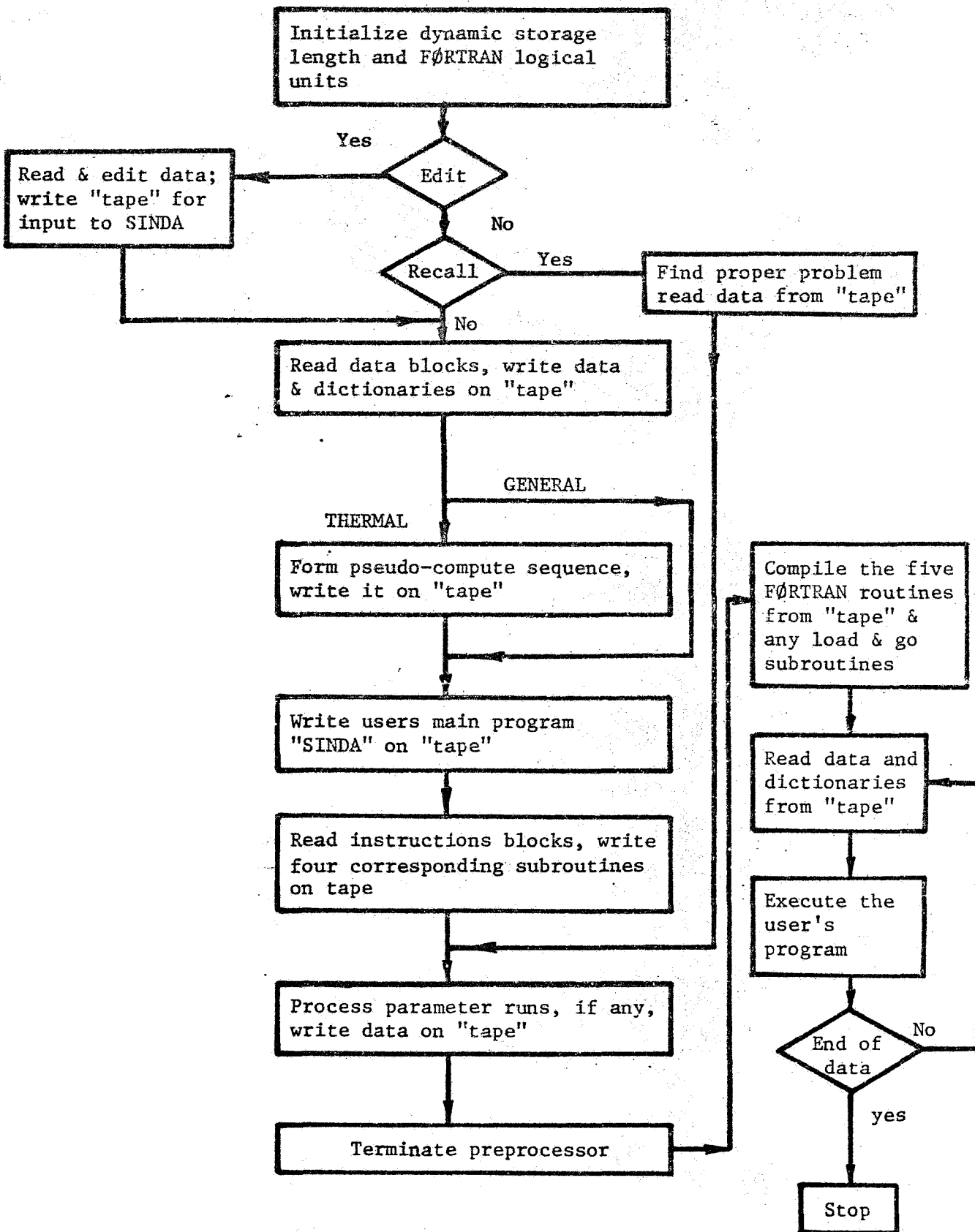


Figure 4.1-1. SINDA Preprocessor - Major Logic and Interface with the User Specified Problem

is requested, control is transferred to subroutine SPLIT for processing. On return a branch is made to the PARAMETER RUNS section. If the RECALL option has not been requested, the second card of the deck is checked for the type of problem, THERMAL or GENERAL.

- (3) THERMAL problem: The type of pseudo-compute sequence requested is noted, the title block is read, the data blocks are read and processed, the pseudo-compute sequence is formed, the driver for the user program (SINDA) is written on "tape," the operations blocks are read and processed and their FORTRAN equivalents are written on "tape," and finally a check is made for the user requests of the PARAMETER RUNS option.
- (4) GENERAL problem: This section is identical to a THERMAL problem except that only constants data and array data of the data blocks are read and processed; a pseudo-compute sequence is not formed.
- (5) PARAMETER RUNS option: A check is made for the user request of the PARAMETER RUNS option. If the PARAMETER RUNS option is requested, the appropriate data blocks are read and processed. If not, the preprocessor is terminated.

Description of SINDA preprocessor routines is presented in the sections to follow. Terminology used in the description is listed and defined in Table 4.1-1.

Table (4.1-1) Terminology Used in Description of
SINDA Preprocessor Routines

- (1) DATA BLOCKS: The five user input blocks which contain data rather than instructions; these DATA BLOCKS are NODE DATA, CONDUCTOR DATA, CONSTANTS DATA, ARRAY DATA and the optional blocks SOURCE DATA.
- (2) OPERATIONS BLOCKS: The four user input blocks which contain instructions on problem solution, as opposed to data contained in the DATA BLOCKS. These OPERATIONS BLOCKS are EXECUTION, VARIABLES 1, VARIABLES 2 and OUTPUT CALLS.
- (3) Non-fatal error: An error that does not terminate the preprocessor immediately. That is, the preprocessor will continue scanning the remaining cards of the input deck for errors. However, the user program will not be executed.
- (4) Fatal error: An error that terminates the run immediately.
- (5) N/A: Means not applicable.
- (6) "TAPE": The term "tape" in quotes is used to signify any external storage device. That is, any piece of computer hardware, excluding the central processor, on which data can be stored and retrieved. The three most familiar examples are: magnetic tape, drum and disk.
- (7) Dictionary: A list of the actual SINDA numbers in relative order. For example, the actual node number corresponding to the kth relative node number is the kth item of the node number dictionary.
- (8) Data group: A data group composed of the pertinent information extracted from a particular data block. For example, the two groups derived from the constants data are: the user constants numbers and the user constants values.
- (9) Bit manipulation: Terminology that implies the ability to store and access information within a computer word. This capability is also called packing and unpacking.
- (10) Routine: A general term used to describe any program element.
- (11) Subroutine: A special type of program element that is callable from a routine.
- (12) Fixed constants: The term used in the preprocessor for control constants.

4.2 Description of Subroutines

Sections 4.2.1 through 4.2.30 below describe the 30 routines of the SINDA preprocessor. The descriptions are based on the UNIVAC 1108 computer under the EXEC II operating system; it should be understood, however, that much of the information is machine-dependent and is dependent upon the facilities operating system. Note that the element named SINDA (Section 4.2.1) and the element named PREPRØ (Section 4.2.2) are not subroutines in the technical sense of the word; hence, these two elements are referred to by the more general term "routine."

Each element of the preprocessor is described by the following eleven subtitles:

- (1) SUBROUTINE NAME - this specifies the name of the element.
- (2) PROGRAMMING LANGUAGE - This may be FØRTRAN, ASM, or MAP. FØRTRAN implies FØRTRAN V, ASM stands for assembly language (sometimes called SLEUTH II) and MAP is a special language which defines the overlay structure.
- (3) PURPOSE - This gives a brief statement of the functional capabilities of the element.
- (4) RESTRICTIONS - This gives an indication of where the input parameters come from, the form of the input parameters and the placement of the output parameters.
- (5) "TAPES" USED - This represents a list of each FØRTRAN logical unit referenced within this element.
- (6) SPECIAL FEATURES - This specifies programming features that are unique to a particular machine.
- (7) OTHER SUBROUTINES CALLED - This represents a list of the external references.
- (8) CALLING SEQUENCE - This gives a list of the subroutine arguments, if any, and a brief discussion of their use.
- (9) ERROR PROCEDURES - This discusses the steps taken when an error is encountered.
- (10) STORAGE REQUIRED - This gives the octal and decimal storage required for this element.
- (11) LABELED COMMON - This represents a list of each labeled common name used in this element.

4.2.1 ROUTINE NAME: SINDA

PROGRAMMING LANGUAGE: MAP

PURPOSE: This routine specifies the overlay structure of the preprocessor to the system allocator (loader).

RESTRICTIONS: N/A

"TAPES" USED: N/A

SPECIAL FEATURES: N/A

OTHER SUBROUTINES USED: N/A

CALLING SEQUENCE: N/A

ERROR PROCEDURES: N/A

STORAGE REQUIRED: N/A

LABELED COMMON: N/A

4.2.2 ROUTINE NAME: PREPRØ

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This routine is the main routine (i.e., the driver) for the pre-processor. It initializes the counters and FORTRAN logical units, sets the length of the dynamic storage array, and controls the major logic. The major logic includes: (1) the EDIT feature (NASA MSC only); (2) the RECALL of a stored problem; (3) setup of a new user problem; (4) and preprocessor termination procedures.

RESTRICTIONS: N/A

"TAPES" USED:

System input "tape"	NIN
System output "tape"	NØUT
Problem data "tape"	LB3D
Problem FORTRAN "tape"	LB4P
Dictionary "tape"	LUT1
Parameter runs "tape"	LUT3
Recall "tape"	LUT7
Internal scratch "tape"	INTERN

SPECIAL FEATURES: System error termination - the problem data unit (LB3D) and the problem FORTRAN unit (LB4P) are flagged to stop before the data scan begins in the event that a system error terminates the preprocessor prematurely. The reason the problem data unit is flagged to stop is that for a RECALL problem the problem FORTRAN unit must not be written on.

OTHER SUBROUTINES USED: CØDERD, GENLNK, PRESUB, PSEUDØ, SINDA4, SPLIT and WRTBLK.

CALLING SEQUENCE: N/A

ERROR PROCEDURES: The error termination procedures are controlled by three flags named ERDATA, PRØGRM, and ENDRUN. The three flags are in the labeled common block named DATA. ERDATA is used to flag non-fatal errors encountered while reading the data blocks, while PRØGRM performs the same function for the operations blocks. See Section 4.7.2.

STORAGE REQUIRED: 443 octal words = 291 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CRDBLK, DATA, LØGIC, PLØGIC, and TAPE.

4.2.3 SUBROUTINE NAME: ALPINT

PURPOSE: This subroutine accepts an integer in the alphanumeric format nAl, and converts it to integer format, determines the relative number of this actual number, and converts the relative number back to an alphanumeric format of the form mAl.

RESTRICTIONS: The input and output is transmitted via the labeled common block named CIMAGE (see Section 4.3). The input must consist exclusively of the ten decimal digits.

"TAPES" USED: System output "tape" NØUT

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: SEARCH

CALLING SEQUENCE: ALPINT(KLET,IST,IEND,J)

KLET is an integer variable that indicates which dictionary is to be used for converting actual to relative.
IST is the starting location of the alphanumeric integer.
IEND is the ending location of the alphanumeric integer.
J points to the last location + 1 of the converted integer.

ERROR PROCEDURES: In the event that a given actual number has no relative number in the dictionary list, an error message will be issued and the relative operations blocks error flag (PRØGRM) will be set to 1.0.

STORAGE REQUIRED: 665 octal words = 437 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CIMAGE, DATA PØINT, and TAPE.

4.2.4 SUBROUTINE NAME: BLKCRD

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine formats the five generated FØRTRAN routines (SINDA, EXEC TN, VARBL1, VARBL2, and ØUTCAL) in 507 word blocks, which are acceptable to the FØRTRAN compiler. This information is stored in labeled common block CRDBLK, array KBLK. A complete discussion of the required tape format is found in UNIVAC 1108, EXEC II, Programmers Reference Manual, UP-4058 C, Appendix D.4 entitled, Program Elements on Magnetic Tape (via CUR).

RESTRICTIONS: The input is Hollerith card images with a 14A6 format. It is transmitted either through the array IMAGE in labeled common CRDBLK, or through "tape" INTERN.

"TAPES" USED: Internal scratch "tape" INTERN

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: STFFB and WRTBLK

CALLING SEQUENCE: BLKCRD

ERROR PROCEDURES: none

STORAGE REQUIRED: 753 octal words = 491 decimal words. See Section 4.7.1.

LABELED COMMON: CRDBLK and TAPE.

4.2.5 SUBROUTINE NAME: CØDERD

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine reads and checks the block header cards for the data blocks. It also performs the following functions: (1) the second data card of the deck is checked for a thermal or general problem, and if it is a thermal problem the type of pseudo-compute sequence specified is noted; (2) the title block is read and processed; (3) the actual array and constant numbers from the automated options are converted into FØRTRAN addresses; and (4) the parameter run block header cards are read and checked.

RESTRICTIONS: None

"TAPES USED:

System input "tape"	NIN
System output "tape"	NØUT
FØRTRAN V reread	30

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

CALLING SEQUENCE: CØDERD

ERROR PROCEDURES: In general, the errors checked for in this subroutine are of the fatal type; for example, data blocks out of order. The result of a fatal error is that the fatal error flag (ENDRUN) is set to 1.0 and control is returned to PREPRO for immediate termination.

STORAGE REQUIRED: 3213 octal words = 1675 words decimal. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, LØGIC, PLØGIC, and TAPE.

4.2.6 SUBROUTINE NAME: CØNVRT

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine converts Hollerith data to integer data.

RESTRICTIONS: The Hollerith data must be contained in one word and consist of only the ten decimal digits.

"TAPES" USED: None

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: ERRMES

CALLING SEQUENCE: CØNVRT(IST,IEND,ITEMP,CRDERR)

IST is the pointer to the first bit of the first character.

IEND is the pointer to the first bit of the last character.

ITEMP is the word containing the Hollerith data on entry and the integer number on return.

CRDERR is a logical error flag which is set true if an error is encountered during the conversion.

ERROR PROCEDURES: If a non-integer is encountered, an error message is printed and CRDERR is set to true.

STORAGE REQUIRED: 150 octal words = 104 decimal words. See Section 4.7.1.

LABELED COMMON: None

4.2.7 SUBROUTINE NAME: DATARD

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine scans the data block card images under an A format and determines the appropriate format (of the form Fn, In, or An) to reread the card image. The card images are then reread under the generated format. In addition, the constants data block and the array data block are processed.

RESTRICTIONS: None

"TAPES" USED:

System input "tape"	NIN
System output "tape"	NØUT
FØRTRAN V reread	30

SPECIAL FEATURES: The FORTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: ERMES, FINDRM, GENUK, NØDEDA (and its entry point CØNDDA), SETFMT, SQUEEZ, and TYPCHK.

CALLING SEQUENCE: DATARD

ERROR PROCEDURES: All errors checked for in this subroutine are non-fatal. An error message is printed either internally or from subroutine ERMES and the data blocks error flag (ERDATA) is set to 1.0.

STORAGE REQUIRED: 5344 octal words = 2788 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CHECKD, DATA, FLAGS, LØGIC, PLØGIC, PØINT, and TAPE.

4.2.8 SUBROUTINE NAME: ERRMES

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine prints most of the error messages that can be generated within the data blocks.

RESTRICTIONS: None

"TAPES" USED: System output "tape" NØUT

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: System subroutine EXIT

CALLING SEQUENCE: ERRMES (JUMP, I, J, K)

JUMP is an integer that points to the appropriate error message via a computed GØ TØ statement

I }
J } are data words which allow a maximum of three printed
K } data words per error message.

ERROR PROCEDURES: If the number of error messages printed exceeds 199, the preprocessor is terminated by a call to EXIT.

STORAGE REQUIRED: 2166 octal words = 1142 decimal words. See Section 4.7.1.

LABELED COMMON: DATA and TAPE

4.2.9 SUBROUTINE NAME: FINDRM

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine moves the data in the dynamic storage array either up or down by 100 words. In the process it may delete certain groups of data that are no longer needed.

RESTRICTIONS: None

"TAPES" USED: System output "tape" NØUT

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: SQUEEZ, and system subroutine EXIT.

CALLING SEQUENCE: FINDRM(LØCNØ,M)

LØCNØ is a pointer to a portion of the dynamic storage array where the data group that needs more room resides.

M is the address where the next data value is to be stored.

ERROR PROCEDURES: If the dynamic storage array is full, an error message is printed and the preprocessor is terminated via CALL EXIT.

STORAGE REQUIRED: 407 octal words = 263 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET LØGIC PØINT, and TAPE.

4.2.10 SUBROUTINE NAME: GENLNK

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine generates the driver, FORTRAN routine name and SINDA, for the user's program.

RESTRICTIONS: None

"TAPES" USED: Internal Scratch "tape" INTERN

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: BLKCRD

CALLING SEQUENCE: GENLNK

ERROR PROCEDURES: None

LABELED COMMON: CRDBLK, DATA, LOGIC, PLOGIC, and TAPE.

4.2.11 SUBROUTINE NAME: GENUK

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine is used to generate user constants.

RESTRICTIONS: The input data is taken from array TEMP in labeled common CHECKD and the output data (i.e., the generated user constants) is put into array B in labeled common BUCKET.

"TAPES" USED: None

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: ERRMES, FINDRM, and TYPCHK.

CALLING SEQUENCE: GENUK(IWRDS)

IWRDS is the number of words to be processed in array TEMP.

ERROR PROCEDURES: The input data is checked and if an error is found, control is transferred to subroutine ERRMES.

STORAGE REQUIRED: 451 octal words = 297 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CHECKD, DATA, and PØINT.

4.2.12 SUBROUTINE NAME: INCØRE

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine reads data into the dynamic storage array for the parameter runs option.

RESTRICTIONS: None

"TAPES" USED: Dictionary "tape" LUT1
Parameter runs "tape" LUT3

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: INCØRE(ITEST)

ITEST is an integer flag which determines the data group group to be read.

ERROR PROCEDURES: None

STORAGE REQUIRED: 600 octal words = 384 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, LØGIC, PLØGIC, PØINT, and TAPE.

4.2.13 SUBROUTINE NAME: MXTØFN

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine processes the data for the "M" option. That is, it converts card images from mixed FØRTRAN/SINDA notation to FØRTRAN notation.

RESTRICTIONS: The input (array IHØLL) and output (array JHØLL) are both in labeled common CIMAGE and they are both in an 80A1 format. The FØRTRAN from array JHØLL is copied to array IMAGE under a 14A6 format for processing by the FØRTRAN compiler.

"TAPES" USED: None

SPECIAL FEATURES: The FORTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: ALPINT and BLKCRD

CALLING SEQUENCE: MXTØFN

ERROR PROCEDURES: None

STORAGE REQUIRED: 522 octal words = 338 decimal words. See Section 4.7.1.

LABELED COMMON: CIMAGE and CRDBLK

4.2.14 SUBROUTINE NAME: NØDEDA

PROGRAMMING LANGUAGE FØRTRAN

PURPOSE: This subroutine processes the data for the node and conductor data blocks.

RESTRICTIONS: The input is received via labeled common CHECKD: array TEMP and the processed data are stored in the dynamic storage array.

"TAPES" USED: None

SPECIAL FEATURES: This subroutine has a second entry point named CØNDDA. Also, the FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: ERRMES, FINDRM, RELACT, and TYPCHK

GALLING SEQUENCE: NØDEDA(JUMP,IWRDS)

or CØNDDA(JUMP,IWRDS)

JUMP is a flag which indicates which code option (columns 8, 9, and 10 of the data card) the user has selected.

IWRDS is the number of data values in array TEMP to be processed.

ERROR PROCEDURES: If an error is detected while scanning the input data, control is transferred to subroutine ERRMES.

STORAGE REQUIRED: 7030 octal words 3608 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CHECKD, DATA, FLAGS, and PØINT.

4.2.15 SUBROUTINE NAME: PCS2

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine packs the FØRTRAN addresses for the array and constants locations required by the pseudo-compute sequence.

RESTRICTIONS: None

"TAPES" USED: None

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: PCS2 (IB, IPCS, LITA)

IB is the word in the dynamic storage array where the addresses are found.

IPCS is the word into which the addresses are packed.

LITA is a flag that is set to 1 if the array address was input as a literal and therefore has been added to the constants data.

ERROR PROCEDURES: None

STORAGE REQUIRED: 54 octal words = 44 decimal words. See Section 4.7.1.

LABELED COMMON: None

4.2.16 SUBROUTINE NAME: PRESUB

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine reads and checks the block header cards for the operations blocks and generates the non-executable FØRTRAN cards for each of the operations blocks via a call to BLKCRD.

RESTRICTIONS: None

"TAPES" USED: System input "tape" NIN
System output "tape" NØUT

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: BLKCRD

CALLING SEQUENCE: PRESUB(N)

N is an integer from 1 to 4 which indicates which operations block is being processed.

ERROR PROCEDURES: If the card read is not the correct block header card, an error message is printed and the fatal error flag is set to 1.0.

STORAGE REQUIRED: 200 octal words = 128 decimal words. See Section 4.7.1.

LABELED COMMON: CRDBLK, DATA, LØGIC, and TAPE.

4.2.17 SUBROUTINE NAME: PSEUDØ

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine forms the first and second pseudo-compute sequence. See Section 4.6.

RESTRICTIONS: The necessary input is extracted from the dynamic storage array and the output (the two pseudo-compute sequences) is placed in the dynamic storage array.

"TAPES" USED: NØRT

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: FINDRM, PCS2 and WRDTA.

CALLING SEQUENCE: PSEUDØ

ERROR PROCEDURES: If an error is encountered while forming the pseudo-compute sequences, an error message will be printed and the non-fatal error flag (ERDATA) is set to 1.0.

STORAGE REQUIRED: 2244 octal words = 1188 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, LØGIC, PLØGIC, and TAPE.

4.2.18 SUBROUTINE NAME: QDATA

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine checks and processes all data input in the source data block.

RESTRICTIONS: The input is received from the calling sequence and labeled common CHECKD. The processed data is placed in the dynamic storage array.

"TAPES" USED: None

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: ERRMES, FINDRM, RELACT, and TYPCHK.

CALLING SEQUENCE: QDATA(CØDE, IWRDS)

CØDE is the three letter option from columns 8, 9, and 10 of the data card.

IWRDS is the number of words in array TEMP to be processed.

ERROR PROCEDURES: If an error is encountered, control is transferred to subroutine ERRMES.

STORAGE REQUIRED: 1655 octal words = 941 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CHECKD, and PØINT.

4.2.19 SUBROUTINE NAME: RELACT

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine finds the relative node number from the actual node number. In addition, it computes the FØRTRAN address for arrays and user constants from the actual number.

RESTRICTIONS: This subroutine is used in conjunction with the data blocks.

"TAPES" USED: None

SPECIAL FEATURES: The FØRTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINE USED: ERRMES

CALLING SEQUENCE: RELACT(K,MM,J,JJ)

K determines the path through the program via a computed GØ TØ statement.

MM is the actual number on entry, and the FØRTRAN address on return.

J } are print variables for subroutine ERRMES.
JJ }

ERROR PROCEDURES: In the event an error is encountered, control is transferred to subroutine ERRMES.

STORAGE REQUIRED: 311 octal words = 201 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, and PØINT.

4.2.20 SUBROUTINE NAME: SEARCH

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine returns a relative number for nodes, conductors, user constants, and arrays, given the actual number.

RESTRICTIONS: This subroutine is used in conjunction with the operations blocks.

"TAPES" USED: None

SPECIAL FEATURES: The FORTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: SEARCH(N, IA, NDIM, LØC)

N is the actual number.

IA is the first word of the dictionary of actual numbers to be searched.

NDIM is the number of words of IA to be searched.

LØC is the relative number returned to the calling program.

ERROR PROCEDURES: If the input actual number is not found in the dictionary, LØC is set to zero.

STORAGE REQUIRED: 101 octal words = 65 decimal words. See Section 4.7.1.

LABELED COMMON: None

4.2.21 SUBROUTINE NAME: SETFMT

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine processes the cards for the "new format" option; that is, it sets up the format for data cards as specified by the cards with a N in column one.

RESTRICTIONS: The input/output array is passed through the calling sequence argument.

"TAPES" USED: FØRTRAN V reread 30

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: SETFMT(JUMP,B)

JUMP is an integer flag that determines the path through the code.

B is an array which contains the card images.

ERROR PROCEDURES: None

STORAGE REQUIRED: 221 octal words = 145 decimal words. See Section 4.7.1.

LABELED COMMON: None

4.2.22 SUBROUTINE NAME: SINDA4

PROGRAMMING LANGUAGE: FØRTRAN

PURPOSE: This subroutine reads and processes the user input cards from the operations blocks.

RESTRICTIONS: None

"TAPES" USED:

System input "tape"	MIN
System output "tape"	NØUT
Internal scratch "tape"	INTERN
FØRTRAN V reread	30

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: BLKCRD, MXTØFN, and SEARCH.

CALLING SEQUENCE: SINDA4(NAME)

NAME is an integer flag that tells the subroutine which operations block is being processed.

ERROR PROCEDURES: In the event an error is encountered while processing the operations blocks, an error message is printed and the error flag PROGRAM is set to 1.0.

STORAGE REQUIRED: 2372 octal words = 1274 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, CIMAGE, CRDBLK, DATA, LØGIC, PLØGIC, PØINT, and TAPE.

4.2.23 SUBROUTINE NAME: SKIP

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine is used when a problem is being RECALLED. It positions the tape to the proper problem as specified on the first card of the data deck.

RESTRICTIONS: The data is read from tape R. There is no output.

"TAPES" USED: RECALL "tape" LUT7

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: SKIP

ERROR PROCEDURES: None

STORAGE REQUIRED: 324 octal words = 212 decimal words. See Section 4.7.1.

LABELED COMMON: TAPE

4.2.24 SUBROUTINE NAME: SPLIT

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine reads the data from the RECALL tape and splits the RECALL information onto the program data "tape" (LB3D) and the dictionary "tape" (LUT1).

RESTRICTIONS: The input is from the RECALL "tape" and the output is placed on the program data "tape," the dictionary "tape," and the parameter runs "tape."

<u>"TAPES" USED:</u>	RECALL "tape"	LUT7
	Program data "tape"	LB3D
	Dictionary "tape"	LUT1
	Parameter runs "tape"	LUT3

SPECIAL FEATURES: None

OTHER SUBROUTINES CALLED: SKIP

CALLING SEQUENCE: SPLIT(ID)

 ID is the RECALL name punched in the first card of the data deck.

ERROR PROCEDURES: None

STORAGE REQUIRED: 746 octal words = 486 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, and TAPE.

4.2.25 SUBROUTINE NAME: SQUEEZ

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine compresses the specified data groups in the dynamic storage array. The compression is accomplished by placing the data groups sequentially in the dynamic storage array.

RESTRICTIONS: None

"TAPES" USED: None

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: SQUEEZ(IST,IEND)

IST is the data group number where the compression is to start.

IEND is the last data group number for this compression.

ERROR PROCEDURES: None

STORAGE REQUIRED: 115 octal words = 77 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET and POINT

4.2.26 SUBROUTINE NAME: STFFB

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine fills out a card image in array KBLK with Hollerith blanks.

RESTRICTIONS: The pointers to the words to set to blank are in the calling sequence, and the array containing the card images is in labeled common CRDBLK.

"TAPES" USED: None

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: STFFB(I,J)

I is the first word in KBLK to set to blank.

J is the last word in KBLK to set to blank.

ERROR PROCEDURES: None

STORAGE REQUIRED: 41 octal words = 33 decimal words. See Section 4.7.1.

LABELLED COMMON: CRDBLK.

4.2.27 SUBROUTINE NAME: TYPCHK

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine checks the input from the data blocks for the correct type; type means integer, floating point, or alphanumeric. Also, it regulates the conversion of the A's and K's for the automated options via a call to CONVRT.

RESTRICTIONS: The input and output are transferred through the calling sequence arguments and labeled common CHECKD.

"TAPES" USED: None

SPECIAL FEATURES: The FORTRAN V intrinsic function FLD is used for bit manipulation.

OTHER SUBROUTINES USED: CONVRT and ERRMES

CALLING SEQUENCE: TYPCHK(JUMP,IERR,J)

JUMP indicates what type the word should be.

IERR tells subroutine ERRMES which error message to print if the word is not of the type indicated by JUMP.

J is a pointer to the word type in array KFLFX.

ERROR PROCEDURES: If a word is not of the proper type control is transferred to subroutine ERRMES to print an error message and the logical flag CRDERR is set to true.

STORAGE REQUIRED: 233 octal words = 155 decimal words. See Section 4.7.1.

LABELED COMMON: CHECKD

4.2.28 SUBROUTINE NAME: WRDTA

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine writes the program data "tape" in the format required by INPUTT or INPUTG.

RESTRICTIONS: The data to be written on "tape" is found in the dynamic storage array.

"TAPES" USED: Program data "tape" LB3D

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: WRDTA(JUMP)

JUMP is an integer flag that indicates which data block to write and what format to use.

ERROR PROCEDURES: None

STORAGE REQUIRED: 645 octal words 421 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, LOGIC, PLOGIC, POINT, and TAPE

4.2.29 SUBROUTINE NAME: WRTPMT

PROGRAMMING LANGUAGE: FORTRAN

PURPOSE: This subroutine writes the data that is needed for parameter runs on the parameter runs "tape" and writes the dictionary "tape."

RESTRICTIONS: The information that is written on the "tapes" is found in the dynamic storage array.

"TAPES" USED: Dictionary "tape" LUT1
Parameter runs "tape" LUT3

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: WRTPMT(JUMP)

JUMP is an integer flag that indicates to WRTPMT which set of information to write.

ERROR PROCEDURES: None

STORAGE REQUIRED: 401 octal words = 257 decimal words. See Section 4.7.1.

LABELED COMMON: BUCKET, DATA, LOGIC, POINT, and TAPE.

4.2.30 SUBROUTINE NAME: WRTBLK

PROGRAMMING LANGUAGE: Assembly Language

PURPOSE: This subroutine writes the 507 word blocks contained in array KBLK on the program FORTRAN "tape."

RESTRICTIONS: None

"TAPES" USED: Program FORTRAN "tape" LB4P

SPECIAL FEATURES: None

OTHER SUBROUTINES USED: None

CALLING SEQUENCE: WRTBLK

ERROR PROCEDURES: None

STORAGE REQUIRED: 14 octal words = 12 decimal words. See Section 4.7.1.

LABELED COMMON: CRDBLK

4.3 Labeled Common Variables

The SINDA preprocessor uses nine labeled common blocks to pass data and flags between the various subroutines. Labeled common names, in alphabetical order, are:

BUCKET	CHECKED	CIMAGE
CRDBLK	DATA	FLAGS
LOGIC	PLAGIC	PPOINT

Note that the UNIVAC 1108 version does not utilize blank common. The two sections that follow give: 1) a map of the labeled common usage by subroutine name and by overlay link; 2) a definition of the variables used within each labeled common block; and 3) dynamic storage structure.

4.3.1 Labeled Common Map

The map below gives the labeled common name, a list of the overlay links that use it by link number and a list of the routines that use it.

<u>Labeled Common Name</u>	<u>Overlay Link Names</u>	<u>Routine Names</u>
BUCKET	0, 1, 2, 4, 5	ALPINT
		CHDERD
		DATARD
		FINDRM
		GENUK
		INCORE
		NODEDA
		PREPRO
CHECKD	1	PSEUDO
		QDATA
		RELACT
		SINDA4
SPLIT	4	SQUEEZ
		WRTPMT
CRDBLK	0, 3, 4	WRTPMT
		DATARD
		GENUK
CIMAGE	4	NODEDA
		QDATA
ALPINT	0, 3, 4	TYPCHK
		MXTOFN
		SINDA4
		BLKCRD
CRDBLK	0, 3, 4	GENLNK
		PREPRO
		PRESUB
STFFB	0, 3, 4	SINDA4
		WRTBLK

<u>LABELED COMMON NAME</u>	<u>OVERLAY LINK NAMES</u>	<u>ROUTINE NAMES</u>	
DATA	0, 1, 2, 3, 4, 5	ALPINT	CØDERD
		DATARD	ERRMES
		GENLINK	GENUK
		INCØRE	NØDEDA
		PREPRØ	PRESUB
		PSEUDØ	RELACT
		SINDA4	SPLIT
		WRTDTA	WRTPMT
FLAGS	1	DATARD	NØDEDA
LØGIC	0, 1, 2	CØDERD	DATARD
		FINDRM	INCØRE
		PREPRØ	PSEUDØ
		WRTDTA	WRTPMT
PLØGIC	0, 1, 3, 4	CØDERD	DATARD
		GENLNK	INCØRE
		PREPRØ	SINDA4
		WRTDTA	
PØINT	0, 1, 2, 4	ALPINT	CØDERD
		DATARD	FINDRM
		GENUK	INCØRE
		NØDEDA	PREPRØ
		PSEUDØ	QDATA
		RELACT	SINDA4
		SQUEEZ	WRTDTA
		WRTPMT	
TAPE	0, 1, 2, 3, 4, 5	ALPINT	BLKCRD
		CØDERD	DATARD
		ERRMES	FINDRM
		GENLNK	INCØRE
		PREPRØ	PRESUB
		PSEUDØ	SINDA4
		SKIP	SPLIT
		WRTDTA	WRTPMT

4.3.2

Definition of Labeled Common Variables

(1) Labeled common name BUCKET.

BUCKET is the dynamic storage array (see Section 4.3.3).

(2) Labeled common name CHECKD

CHECKD is used to temporarily store and check the user's input data.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
TEMP(35) or ITEMP	A temporary storage array, which contains the user's input data as read from the data cards.
XGEN(35) or IGEN	A temporary storage array used to store a copy of TEMP when the user is generating data.
KFLFX(35)	An indicator used to check the data array which contains one of the following numbers: 1 = floating point number 0 = integer number -1 = Hollerith word
CRDERR	A logical flag: Set true if and only if an error was found on the data card now being processed.

(3) Labeled common name CIMAGE

CIMAGE is used to store and manipulate Hollerith card images for the 'M' option.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
IHOLL(80)	The input card image, read under an 80A1 format.
JHOLL(160)	The constructed card image, also an 80A1 format.

(4) Labeled common name CRDBLK

CRDBLK is used to construct the five generated FØRTRAN routines.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
LSTART	A logical flag that signals the start of a new routine if set to true.
LECARD	A logical flag that signals the end of a routine.
LCØPY	A logical flag that tells the program to copy the card image (14A6) found in IMAGE to the next available slot in KBLK.
NW	Is a counter whose value is the next available word in KBLK.
KBLK(507)	An array that contains FØRTRAN card images of the generated routines.
IMAGE(14)	An array that contains one card image to be copied into KBLK.

(5) Labeled common name DATA

DATA is used to store the counters that indicate (to the program) how many of each data type has been encountered. In addition, it contains three error flags.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
NND	The number of diffusion nodes.
NNA	The number of arithmetic nodes.
NNB	The number of boundary nodes.
NNT	The total number of nodes.
NGL	The number of linear conductors.
NGR	The number of radiation conductors.
NGT	The total number of conductors.
NUC	The number of user constants.
NEC1	The number of added constants from automated options in the node data block.

NEC2	The number of added constants from the automated options in the conductor data block.
NCT	The total number of constants.
LENA	The total number of words used in the array data block.
ERDATA	The non-fatal error flag for the data blocks. ERDATA \neq 0.0 means an error has been found.
PRØGRAM	The non-fatal error flag for the operations blocks. PRØGRAM \neq 0.0 indicates an error condition.
ENDRUN	The fatal error flag for the preprocessor. ENDRUN \neq 0.0 signals the program to terminate immediately.
LSEQ1	The length of the first pseudo-compute sequence.
LSEQ2	The length of the second pseudo-compute sequence.
LØNG	A logical flag set to true if the user is requesting the long pseudo-compute sequence.

(6) Labeled common name FLAGS

FLAGS contains three flags that are used to go to the proper block of coding in subroutine NØDEDA.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
LEAP	Used with the GEN option.
NØNLIN	Flags a set of multiply connected conductors as radiation, if set to true.
INDX	Determines path when multiply connected conductors require more than one data card.

(7) Labeled common name LØGIC

LØGIC contains a number of logical flags and the fifty fixed constants.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
LNØDE	Set to true if any node data was processed.
LCØND	Set to true if any conductor data was processed.
LCØNST	Set to true if any user constants were processed
LARRAY	Set to true if any array data was processed.
LPRINT	Debug print flag, set to true if there is an asterisk in column 80 of the BCD 3 THERMAL/GENERAL card.
KBRNCH	An integer that specifies which data block is being processed.
FIXC(50) or IFIXC	The array that contains the fixed (control) constants.
KTPRNT	Optional print flag for a list of relative versus actual user constant numbers. Set true if there is an asterisk in column 80 of the BCD 3CØNSTANTS DATA card.
AYPRNT	Optional print flag for a list of actual array numbers versus FØRTRAN address. Set true if there is an asterisk in column 80 of the BCD 3ARRAY DATA card.
GENERL	Set true for a general problem.
LQ	Set true if any data was processed from the source data block.

(8) Labeled common name PLØGIC

PLØGIC contains a number of logical flags that are used in conjunction with parameter runs.

<u>VARIABLE NAME</u>	<u>DESCRIPTION</u>
PARINT	Set true for initial parameters run.
PARFIN	Set true for final parameters run.
PNØDE	Set true if node data was processed.
PCØND	Set true if conductor data was processed.
PCØNST	Set true if user constants data was processed.

PARRAY	Set true if array data was processed.
PTITLE	Set true if a new title was input.
PCHGID	Contains the alphanumeric word INITIAL or FINAL to be used as the run identification on "tape" LB3D.

(9) Labeled common name PØINT

PØINT is used in conjunction with dynamic storage array BUCKET. See Section 4.3.3.

4.3.3. Dynamic Storage Structure

Dynamic storage represents one of the techniques of maximizing problem size with a computer with finite core. In dynamic storage each data set is placed sequentially into one array end-to-end. This eliminates the wasted core inherent with the traditional system of dimensioning each variable at some fixed length. However, the price paid for the additional core is the extra time required to compute the address of a variable.

The SINDA preprocessor used three arrays to store and address the data sets. The data sets are stored in an array named B, or IB, or BB. This array resides in labeled common BUCKET. The length at which B can be dimensioned depends on the system that the computer facility uses. At NASA MSC approximately 30,000 words are allocated to B. In addition, in labeled common PØINT there are two arrays named LØC and LEN, each dimensioned at 20. LØC (I) contains the starting location in B for the Ith data set and LEN (I) contains the length of the Ith data set.

The information below gives, in detail, the contents of the dynamic storage array for each data block as it exists just after the data block has been processed.

(1) Node data block

data set 1:

bit 1,	automated option flag
bit 2,	Q from SOURCE DATA flag
bits 16-35,	actual node number

data set 2:
 bits 0-35, temperature value

data set 3:
 bits 0-35, capacitance value

data set 4:
 bits 0-5 non-linear capacitance type
 bit 6, literal array flag
 bits 7-20, actual array number
 bit 21, literal constant flag
 bits 22-35, actual constant number

data set 5:
 bits 0-35, literals encountered in 4.

(2) Source data block

data set 2 (first word of group):
 bits 0-5, source option type
 bits 6-20, relative node number
 bits 21-35, not used

data set 2 (second word of group):
 bits 0-5, not used
 bit 6, literal array flag
 bits 7-20, actual array number
 bit 21, literal constant flag
 bits 22-35, actual constant number

data set 3:
 bits 0-35, literals encountered in 2

(3) Conductor data block

data set 6:
 bits 0-35, actual conductor number

data set 7:
 bit 0, multi connections flag
 bit 1, radiation flag
 bit 2, automated option flag

- | | |
|-------------|-------------------------|
| bits 3-5, | not used |
| bit 6, | 1 way flag for NA |
| bits 7, 20, | relative node number NA |
| bit 21 | 1 way flag for NB |
| bits 22-35, | relative node number NB |
- data set 8:
- | | |
|------------|-------------------|
| bits 0-35, | conductance value |
|------------|-------------------|
- data set 9:
- | | |
|-------------|------------------------|
| bits 0-5, | conductor option type |
| bit 6, | literal array flag |
| bits 7-20, | actual array number |
| bit 21, | literal constant flag |
| bits 22-35, | actual constant number |
- data set 10:
- | | |
|------------|---------------------------|
| bits 0-35, | literals encountered in 9 |
|------------|---------------------------|
- (4) Constants data block
- data set 11:
- | | |
|------------|------------------------|
| bits 0-35, | actual constant number |
|------------|------------------------|
- data set 12:
- | | |
|------------|----------------|
| bits 0-35, | constant value |
|------------|----------------|
- (5) Array data block
- data set 13:
- | | |
|------------|---------------------|
| bits 0-35, | actual array number |
|------------|---------------------|
- data set 14:
- | | |
|------------|--------------|
| bits 0-35, | array length |
|------------|--------------|
- data set 15:
- | | |
|------------|-------------|
| bits 0-35, | array value |
|------------|-------------|
- (6) Pseudo-compute sequences
- data set 16 (1st pseudo-compute sequence):
- | | |
|--------|----------------------------|
| bit 0, | last conductor flag |
| bit 1, | automated capacitance flag |
| bit 2, | automated conductance flag |

bit 3,	radiation conductance flag
bit 4,	Q from source block flag
bits 5-20,	relative conductor number
bit 21	1 way conductor flag
bits 22-35,	relative adjoining node number

data set 17 (second pseudo-compute sequence):

bits 0-4,	automated option type
bit 5,	not used
bits 6-21,	FØRTRAN address for array
bit 22,	not used
bits 23-35,	relative constant number

The bit numbering convention above conforms to the UNIVAC standard notation, where each 36 bit word is numbered 0 through 35 from left to right. Each of the 1 bit flags above is queried in the following manner: 0 means NO, and 1 means YES. If the literal array flag or the literal constant flag is set to 1, then the bits immediately to the right of the flag do not contain the actual array or constant number. Instead, they contain a pointer to the next data set where the literal value is stored. In those data sets that store information for the automated options it is sometimes necessary to use more than one word per option. When this is the case, the automated option type (bits 0-5) is set to 0.

4.4 SINDA "Tapes" and Their Formats

The SINDA program in its normal operating mode utilizes six "tapes." Five of these "tapes" are assigned by the program and the remaining one contains the program; it is assigned via control cards. The store and recall options require one additional "tape" each and the NASA edit feature requires two additional "tapes." The following paragraphs contain information on the five normal SINDA "tapes."

4.4.1 LB3D - Program Data "Tape"

This "tape" is set up by the preprocessor (WRTDTA) and read by INPUTT, for a thermal problem, or INPUTG, for a general problem, just prior to performing the instructions of the execution block. The contents of this unit are:

- (1) Problem identification.

```
WRITE(LB3D)RUNID
```

- (2) Title information (20 words).

```
WRITE(LB3D)(TITLE(I),I=1,20)
```

- (3) The number of: diffusion nodes, arithmetic nodes, and total nodes; followed by a temperature value for each node; then a capacitance value for each diffusion, if any.

```
WRITE(LB3D)NND,NNA,NNT,(T(I),I=1,NNT)
```

```
IF(NND.GT.0)WRITE(LB3D)(C(I),I=1,NND)
```

- (4) The total number of conductors followed by a conductor value for each one.

```
WRITE(LB3D)NGT,(G(I),I=1,NGT)
```

- (5) The total number of user constants are followed by the 50 control constant values; then the user constants values, if any.

```
WRITE(LB3D)NCT,(FIXC(I),I=1,50)
```

```
IF(NCT.GT.0)WRITE(LBJD)(K(I),I=1,NCT)
```

- (6) The total number of arrays and the overall length of the array data; then the array values, if any.

```
WRITE(LB3D)NAT,LENA
```

```
IF(LENA.GT.0)WRITE(A(I),I=1,LENA)
```


- (7) The lengths of the first and second pseudo-compute sequences, followed by the data for the first pseudo-compute sequence; then the data for the second pseudo-compute sequence, if any.

```
WRITE(LB3D)LSEQ1,LSEQ2,(P1(I),I=1,LSEQ1)
IF(LSEQ2.GT.0)WRITE(LB3D)(P2(I),I=1,LSEQ2)
```

Note that (3), (4), and (7) above apply only to a thermal problem.

4.4.2 LB4P - Program FORTRAN "Tape"

This "tape" is especially formatted in 507 word blocks as required by the FORTRAN compiler. Where:

WORD 1 on the first block of each routine contains the name of the routine.

WORD 2 contains the integer number of card images in the block.

WORDS 3 - 506 contain the card images

WORD 507 is set to +0 except on the last block of each routine where it is set to -0.

4.4.3 INTERN - Preprocessor Scratch "Tape"

Generally INTERN is used to pass card images to subroutine BLKCRD under a 14A6 format.

4.4.4 LUT1 - Dictionary "Tape"

This "tape" contains a list of the actual SINDA numbers in a relative order. That is, the actual node number corresponding to the kth relative node number is the kth item of the node number dictionary. The format of this "tape" is:

- (1) The total number of nodes, followed by an actual node number for each node.

```
WRITE(LUT1)NNT,(NN(I),I=1,NNT)
```

- (2) The total number of conductors, followed by the list of actual conductor numbers.

```
WRITE(LUT1)NGT,(NG(I),I=1,NGT)
```

- (3) The number of user constants, the total number of constants, followed by a list of the actual constant numbers.

```
WRITE(LUT1)NUC,NCT,(NK(I),I=1,NCT)
```

- (4) The total number of arrays followed by a list of the actual array numbers, then the total number of arrays followed by a list of the length of each array.

```
WRITE(LUT1)NAT,(NA(I),I=1,NAT)
```

```
WRITE(LUT1)NAT,(LA(I),I=1,NAT)
```

4.4.5 LUT3 - Parameter Runs "Tape"

This "tape" contains some data from the original problem. It is required by the initial parameters capability. The format of "tape" LUT3 is:

- (1) The original title.

```
WRITE(LUT3)(TITLE(I),I=1,20)
```

- (2) A list of original temperature and capacitance values.

```
WRITE(LUT3)NND,(T(I),I=1,NNT)
```

```
IF(NND.GT.0)WRITE(LUT3)(C(I),I=1,NND)
```

- (3) A list of the original conductor values.

```
WRITE(LUT3)(G(I),I=1,NGT)
```

- (4) Lists of the original fixed and user constants.

```
WRITE(LUT3)NUC,NCT,(FIXC(I),I=1,50)
```

```
IF(NCT.GT.0)WRITE(LUT3)(K(I),I=1,NCT)
```

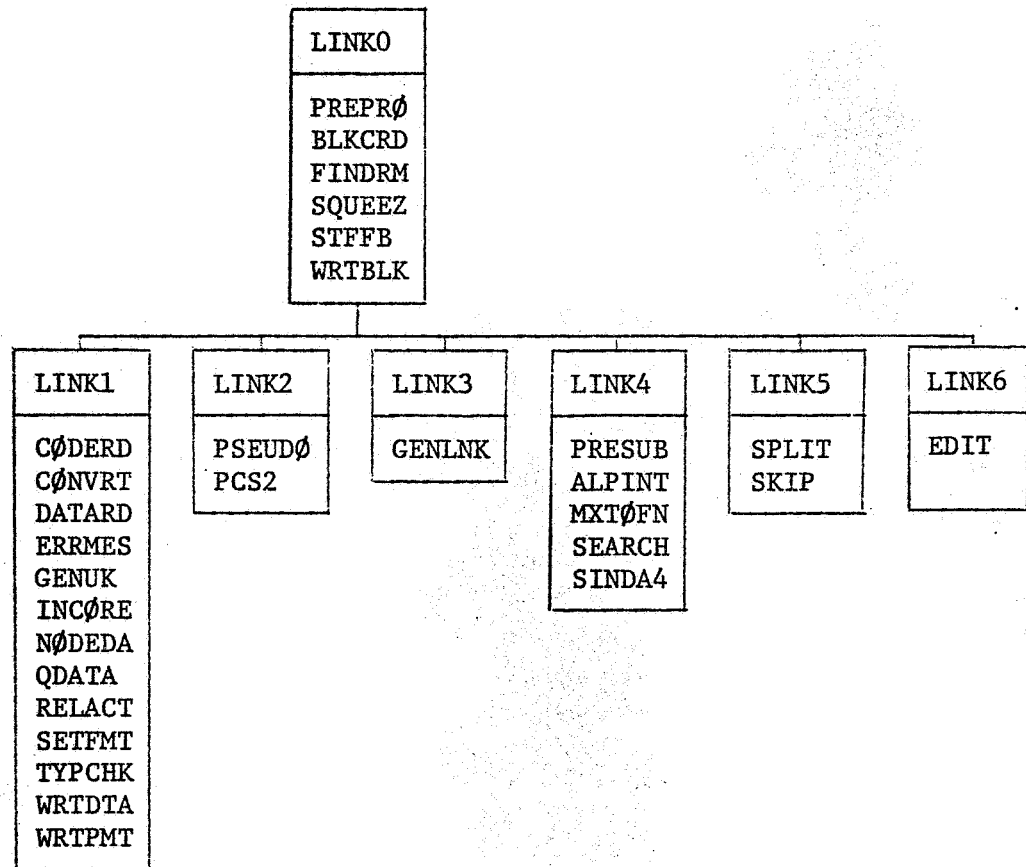
- (5) The original array values.

```
WRITE(LUT3)NAT,LENA
```

```
IF(LENA.GT.0)WRITE(LUT3)(A(I),I=1,LENA)
```

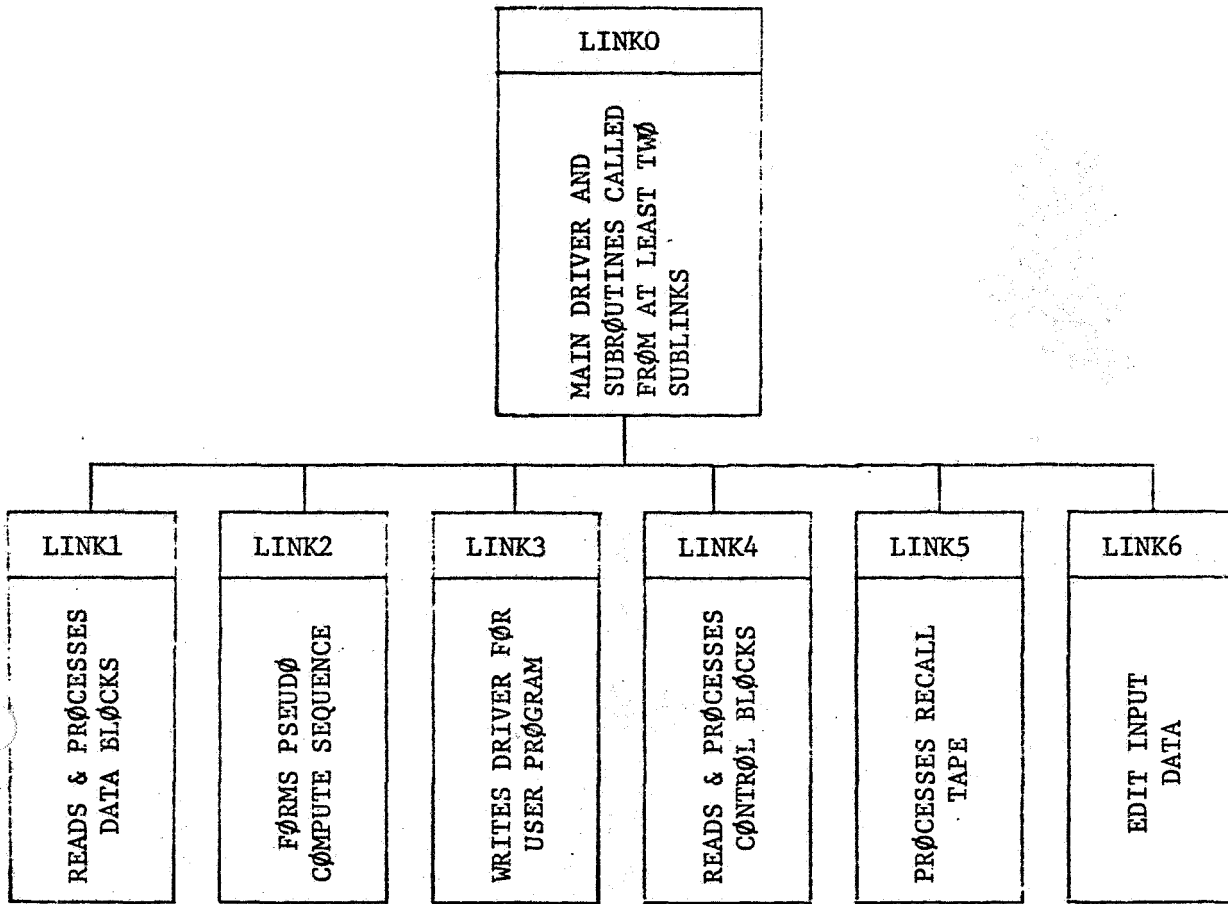
4.5 Overlay Structure

The SINDA preprocessor has an overlay structure composed of a main link (designated LINK0 below) which is always in core and five sublinks (designated LINK1, LINK2, LINK3, LINK4, LINK5, and LINK6 below) which overlay one another as they are brought into core.



Note that the first subroutine listed above in each of the sublinks serves as the driver for that sublink and it is also the subroutine called from PREPRØ.

Another approach to overlay specification is to think of each link as a functional unit, hence the graph below.



4.6 Structure of Pseudo-Compute Sequences

4.6.1 Descriptions

The use and structure of the two pseudo-compute sequences generated by the SINDA preprocessor appear to be rather confusing and mysterious. The term "pseudo" itself leads to immediate interpretation difficulties. Suppose that an element G_k between nodes i and j is to be identified for the i th node; by specifying explicitly i, j , and k the element is completely defined. On the other hand, SINDA explicitly specifies j and k but i is implicit in the DØ-LØØP; hence, the description pseudo-compute sequence (PCS) arises. Confusion also arises from the lack of information regarding the need for the (PCS) and the difficulties in reading the packed information. In short, the PCS as used in SINDA is simply two lists of relative numbers which are ordered in a specific manner. The two lists of relative numbers form the heart of the PCS, although other information pertinent to the computation must also be considered.

The PCS is necessary because the data as input by the user does not lend itself efficiently to the computational capabilities of FØRTRAN. As a result, the preprocessor scans the user input data and places the relative numbers (FØRTRAN addresses) into an array in the order in which the data will be used at a later time by the user selected numerical solution routine.

Packing of the data is a technique that conserves computer storage by placing two or more pieces of information in one computer word. This allows the user to execute a larger problem than the one that can be accommodated if the traditional one computer word for each piece of information approach. The penalty for this larger problem capability is an increase in execution time required for the extraction of information each time it is used.

4.6.2 Structure of PCS1

The first PCS, designated PCS1, contains the following information:

bit 0	last G for this node flag
bit 1	automated C option flag
bit 2	automated G option flag
bit 3	radiation G flag
bit 4	Q from source data flag
bits 5-20	relative G number
bit 21	1 way G flag
bits 22-35	relative adjoining node number

The 36 bits of each computer word are numbered 0 through 35 from left to right. All of the 1 bit flags are set such that 0 means NØ and 1 means YES.

PCSl is stored in an array named NSQ1 and is ordered by relative node number. That is, for relative node number 1 the conductor data is scanned and each time a conductor connected to node number 1 is encountered the PCS1 information is stored in NSQ1. When all of the conductor data has been scanned, the 0 bit of the latest word of PCS1 information is set to 1. This process is repeated for relative node numbers 2, 3, etc., until all diffusion and arithmetic nodes have been processed. PCS1 will be formed as either long or short as specified by the user on the BCD 3THERMAL card. This option is applied to diffusion nodes only since arithmetic nodes are always formed under the long option. The difference between the long PCS and the short PCS is that in the long PCS, each conductor will be listed twice, whereas in the short PCS each conductor will be listed once. This assumes the conductor connects two diffusion nodes. If one or both of the nodes is arithmetic, then the conductor will be listed twice, and if one of the nodes is a boundary the conductor will only be listed once. For example, given conductor number k which connects diffusion nodes i to j, where $i < j$. The long PCS would contain the k, j information for the processing of node i, and the k,i information for the processing of node j; whereas, the short PCS would only contain the k,j information. The short PCS thus has the advantage of requiring less computer storage than the long PCS, but a block iterative method (refer to Section 5.2.2) must be used; in general, the short PCS requires more iterations to converge than the successive point iterative (refer to Section 5.2.2) method which requires the long PCS.

4.6.3 Structure of PCS2

The second PCS is designated PCS2. The following information is stored whenever bit 1, bit 2, or bit 4 of PCS1 is set to one.

bits 0-4	automated option code
bit 5	not used
bits 6-21	FØRTRAN address of the array or relative constant number.
bit 22	not used
bits 23-35	relative constant number

If the automated option is a doublet type, like DIV, and therefore requires two words to store the information, the automated option code on the second word is set to zero. In the event that more than one of the flag bits (bits 1, 2, or 4) of PCS1 is set to one, then the following order is imposed on PCS2: the capacitance information is stored first, the source block information second and finally the conductor information.

The PCS2 information is stored in array named NSQ2. This array is the same under the long PCS1 or the short PCS1 since automated conductors are only flagged on their first encounter.

4.7 Other Information

This section contains miscellaneous information that may be of interest to the user.

4.7.1 Subroutine Lengths

The storage required by a particular routine will vary depending on the type of computer and the system being used. The routine lengths given in Section 4.2 are based on compiler listings made on 23 January 1971 at Jacobi Computation Center.* The machine is a UNIVAC 1108 with a highly modified system. The numbers represent the sum of the computer storage for computer instructions, constants, and simple variables.

4.7.2 Maximum Thermal Problem Size and Maximum Data Value Size

A short formula for estimating the maximum thermal problem size that can be run on SINDA, and a list of the maximum size of the various data values is given below.

* Now called Computation and Systems Corporation, Los Angeles, California.

Estimation of Maximum Problem Size

$$NNT + 3*NGT + NCT + 4*NA\emptyset < \underline{LENBKT}$$

where,

- NNT is the total number of nodes.
- NGT is the total number of conductors.
- NCT is the total number of constants (user constants plus literals from automated options).
- NA \emptyset is the number of automated options specified.
- LENBKT is the length of the dynamic storage array as set in routine PREPR \emptyset .

Maximum Size of Data Values

Actual node number	
Core storage	$2^{33}-1$
Print out	999,999
Relative node number	
Core storage	16,383
Temperature	
Core storage	$\pm 10^{38}$
Capacitance	
Core Storage	$\pm 10^{38}$
Relative conductor number	
Core storage	$2^{35}-1$
Actual user constants number	
Core storage	32,767
Automated options	16,383
Relative user constants number	
Core storage	32,767
Automated options	8,191

User constant values	
Integer	$\pm 2^{35}-1$
Floating point	$\pm 10^{38}$
Alphanumeric	6 characters
Actual array number	
Core storage	$2^{35}-1$
Automated options	16,383
Print out	99,999
Relative array number	
Core storage	$2^{35}-1$
Automated options	65,535
Print out	99,999
Array values	
Integer	$\pm 2^{35}-1$
Floating point	$\pm 10^{38}$
Alphanumeric	6 characters

Note that some of the maxima, such as the relative conductor number of $2^{35}-1$, are strictly academic since the dynamic storage array is considerably smaller than the indicated maximum data value size.

5. REVIEW OF LUMPED PARAMETER EQUATIONS AND BASIC NUMERICAL SOLUTIONS

The use of SINDA as mentioned in a previous section is based on a lumped parameter representation of a physical system.⁷ Thus SINDA solves numerically a set of ordinary (in general nonlinear) differential equations that represent the transient behavior of a lumped parameter system or a set of nonlinear algebraic equations representing steady state conditions. Numerous numerical solution techniques are reported in literature; a few of these are listed in the Reference Section.⁸⁻²⁴ These numerical methods are based on finite difference algorithms as opposed to finite element methods which have received considerable attention recently.²⁵⁻²⁹ For problems that are generally encountered in spacecraft thermal design, use of the finite element method appears to be inappropriate because of the nonlinearity presented with radiation heat transfer and because of complex geometric configurations.

Variations of the basic finite difference algorithms are numerous because no single numerical solution technique is optimum for all the endless types of thermal problems that can be encountered. Furthermore, because of the nonlinearity of the problems, a specific set of criteria to indicate solution accuracy and stability is not available and does not appear to be forthcoming. As a result, the user is placed in a rather awkward and confused position of not knowing which subroutine to use if a choice is available. Some thermal analyzer-type computer programs allow no choice, as a result, user decision is not necessary. SINDA represents a computer program at the other extreme of user decision flexibility by providing a number of numerical solution methods.

The intent of this Section 5 is to review and formulate the basic numerical solution methods with the presentation (from an engineering standpoint) of the characteristics of each SINDA numerical solution routine deferred to Section 6. In addition to place the use of SINDA in a proper perspective relative to accurate temperature prediction of a physical system, difficulties associated with lumped-parameter representation are discussed here.

5.1 Lumped Parameter Representation

Reduction of a distributive (physical) system to a lumped system which can be represented as an equivalent thermal network is a rather important phase of thermal analysis. From a temperature accuracy standpoint lumping (or nodalization) of the physical system may be far more important than a numerical solution technique that is used in a computer program. The latter is often given undue attention with apparent ignorance of other error sources which may be far more important. A general discussion on lumped parameter representation is not intended for presentation here since the subject material is extensively covered in technical literature, but it is convenient for continuity to indicate basic considerations.

For simple geometries and linear problems, it is rather straightforward to solve the partial differential equations of the type,

$$\frac{\partial T}{\partial t} = \alpha \nabla^2 T + Q \quad (5.1-1)$$

where, α = thermal diffusivity (k/C)

T = temperature

Q = source

$$\nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} \quad (\text{two dimensional})$$

Numerous analytical solutions of (5.1-1) for different types of boundary conditions and geometries are available.^{30, 31} Finite difference algorithms formed directly from the partial differential equations are also abundantly reported in literature.^{12, 14} These finite difference formulations were generally developed for well-defined geometries and symmetrical discretization. For these problems, the so-called nodal connections or resistances are immediately available and, in general, automatically generated by the computer program. Thus, the need for a lumped parameter representation does not exist. For these types of problems, inaccuracies due to truncation and solution stability are specifically established.

For complex geometries and nonlinear problems such as those that include thermal radiation exchange, analytical solutions of thermal problems are limited.^{32, 34} As a result, it is a common practice because of practical considerations to nodalize a physical system directly with-

out undue consideration of inaccuracies. Thus, a user merely represents the heat flow between two connecting nodes by using the basic network building block,

$$q_{ij} = (T_i - T_j)/R_{ij} \quad (5.1-2)$$

where, R_{ij} represents an effective resistance between adjoining nodes i and j .

It should be particularly noted here that SINDA employs the concept of conductance in lieu of resistance which is common with most network-type computer programs. Thus the heat flow is represented as:

$$q_{ij} = G_{ij}(T_i - T_j) \quad (5.1-3)$$

where, G_{ij} is the conductance between node i and node j .

The proper value of G_{ij} (or R_{ij}) for an arbitrary nodalization is (and should be) of concern to the user but because of the multitude of variables that must be considered, any discussion here would be incomplete. An excellent article on asymmetrical finite difference networks is presented in Reference 35.

5.1.1 Some Thoughts on Lumped Parameter Errors

Reduction of a physical system to a topological model consisting of a network with resistors and capacitors requires considerable engineering judgment. More often than not, nodal size of a model is governed by budget and schedule constraints. As a result, the discrete areas larger than desired are often used. This does not necessarily mean, however, that the use of a large number of nodes will always yield realistic results since the uncertainties of the input parameters can be appreciable.³⁶

Spatial truncation errors are controlled by selecting the grid size so that nonlinear temperature distributions lie within required accuracy by linear interpolation between nodal points, and that variation of temperature-dependent properties over the volume of each node is within required limits of the average values determined for the nodal point. The assumption of linear temperature distribution assumed for the lumped-parameter equations (equation 5.1-6) leads to a spatial truncation error of the order $O(\Delta x^2)$ only if all nodes are symmetrically located.²⁷ If a non-uniform grid is used, the accuracy of computation is only $O(\Delta x)$. Spatial truncation errors are thus inherent in the mathematical model and beyond user control once inputted into

SINDA. For a two-dimensional problem with symmetrical grids, the spatial truncation error can be expressed for typical explicit and implicit methods^{38, 12, 27}

$$E = -\frac{(\Delta x)^2}{12} \frac{\partial^4 T}{\partial x^4} - \frac{(\Delta y)^2}{12} \frac{\partial^4 T}{\partial y^4} + O(\Delta x^4) \quad (5.1-4)$$

Temperature distribution other than linear can also be formulated;³⁸ however, most thermal analyzer-type computer programs such as TRUMP³⁹ and including SINDA are based on the linear assumption.

Time truncation errors are directly dependent upon the time-step since the error for the typical explicit and implicit method is,

$$E = -\frac{\Delta t}{2} \frac{\partial^2 T}{\partial t^2} \quad (5.1-5)$$

Normally the time step is dependent upon a particular criterion chosen by the user. A more detailed discussion on user control of the time step will be given for each numerical solution routine within the SINDA sub-routine library.^{3, 4}

Another approximation error which is due to discretizing is the assumption of constant radiosity for the discrete areas. Inaccuracies can be expected to affect the level and distribution of temperature. The analysis of thermal radiation exchange has received considerable attention in recent years because of its importance in spacecraft thermal design.^{40, 52} The influence of non-uniform local heat flux on overall heat transfer between a gray differential area parallel to a gray infinite plane is examined in Reference 43; the assumption of uniform local heat flux appears to be reasonable for this geometry and for the evaluation of the overall heat flux calculations. A method of analysis suitable for engineering applications is developed in Reference 50 for computing local radiant flux and local temperature of opaque surfaces in a space environment. A study evaluating the validity of commonly used simplified methods of radiant heat transfer analysis is reported in Reference 48. A study directed at improving the understanding and prediction of orbiting spacecraft thermal performance is presented in References 46 and 49. A method presented in Reference 51 provides a means of evaluating the uncertainties associated with thermal radiation exchange. For an excellent status review (as of 1969) on radiation exchange between surfaces and in enclosures, the reader should consult Reference 52.

The above discussion merely serves to indicate that considerable care must be given when nodalizing a physical system and that the numerical evaluation of the finite difference equations must be considered from the total temperature error context. This means that user attention to a given numerical solution must be placed in a proper perspective.

5.1.2 Lumped-Parameter Equations

Using the network building block as expressed by equation (5.1-3) the lumped parameter system is identified as a set of ordinary non-linear differential equations by taking a heat balance as an *i*th node,

$$\frac{dT_i}{dt} = \frac{1}{C_i} \left[q_i + \sum_{j=1}^p a_{ij} (T_j - T_i) + \sum_{j=1}^p \sigma b_{ij} (T_j^4 - T_i^4) \right] \quad (5.1-6)$$

$$i = 1, 2, \dots, N \text{ (number of variable temperatures)}$$

$$T_j = \text{constant, } N < j \leq p$$

- where,
- C_i = the *i*th nodal capacity which may be a function of temperature
 - q_i = the heat into node *i* and may be a function of time and temperature (impressed)
 - a_{ij} = the conduction coefficient between nodes *i* and *j*; it may be a function of time and temperature
 - b_{ij} = the radiation coefficient between nodes *i* and *j*; it may be a function of time and temperature
 - σ = Stefan-Boltzmann constant

Coefficients a_{ij} and b_{ij} are SINDA input quantities with the temperature factor of equation (5.1-6) calculated internally by the program. Both a_{ij} and b_{ij} may be variables. Conductance updating is a subject for discussion in a later paragraph. The user requirement to input the coefficients, a_{ij} and b_{ij} , provides considerable program flexibility, but at the same time user generation of these input quantities presents, in some instances, rather difficult engineering judgment decisions.

Radiation coefficient b_{ij} is, in essence, a radiation interchange factor, F_{ij} ,⁵³⁻⁵⁵ (also known as script F) between nodes *i* and *j*. Generation of this quantity analytically can be quite difficult and inaccurate. A number of methods and computer programs (see, for example,

Reference 56) are available for evaluating the shape factors which represent an important part of determining script F. A direct generation of script F is normally through the use of the Monte Carlo technique,⁴⁹ but a recent development utilizes a matrix formulation for determining the script F in an enclosure containing surfaces with arbitrary emission and reflection characteristics.^{57, 58} An experimental technique is reported in Reference 59.

5.2 Basic Finite Difference Formulations

The various numerical solution techniques differ in the finite difference formulations for the time-derivative (refer to equation 5.1-6); since the thermal equation is of the parabolic type, the transient heat transfer problems are of the initial value type. This means that at some time point, $t = n\Delta t$ (n is the number of time steps, Δt) all values of T_i are known. Thus,

$$T_{i,n+1} = T_{i,n} + \left[\frac{dT_{i,n}}{dt} \right]_{t_{n,n+1}} \Delta t \quad (5.2-1)$$

$$i = 1, 2, \dots, N$$

where $t_{n,n+1}$ represents the time interval between $t = n\Delta t$ and $t = (n+1)\Delta t$

It is apparent that the selection of the proper value to $(dT_{i,n}/dt)_{t_{n,n+1}}$ cannot be explicit and its selection identifies one numerical method from another. Although many finite difference formulations of the parabolic differential equation are available, two general classifications are commonly denoted as explicit or implicit. These numerical methods are well-documented in literature; the reader should refer to Reference 12 for a comprehensive discussion on various finite difference approximations. Explicit methods also discussed in References 14, 17, 19 and 20, among others, are step-by-step in time and equations.

Explicit methods include:

- (1) Forward-difference explicit approximation^{12, 14}

This is an Euler method that computes temperatures in a step-by-step fashion. The requirement of stability places an upper limit on the time increment. SINDA subroutines CNFRWD, CNFRDL and CNFAST fall within this

category. CNFAST is a modified version of CNFRWD which allows the user to specify the minimum time step to be taken. Refer to Sections 6.3.1 and 6.3.2 for details.

(2) Dufort-Frankel approximation^{9, 12, 17}

The Dufort and Frankel finite difference formulation is a three level formula that appears to be unconditionally stable. SINDA subroutine CNDUFU uses the Dufort-Frankel finite difference algorithm (refer to Section 6.3.4 for a detailed discussion).

(3) Exponential approximation^{1, 17}

The exponential approximation is found by integrating the heat balance equations after making linear and constant coefficient assumptions. This method is unconditionally stable for linear systems but may be unstable for some types of nonlinear problems. SINDA subroutine CNEXPN employs this method and is discussed at length in Section 6.3.3.

(4) Alternating direction approximations¹⁷

This technique employs two formulations, one on odd time levels and the other on even time levels and is unconditionally stable.

The implicit finite difference formulations require a simultaneous computational procedure. In addition to Reference 12, implicit methods are also discussed in References 8, 10, 17, and 20, among others. Implicit methods include:

(1) Backward difference implicit approximation¹²

The backward difference weights only the flux terms at $t = (n+1)\Delta t$. As a result, the method is stable for all values of Δt . SINDA subroutine CNBACK employs this method and is detailed in Section 6.4.1.

(2) Crank-Nicolson approximation⁸

The Crank-Nicolson method uses the arithmetic average of the heat flux at the two time levels, $t = n\Delta t$ and

$t = (n+1)\Delta t$. The method is unconditionally stable. SINDA CNFWBK uses this method and is discussed in Section 6.4.2.

Steady state analysis also requires an implicit method of solution. SINDA steady state subroutines are called CINDSS, CINDSL and CINDSM which are detailed in Sections 6.5.1, 6.5.2, and 6.5.3.

5.2.1 Forward Finite Difference Explicit Method

By replacing the first derivative of temperature with respect to time, dT/dt , with the forward first difference quotient, equation (5.1-6) becomes,

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_i - \sum_{j=1}^P a_{ij} (T_{j,n} - T_{i,n}) + \sum_{j=1}^P \sigma b_{ij} (T_{j,n}^4 - T_{i,n}^4) \quad (5.2-1)$$

$$t = n\Delta t$$

$$i = 1, 2, \dots, N$$

$$T_{j,n} = \text{constant}, N < j \leq p$$

where, the second subscript on T represents the time level such that

$$T_{i,n} = T_i(n\Delta t)$$

Equation (5.1-6) is represented in the form expressed by equation (5.1-3) by letting,

$$G_{ij} = a_{ij} + \sigma b_{ij} (T_i^2 + T_j^2)(T_i + T_j) \quad (5.2-2)$$

It is interesting to note that the finite difference form of (5.1-6) (and thus 5.2-1) represents a second central-difference quotient of $\nabla^2 T$ (refer to (5.1-1)).

The computational procedure for the forward difference formulation is rather straightforward since only a single unknown temperature at each time step, $T = n\Delta t$ for each equation is present. Note that the averaging of dT_i/dt assigns a weighting factor to the heat flux terms only (terms on the right side of equation (5.1-6) at $t = n\Delta t$). Along with the computational simplicity, however, is the stability constraint which places an upper limit on the time increment, Δt , that can be used in the numerical procedure. The stability criterion for the explicit finite difference method

is (for the most limiting node),^{12, 14}

$$\Delta t < C_i / \sum_{j=1}^P G_{ij} \quad (5.2-3)$$

$$i = 1, 2, \dots, N$$

A modified stability criterion that allows for a larger time step which results in a conditionally stable temperature for the most limiting node is reported in Reference 23. Since the stability criterion will govern the maximum time step that can be used, it is thus particularly important that a user gives some attention to those factors that compose the condition of stability when nodalizing a physical system.

In the discussion presented so far, arithmetic nodes (nodes with no heat capacity) have not been mentioned. Normally, the computational procedure treats arithmetic nodes separately from the diffusion nodes; arithmetic-node temperatures are solved implicitly. Detailed discussion on the general procedure will be presented in a later paragraph as well as in Section 6 which discusses the various SINDA numerical solution routines.

5.2.2 Implicit Finite Difference Method

The implicit difference equations can be constructed for heat transfer problems in many ways (see, for example, References 12 and 20).

Replacement of equation (5.1-6) with the backward time difference yields,

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_i + \sum_{j=1}^P a_{ij} (T_{j,n+1} - T_{i,n+1}) + \sum_{j=1}^P \sigma b_{ij} (T_{j,n+1}^4 - T_{i,n+1}^4) \quad (5.2-4)$$

$$i = 1, 2, \dots, N$$

$$T_{j,n+1} = \text{constant}, N < j \leq p$$

$$T_{i,n} = T_i (n\Delta t)$$

The computational procedure for the backward difference formulation must necessarily be re-iterative because of the need to solve a set of simultaneous non-linear equations.

In view of the importance of iteration techniques (such as method of

successive approximation), it may be of interest to formulate equation (5.2-4) into an iterative form. If we let $C_i/\Delta t \equiv \bar{C}_i$, use equation (5.2-2) in equation (5.2-4) and solve the resultant expression for $T_{i,n+1}$, this yields the recurrent equation for a given time increment, Δt , and time-step, n ,

$$T_{i,k+1} = \frac{\bar{C}_{i,k} T_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k} + q_{i,k}}{\bar{C}_{i,k} + \sum_{j=1}^P G_{ij,k}} \quad (5.2-5)$$

$i = 1, 2, \dots, N$

where, $\bar{C}_{i,k} = C_{i,k}/\Delta t$

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,k}^2 + T_{i,k}^2)(T_{j,k} + T_{i,k}) \quad (5.2-6)$$

$T_{j,k} = \text{constant}, N < j \leq P$

$k = k\text{th iteration (note that } \bar{C}_{i,k}, q_{i,k}, a_{ij} \text{ and } b_{ij} \text{ are shown to be updated every iteration; SINDA routines update these quantities once each time-step)}$

The iterative pattern is initiated by assuming "old" temperatures ($T_{i,k}$ and $T_{j,k}$) on the right side of equation (5.2-5) to evaluate a "new" set of temperatures ($T_{i,k+1}$) on the left side of the equation (5.2-5); this single set of calculations represents an iteration. By replacing all of the "old" temperatures ($T_{i,k}$) on the right side of equation (5.2-5) with the just calculated "new" set of temperatures ($T_{i,k+1}$), a second iteration can be made. The iteration procedure is continued until a termination criterion such as the number of iterations or the maximum absolute difference between $T_{i,k}$ and $T_{i,k+1}$ is less than some prespecified value has been satisfied. It should be noted that G_{ij} , C_i and q_i are shown to be updated every iteration. This iterative process is termed "block" iteration since the "old" temperatures on the right side are replaced in a "block" (a set of temperatures) fashion with the "new" temperatures.

Another iterative technique is to utilize on the right side of equation (5.2-5) each "new" temperature as soon as it is calculated. This iterative method is termed "successive point" iteration and appears to yield solutions about 25% faster than the "block" iteration method.

Equation (5.2-5) can be expressed in a "successive point" form as follows:

$$T_{i,k+1} = \frac{\bar{C}_{i,k} T_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^p G_{ij,k} T_{j,k} + q_{i,k}}{\bar{C}_i + \sum_{j=1}^p G_{ij,k}} \quad (5.2-7)$$

where, $G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$
 $(\ell = k \text{ if } j \geq i \text{ and } \ell = k+1 \text{ if } j < i)$

$T_{j,k} = \text{constant, } N < j \leq p$

$k = k\text{th iteration (note that } \bar{C}_{i,k}, q_{i,k}, a_{ij}, \text{ and } b_{ij} \text{ are shown to be updated every iteration)}$

The iterative method as used in SINDA follows a fixed, pre-determined sequence of operations in contrast with a relaxation procedure which is also one of successive approximations but is not processed out in a predetermined sequence. The relaxation procedure seeks and operates on the node with the maximum temperature difference between the "old" and the "new." From a programming standpoint, the search operation requires as much computational time as the temperature calculation itself.

5.2.3 Steady State Method

Standard steady state equations follows directly from equation (5.2-5) for block iteration or from equation (5.2-7) for successive point iteration by letting $\bar{C}_i = 0$ in these equations. The comments made in Section 5.2.2 are equally applicable here.

5.2.4 Some Comments

The finite difference expressions presented in this Section 5.2 represent standard formulations and thus do not show computational techniques and artifices which are used, some more or some less, in all programs. SINDA numerical solution routines contain many computational features (many original with J. D. Gaski) which enhance problem solution. The various computational aspects of the numerical solution methods as used in the SINDA routines are discussed in rather lengthy detail in Section 6.

6. SINDA NUMERICAL SOLUTION ROUTINES

6.1 Objective and Presentation Format

SINDA has available to the user a number of numerical solution routines which employ various numerical methods. A brief description of these routines with the required SINDA input quantities and format are contained in the SINDA users manual;^{3,4} a general review of numerical methods was presented in Section 5. Unfortunately, the brief description is not sufficient for a casual SINDA user to make a selection decision from among several routines that are available and for a serious user to fully understand the computational procedure as well as to understand the role of the various control constants that are employed in each routine.

It is the intent of this section to fill wherever possible and practical the description void that presently exists with the numerical solution routines by detailing the characteristics of each. It is not the intent here to provide sufficient detailed information for a user to make modifications and/or additions to the existing subprograms, but rather to provide information that will aid the user in assessing the various numerical solution routines and in evaluating the numerical results.

Each of the numerical solution routines is detailed from a theoretical as well as from a computational standpoint. Control constants and their role are described and indicated in a step-by-step verbal flow computational procedure. Details of many of the numerous computational checks have purposely been omitted because of the complex interactions. Minute details of each routine can be obtained only from the individual computer listings; a computer listing of each of the SINDA numerical solution routines is presented in Appendices A, B and C. General computational procedure and features that apply to most, if not all of the SINDA numerical solution routines are assembled in a single section (6.2) in order to eliminate undue repetition. The description of each routine is heavily dependent upon, and coupled to, the general description of Section 6.2. The routines have been categorized as steady state or transient with the latter subcategorized as explicit or implicit in order to allow for an orderly presentation as well as to simplify future additions.

6.2 General Computational Procedure and Features

Each of the SINDA numerical solution routines employs a particular finite difference approximation of the lumped parameter heat balance equations. In spite of the uniqueness of each routine, portions of the computational procedure used in each are similar. Also, many of the routines have identical features such as the acceleration of convergence and the use of control constants. As a result, it is convenient to place in this section repetitious material. In some instances material presented here is repeated in the discussion of a particular numerical solution routine.

6.2.1 Order of Computation

It was reported in Section 3.5 that the order of computation depends on the sequence of subroutine calls placed in the EXECUTION block by the program user. No other operations block is performed unless called upon by the user either directly by name or indirectly from subroutines which internally call upon them. Numerical solution subroutines internally call upon operations blocks VARIABLES 1, VARIABLES 2, and OUTPUT CALLS. The internal order of computation for these numerical solution routines is similar with the primary difference between one routine and another being the finite difference approximation employed in a particular routine. A flow diagram indicating the general order of computation for the numerical solution routines is depicted in Figure 6.2-1.

6.2.1.1 Finite Difference Algorithm

Although each of the SINDA numerical solution routines employs a particular finite difference approximation which is detailed for each numerical solution routine, the computational pattern is similar. Within the box depicted as SFDA in Figure 6.2-1, solution of the finite difference algorithm occurs. The computational sequence for transient solutions follows one of two patterns: (1) one for explicit finite difference methods; and (2) one for implicit finite difference methods; steady state solutions follow closely the implicit pattern. Both numerical flow pictures are depicted in Figure 6.2-2; details within the flow pictures are different for each routine and are described separately under the individual SINDA numerical solution routines (refer to Sections 6.3 - 6.5).

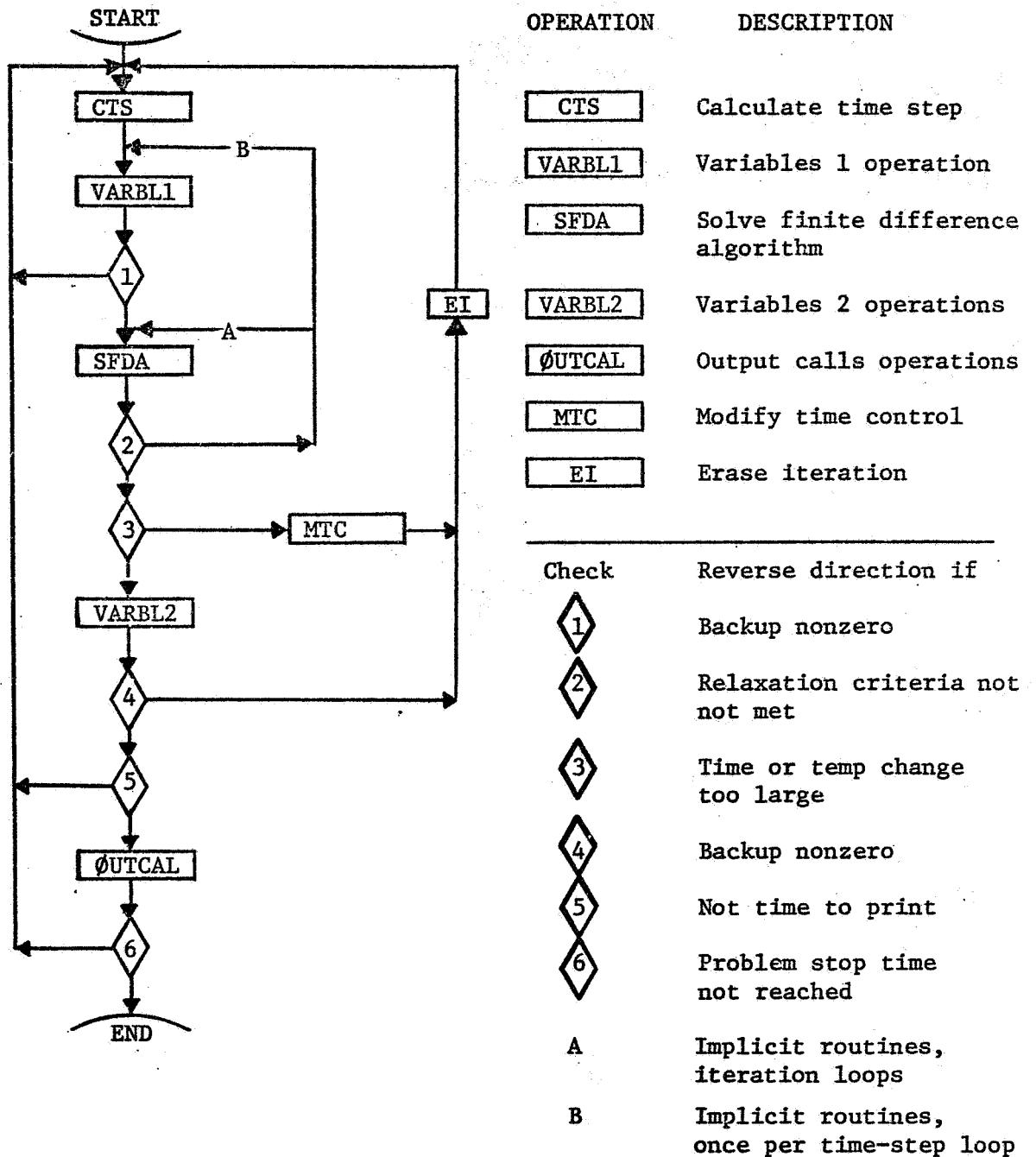
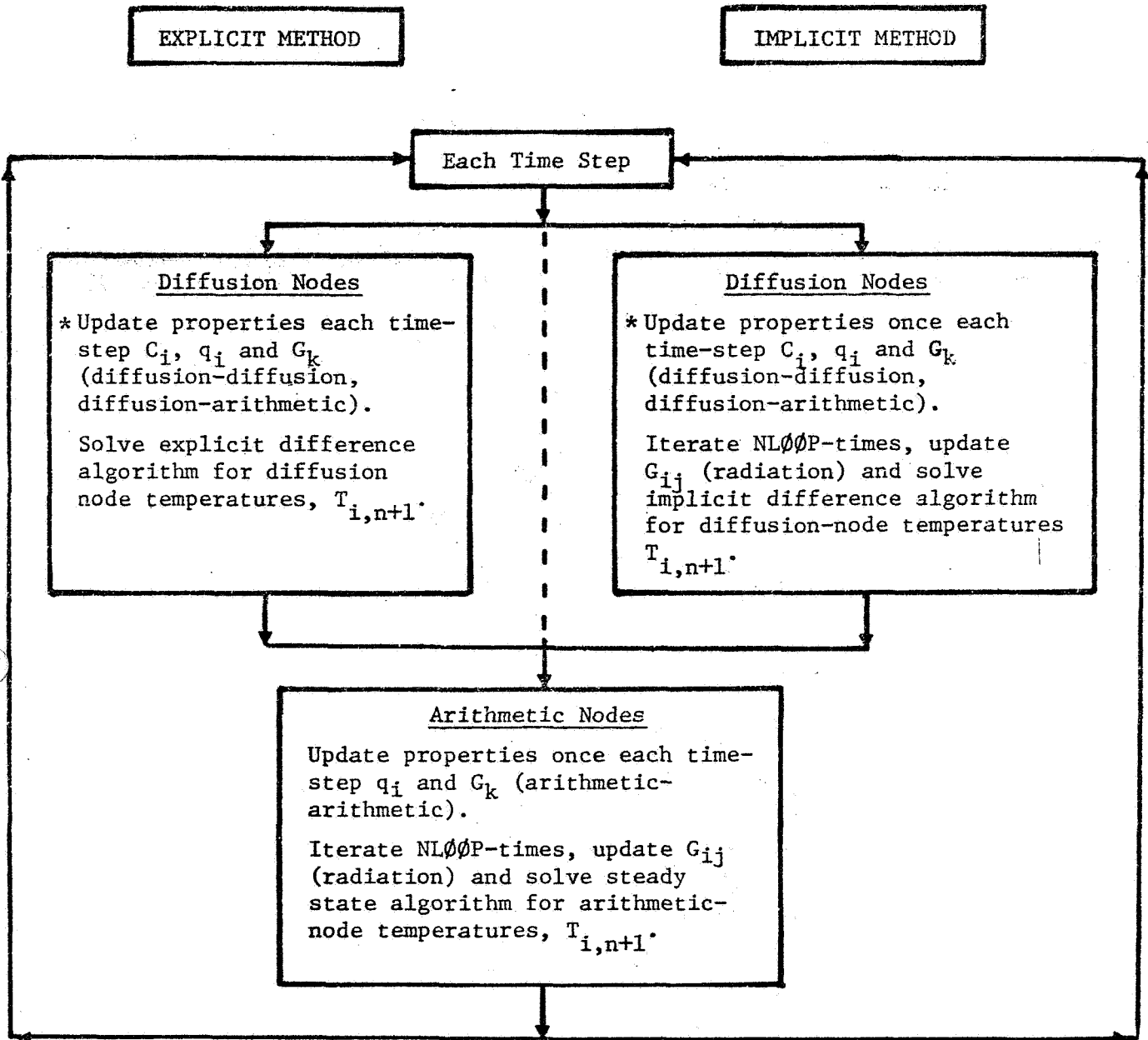


Figure (6.2-1) General Order of Computation for Numerical Solution Routines



* For CINDA -3G users, it should be noted that the updating of properties occurs within the numerical solution routine after VARIABLES 1 call. CINDA-3G evaluates the variable properties before VARIABLES 1 call.

Figure 6.2-2. Numerical Computational Pattern for Explicit and Finite Difference Algorithms

6.2.1.2 Updating of Optionally Specified Properties

Optionally specified properties are defined here as those items which result in pointers being set in the second pseudo compute sequence (refer to Section 3.3.4). The term optional refers to mnemonic options that are available for different types of variable properties.^{3, 4} The properties are updated in all SINDA numerical solution routines the same way. This definition is used here in lieu of stating that optionally specified properties are time and/or temperature varying properties since source data may be specified to be constant. The pointers are set by one or more of the following user input quantities:

- (1) All capacitances, C_i , specified as $f(T)$ or $f(t,T)$ in NODE DATA BLOCK:
- (2) All data, q_i , entered in the SOURCE DATA BLOCK:
- (3) All coefficients, G_k , specified as $f(T)$ or $f(t,T)$ in CONDUCTOR DATA BLOCK. It should be noted here that the term coefficient as used here requires amplification. The conductance, G_{ij} , may be for conduction or for radiation; that is,

$$G_{ij} \equiv G_k = a_{ij} \quad (\text{for conduction conductance})$$

$$G_{ij} = ob_{ij} (T_i^2 + T_j^2)(T_i + T_j) \quad (\text{for radiation conductance})$$
$$= G_k (T_i^2 + T_j^2)(T_i + T_j)$$

Thus, note that the calculated conduction conductance G_{ij} is identical to the updated G_k , whereas for the calculated radiation conductance only ob_{ij} is equivalent to the updated G_k .

The type of optional properties is identified by the integer stored in the first six bits of the second pseudo compute sequence which indicates to the program which option is in effect. Optional property types are listed and described for the three categories of input quantities in Table 6.2-1 for capacitance, Table 6.2-2 for impressed source, and Table 6.2-3 for coefficients with the definition of symbols listed in Table 6.2-4.

6.2.2 Operations Blocks

In a previous paragraph, it was mentioned that the sequence of subroutine calls placed in the EXECUTION block by the user determines the

TABLE 6.2-1 OPTIONALLY SPECIFIED CAPACITANCE EXPRESSIONS

<u>Option</u>	<u>Type</u>	<u>Expression</u>
SIV	1	$C_i = F(A^i:T_i)$
DIV	2	$C_i = F1(A_1^i:T_i) + F2(A_2^i:T_i)$
DIV	3	$C_i = F1(L) + F2(A^i:T_i)$
DIV	4	$C_i = F1(A^i:T_i) + F2(L)$
SPV	5	$C_i = F(A^P:T_i)$
DPV	6	$C_i = F1(A_1^P:T_i) + F2(A_2^P:T_i)$
DPV	7	$C_i = F1(L) + F2(A^P:T_i)$
DPV	8	$C_i = F1(A^P:T_i) + F2(L)$
BIV	9	$C_i = F(A^b:T_i, t_m)$

Notation: Refer to Table 6.2-4.

TABLE 6.2-2 OPTIONALLY SPECIFIED IMPRESSED SOURCE EXPRESSIONS

<u>Option</u>	<u>Type</u>	<u>Expression</u>
blank	1	$q_i = q_i + F$
SIV	2	$q_i = q_i + F(A^i:T_i)$
SIT	3	$q_i = q_i + F(A^i:t_m)$
DIT	4	$q_i = q_i + F1(A_1^i:t_m) + F2(A_2^i:t_m)$
DIT	5	$q_i = q_i + F1(L) + F2(A^i:t_m)$
DIT	6	$q_i = q_i + F1(A^i:t_m) + F2(L)$
DTV	7	$q_i = q_i + F1(A_1^i:t_m) + F2(A_2^i:T_i)$
DTV	8	$q_i = q_i + F1(L) + F2(A^i:T_i)$
DTV	9	$q_i = q_i + F1(A^i:t_m) + F2(L)$

Notation: Refer to Table 6.2-4.

Table 6.2-3. Optionally Specified Coefficient Expressions for Conduction and Radiation

Mnemonic Options	Type	Expression
SIV	1	$G_k = F(A^i:T_m)$
SIV	2	$G_k = F(A^i:T_i)$
DIV (conduction)	3	$G_k = 1.0/[1.0/F1(A_1^i:T_i) + 1.0/F2(A_2^i:T_j)]$
(radiation)		$G_k = [F1(A_1^i:T_i)][F2(A_2^i:T_j)]$
DIV (conduction)	4	$G_k = 1.0/[1.0/F1(L) + 1.0/F2(A^i:T_j)]$
(radiation)		$G_k = [F1(L)][F2(A^i:T_j)]$
DIV (conduction)	5	$G_k = 1.0/[1.0/F1(A^i:T_i) + 1.0/F2(L)]$
(radiation)		$G_k = [F1(A^i:T_i)][F2(L)]$
SPV	6	$G_k = F(A^P:T_m)$
SPV	7	$G_k = F(A^P:T_i)$
DPV (conduction)	8	$G_k = 1.0/[1.0/F1(A_1^P:T_i) + 1.0/F2(A_2^P:T_j)]$
(radiation)		$G_k = [F1(A_1^P:T_i)][F2(A_2^P:T_j)]$
DPV (conduction)	9	$G_k = 1.0/[1.0/F1(L) + 1.0/F2(A^P:T_j)]$
(radiation)		$G_k = [F1(L)][F2(A^P:T_j)]$
DPV (conduction)	10	$G_k = 1.0/[1.0/F1(A^P:T_i) + 1.0/F2(L)]$
(radiation)		$G_k = [F1(A^P:T_i)][F2(L)]$
BIV	11	$G_k = F(A^b:T_m, t_m)$
SIV	12	$G_k = F(A^i:T_j)$
SPV	13	$G_k = F(A^P:T_j)$

Notation: Refer to Table 6.2-4; note $G_k \equiv ob_{ij}$ (for radiation)
 $\equiv a_{ij}$ (for conduction)

Table 6.2-4. Definition of Symbols for Tables 6.2-1 -- 6.2-3

<u>Symbols</u>	<u>Definition</u>
C_i	Capacitance of ith node.
F, F1, F2	Multiplying factors, either user constants or literal
$G_k (=a_{ij})$	Conduction coefficient.
$G_k (=ob_{ij})$	Radiation coefficient.
L	A literal multiplying factor.
q_i	Heat load into the ith node. (impressed)
Δt	Time-step
t_m	Mean time, $(TIME\phi + TIMEN)/2.0$
T_m	Mean temperature, $(T_i + T_j)/2.0$
$(A^i:t_m)$	Interpolated value of array A using t_m as the independent variable.
$(A^i:T_i)$	Interpolated value of array A using T_i as the independent variable.
$(A^b:T_i, t_m)$	Interpolated value of the bivariate array A using T_i and t_m as independent variables.
$(A^b:T_m, t_m)$	Interpolated value of the bivariate array A using T_m and t_m as independent variables.

Mnemonic Options

BIV	<u>B</u> ivariate <u>I</u> nterpolation <u>V</u> ariable
DIT	<u>D</u> ouble <u>I</u> nterpolation with <u>T</u> ime as variable
DIV	<u>D</u> ouble <u>I</u> nterpolation <u>V</u> ariable
DPV	<u>D</u> ouble <u>P</u> olynomial <u>V</u> ariable
DTV	<u>D</u> ouble interpolation with <u>T</u> ime and <u>T</u> emperature as <u>V</u> ariables
SIT	<u>S</u> ingle <u>I</u> nterpolation with <u>T</u> ime as variable
SIV	<u>S</u> ingle <u>I</u> nterpolation <u>V</u> ariable
SPV	<u>S</u> ingle <u>P</u> olynomial <u>V</u> ariable

Subscripts

i	Indicates the ith node.
j	Indicates the jth node.
2	Indicates two (array).

order of computation. Operations blocks number four, EXECUTION, VARIABLES 1, VARIABLES 2, and OUTPUT CALLS. These operations blocks are described in the SINDA Users Manual^{3, 4} but their role insofar as the numerical solution routines are concerned may be of particular interest.

6.2.2.1 EXECUTION Operations Block

The EXECUTION operations block provides the user considerable flexibility in the use of SINDA calls and FORTRAN operations. Combinations of SINDA calls and FORTRAN operations are innumerable since the user is actually programming. Now all instructions contained in the VARIABLES 1, VARIABLES 2 and OUTPUT CALLS are performed each iteration or on the output call interval. Thus, if an operation being performed in VARIABLES 1 utilizes and generates non-changing constants, the operation should be placed in the EXECUTION block (prior to the numerical solution call) so that it will be performed only once and thus eliminate repetitious non-changing calculations. Operations of this type are conveniently performed in the EXECUTION operations block. Note, however that a constant impressed source should be placed in the optional source data block for SINDA and VARIABLES 1 block for CINDA-3G.

6.2.2.2 VARIABLES 1 Operations Block

The VARIABLES 1 operations block provides the user with a means of specifying at a point in the computational sequence, as shown in Figure 6.2-1, the evaluation of nonlinear network elements, coefficients and boundary values not considered by the various mnemonic codes utilized for node, conductor and source data. It is seen from Figure 6.2-1 that VARIABLES 1 operations occur just prior to entering the numerical solution phase in order to define the network completely.

6.2.2.3 VARIABLES 2 Operations Block

VARIABLES 2 operations are post-solution operations in contrast to the VARIABLES 1 operations which are pre-solution operations as shown in Figure 6.2-1. VARIABLES 2 provides the user with a means to examine the characteristics of the numerical solution and make corrections. For example, the heat flow from one node to another can be evaluated or a temperature(s) determined without material phase change can be corrected to account for the phase change by using the VARIABLES 2 operations block.

6.2.2.4 OUTPUT CALL Operations Block

The OUTPUT CALL operations block provides the user with a means of calling any desired subroutine with the operation performed on the output interval. In addition to various subroutines for printing output, several plotting subroutines are available.^{3,4}

6.2.3 Control Constants

Control constants number forty-nine and have alphanumeric names. Control constant values are communicated through program common to specific subroutines which require them. Whenever possible, control constant values not specified are set internally to acceptable values. If a required control constant value is not specified, an appropriate error message is printed and the program terminated. Each of the SINDA numerical solution routines employs a number of control constants which fall under the categories as: (1) user specified; (2) optionally user specified; (3) internally set by program; and (4) dummy. These control constants are listed alphabetically with a brief description of each in Section 6.2.3.1 followed by a detailed description of user specified control constants in Section 6.2.3.2; nominal values of these control constants that must be specified or are optionally specified for each SINDA numerical solution routine are indicated in Table 6.2-5. Specification of these control constants is detailed under the discussion of each SINDA numerical solution routine.

6.2.3.1 Alphabetical Listing and Brief Description of Control Constants

ARLXCA (control constant 19)

Maximum arithmetic node relaxation temperature change allowed between iterations; this check occurs after each iteration. Specification is required for the implicit and steady state routines (except CINDSM) and if not specified an error message is printed if the number of arithmetic nodes is greater than zero. Specification is not required for explicit routines and if not specified, ARLXCA is set to 1.E+8.

ARLXCC (control constant 30)

Maximum arithmetic node relaxation temperature change calculated by program; $ARLXCC \leq ARLXCA$ check is made.

ATMPCA (control constant 11)
 Maximum arithmetic node temperature change allowed between time steps for transient routines; the check occurs after the specified number of iterations. If not specified or if specified to be ≤ 0.0 , ATMPCA is set to 1.E+8.

ATMPCC (control constant 15)
 Maximum arithmetic temperature change calculated by program; $ATMPCC \leq ATMPCA$ check is made.

BACKUP (control constant 12)
 Backup switch that is checked after VARIABLES 1 and VARIABLES 2 calls. Initialized at zero. If specified to be non-zero, the completed time step is erased and repeated.

BALENG (control constant 33)
 A user specified system energy balance to be maintained; this control constant is presently used only in CINDSM. If not specified, an error message will be printed.

CSGFAC (control constant 4)
 Time step factor for explicit routines except CNFAST. If not specified or if specified to be less than 1.0, CSGFAC is set internally to 1.0.

CSGMAX (control constant 23)
 Maximum value of $C_i / \sum G_{ij}$; this value aids in the checkout of the thermal network and is calculated only by the output sub-routines, CSGDMP and RCDUMP.

CSGMIN (control constant 17)
 Minimum value of $C_i / \sum G_{ij}$; this value is used to limit the computational time step for explicit methods of solution. If CSGMIN is calculated to be ≤ 0.0 , an error message is printed.

CSGRAL (control constant 24)
 Allowable range between CSGMIN and CSGMAX: this control constant is not presently used but is included for future considerations.

DAMPA (control constant 9)
 Arithmetic-node damping factor for all numerical solution routines; if not specified, or if specified to be ≤ 0.0 , DAMPA is set to 1.0. (Refer to equation 6.2-6.)

DAMPD (control constant 10)
 Diffusion node damping factor for implicit and steady state routines; if not specified or is specified to be ≤ 0.0 , DAMPD is set to 1.0. (Refer to equation 6.2-20.)

DRLXCA (control constant 26)
 Maximum diffusion node relaxation temperature change allowed between iterations for implicit and steady state routines; this check occurs after each iteration. If not specified an error message will be printed when the number of diffusion nodes is greater than zero.

DRLXCC (control constant 27)
 Maximum diffusion node relaxation temperature change calculated by the program; DRLXCA \leq DRLXCC check is made.

DTIMEH (control constant 8)
 Maximum time step allowed; applies to transient routines. If not specified or if specified to be ≤ 0.0 , DTIMEH is set to 1.E-8.

DTIMEI (control constant 22)
 Specified time step for implicit solutions; if not specified, an error message will be printed and the "run" terminated.

DTIMEL (control constant 21)
 Minimum time step allowed for explicit routines. If not specified for CNFAST, an error message will be printed and the "run" terminated. If DTIMEU is less than DTIMEL the routines will

terminate with an error message, except for CNFAST which will do a steady state solution on the offending node. For all routines DTIMEL is initially set at 0.0 internally.

DTIMEU (control constant 2)

Contains time step used in computational procedure.

DTMPCA (control constant 6)

Maximum diffusion node temperature change allowed between time steps for transient routines. If not specified or if specified to be ≤ 0.0 , DTMPCA is set to 1.E+8.

DTMPCC (control constant 15)

Maximum diffusion node temperature change calculated by program; DTMPCA \leq DTMPCC check is made.

ENGBAL (control constant 32)

Calculated energy balance of the system; presently used only in CINDSM.

LAXFAC (control constant 49)

Specified number of iterations to be performed on a linearized system with no updating of elements during a set of LAXFAC iterations for CINDSM only; if not specified, an error message is printed and the "run" terminated.

LINECT (control constant 28)

A line counter location for program output (integer).

LØØPCT (control constant 20)

Contains number of iterations performed (integer).

NØCØPY (control constant 34)

Contains the no copy switch for matrix users.

NLØØP (control constant 5)

Number of specified iteration loops. Must be specified for the steady state and implicit routines; if not specified, an

error message is printed and the "run" is terminated. Optional specification for solution of the arithmetic nodes in the explicit routines; if not specified, $NL\emptyset\emptyset P$ is set to integer 1.

$\emptyset PEITR$ (control constant 7)

Output each iteration if $\emptyset PEITR$ is specified to be non-zero; if not specified, $\emptyset PEITR$ is set at zero. May be switched on and off during a run.

$\emptyset OUTPUT$ (control constant 18)

Time interval for activating $\emptyset OUTPUT$ CALLS of transient routines; if not specified, error message is printed and the "run" terminated. May be addressed by user and modified during a run in VARIABLES 2. Can be used in steady state routines for a series of steady state solutions.

PAGECT (control constant 29)

A page counter location for program output (integer).

TIMEM (control constant 14)

Mean time for a computation interval; $TIMEM = \frac{TIME\emptyset + TIMEN}{2.0}$.

TIMEN (control constant 1)

New time at the end of the computational interval.

$TIMEN = TIME\emptyset + DTIMEU$.

TIMEND (control constant 3)

Problem stop time for transient analysis. Must be $> TIME\emptyset$ for all routines; if not, an error message is printed and "run" terminated. May be addressed by the user and modified during a run.

$TIME\emptyset$ (control constant 13)

Old time at the start of the computational interval. Also used as the problem start time and may be negative; if not specified, $TIME\emptyset$ is set at zero.

I_{TEST}, J_{TEST}, K_{TEST}, L_{TEST}, M_{TEST} (control constants 39, 40, 41, 42 and 43, respectively)

Contain dummy integer constants.

R_{TEST}, S_{TEST}, T_{TEST}, U_{TEST}, V_{TEST} (control constants 44, 45, 46, 47, and 48, respectively)

Contain dummy floating point constants.

(Control constant 31)

Problem type indicator, 0 = THERMAL SPCS, 1 = THERMAL LPCS, 2 = GENERAL.

(Control constant 35)

Contains relative node number of CSGMIN.

(Control constant 36)

Contains relative node number of DTMPCC.

(Control constant 37)

Contains relative node number of ARLXCC.

(Control constant 38)

Contains relative node number of ATMPCC.

6.2.3.2 User Specified and Optionally User Specified Control Constants

The availability of control constants which must be specified or which can optionally be specified provides the user with considerable flexibility to alter the computational criteria and hence the calculated temperatures. On the other hand, this flexibility presents the user with the problem of inputting control constant values if the nominal values are not suitable. An attempt will be made here to provide some guidelines on control constant values based on rather limited data presently available, but it should be recognized that suitable values to be used are dependent on the problem to be solved and often a trade-off must be made between accuracy and computational time. This normally can be obtained only through the use of the numerical solution routines.

ARLXCA (Allowable Arithmetic Node Relaxation Temperature Change)

This control constant must be specified for the implicit routines if any arithmetic node is present and for the steady state routines except CINDSM. For the explicit solution routines, ARLXCA may be optionally specified; if not specified ARLXCA is set to 1.E+8. ARLXCA represents a maximum temperature change convergence criterion for the arithmetic nodes; ARLXCA is checked each iterative step. It is used in conjunction with control constant NLOOP. Satisfaction of either NLOOP or ARLXCA during any iterative step terminates the arithmetic node temperatures calculation for that time-step with computation proceeding on to the next one. Typically, an ARLXCA value is 0.01, but its value is dependent upon the magnitude of expected temperatures. The 0.01 value tries for 5th digit accuracy for temperatures in the hundreds. An ARLXCA value of 0.0001 would try for seventh digit accuracy. Since the computer will not yield 8 digit accuracy, an ARLXCA value < .0001 will always result in NLOOP iterations being performed.

ATMPCA (Allowable Arithmetic Node Temperature Change)

This control constant may be optionally specified by the user for the implicit routines and for the explicit routines except CNFAST. If not specified, ATMPCA is internally set at 1.E+8. ATMPCA represents an allowable arithmetic-node temperature change criterion between one time-step and another with the calculated temperature change stored in control constant ATPCC. If the maximum arithmetic-node temperature change is greater than ATMPCA, the time-step, Δt , is shortened to,

$$\Delta t = .95 * \Delta t \text{ (ATMPCA/ATPCC)}$$

and the arithmetic-node and diffusion-node temperatures re-set to former values. The computational procedure is repeated with the smaller time-step. Specification of ATMPCA prevents a rapid temperature change between time-steps with the value to be specified dependent upon the problem. Thus, the user should estimate the number of time-steps and the range of the temperature to arrive at a reasonable value. For typical spacecraft-type thermal problems an ATMPCA of about 10°F is typical.

BACKUP (Backup Switch)

Control constant BACKUP provides the SINDA user with the means to utilize any thermal numerical solution subroutine as a predictor program. All of the numerical solution subroutines set control constant BACKUP to zero, just prior to the call on VARIABLES 2. Then immediately after the return from VARIABLES 2, a nonzero check on BACKUP is made. If BACKUP is nonzero, all temperature calculations for the just completed time-step are eliminated, the old temperatures (temperatures calculated at the previous time-step) are placed in the temperature locations and the control is routed to the start of the computational sequence.

It should be noted that the user must provide the necessary check and criterion in VARIABLES 2 if the iteration is to be repeated. Thus, if the iteration is to be repeated, BACKUP must be nonzero and a criterion that can be met in subsequent passes established. For example, the criterion may require the correction of a parameter used by the network solution. Further, if other calls in VARIABLES 2 are not to be performed FORTRAN instructions must be generated to bypass these calls.

It should be noted that BACKUP is sometimes checked after VARIABLES 1. However, for the present this use should be ignored since BACKUP check after VARIABLES 1 is planned for future additions of special boundary calculation subroutines.

BALENG (User Specified System Energy Balance)

This control constant is presently used in the steady state routine CINDSM but not in the other SINDA numerical solution routines. BALENG must be specified, otherwise the "run" is terminated with an error message printout; the value of BALENG is a criterion that represents an acceptable net energy balance (energy in minus energy out) of the system in the calculation of steady state temperatures. A value for BALENG depends upon the magnitude of system energy under consideration. As a guideline 1/2% of the total energy into the system (including heat flow from the boundary) is a reasonable value.

CSGFAC (Time Step Factor)

This control constant may be optionally specified by the user

for the explicit routines except CNFAST and it provides the user with some control on the compute time-step as indicated in Section 6.2.4. If CSGFAC is not specified or is specified to be less than one by the user, it is internally set at 1.0. For subroutines CNFRWD and CNFRDL which are conditionally stable CSGFAC is a divisor; a value of CSGFAC greater than one is used to obtain higher accuracy. For subroutines CNEXPN, CNDUFR and CNQUIK, which are unconditionally stable, CSGFAC is a multiplier (refer to page 6-24); a value of CSGFAC greater than one is used to decrease the computational time. A question may be raised, why a value of CSGFAC less than one is not allowed for CNEXPN, CNDUFR and CNQUIK? The reason for this is that it is more accurate to use CNFRWD (or CNFRDL) if a smaller time-step than the one associated with CSGFAC equal to one is desired.

DAMPA (Damping Factor for Arithmetic Nodes)

This control constant may be optionally specified for all of the SINDA numerical solution routines; if not specified, or if specified to be ≤ 0.0 , DAMPA is set to 1.0. In the development of the finite difference expressions as reported in technical literature, little (if any) mention is made about the so-called damping factor. The damping factor does nothing more than to allow a certain fraction ($1.0 - \text{DAMPA}$) of the "old" temperature (temperature at the previous time-step or iteration) to be included as part of the temperature change for the current time-step or iteration. The value to be used is dependent upon the problem and to some extent upon the routine. Typically, a value of 0.6 is used but a value as small as 0.01 has been used with CINDSL for a thermal radiation-dominated problem. In general, a choice for DAMPA becomes a trial and error procedure. DAMPA is used only with arithmetic nodes (refer to equation 6.2-6).

DAMPD (Diffusion Node Damping Factor)

This control constant may be optionally specified for the implicit and steady state routines; if not specified or if specified to be ≤ 0.0 , DAMPD is set to 1.0. DAMPD serves the same purpose for the diffusion nodes as DAMPA provides for the arithmetic nodes (refer to equation 6.2-21).

DRLXCA (Allowable Diffusion-Node Relaxation Temperature Change)

This control constant must be specified for the implicit routines and for the steady state routines except CINDSM. DRLXCA serves the same

purpose for the diffusion-nodes as control constant ARLXCA does for the arithmetic nodes. Thus, the discussion on ARLXCA equally holds true for DRLXCA. It may be asked, why ARLXCA and DRLXCA? The reason for this is that it provides greater computational flexibility.

DTIMEH (Maximum Time-Step Allowed)

This control constant may be optionally specified for the explicit and the implicit routines. DTIMEH represents the maximum time-step allowed during the computational process. One use of DTIMEH is the prevention of a single large and a single small computational time-step during an output interval by specifying DTIMEH as a fraction of the output interval. If DTIMEH is not specified, DTIMEH is set to 1.0E+8.

DTIMEI (Specified Time-Step for Implicit Routines)

This control constant must be specified for the implicit routines and is not used by the other routines. If not specified, the "run" terminates with an error message printout. DTIMEI represents a specified time-step and is arbitrary, but the governing criterion should be minimum computational time with satisfactory temperature accuracy. This means that DTIMEI should be specified in conjunction with control constant NLØØP which represents the maximum number of computational iterations allowed during each time-step. Since each iterative calculation is essentially equivalent to a time-step calculation, DTIMEI should be normally greater than NLØØP*CSGMIN, where CSGMIN is the time-step used in the explicit routines. If savings in computational time cannot be met with the same accuracy by using the implicit routines, it is more reasonable to use the explicit routines.

DTIMEL (Minimum Time-Step Allowed)

This control constant must be specified for subroutine CNFAST and is optional for other explicit solution routines. If not specified for CNFAST, the "run" terminates with an error message printout. DTIMEL represents the minimum time-step allowed; for all the explicit routines except CNFAST, if the calculated time-step is less than DTIMEL, the "run" terminates with an error message printout. For subroutine CNFAST, if the calculated time-step of any node, as expressed by $C_i/\Sigma G_{ij}$ and stored in CSGMIN, is less than DTIMEL, the temperature of the nodes not satisfying DTIMEL are calculated

using the steady state equations without computational iterations (refer to Section 6.3.3 for details on the CNFAST routine). The purpose of this control constant for CNFAST is to shorten the computational time; the danger in its use is that with a large DTIMEL, a large number of diffusion nodes will receive the steady state equations without iterations. As a result, the temperature inaccuracies can be expected to be large.

DTMPCA (Allowable Diffusion Node Temperature Change)

This control constant may be optionally specified by the user for the implicit routines and for the explicit routines except CNFAST. DTMPCA represents a diffusion-node temperature change criterion between one time-step and another. If the maximum diffusion-node temperature change which is stored in DTMPCG is greater than DTMPCA, the time-step is shortened to,

$$\Delta t = .95 * \Delta t \text{ (DTMPCA/DTMPCG)}$$

and the diffusion-node and arithmetic-node temperatures re-set to former values. The computational procedure is repeated with the smaller time-step. DTMPCA serves the same purpose for the diffusion nodes as control constant DRLXCA provides the arithmetic nodes.

LAXFAC (Number of Iterations for Linearized Lumped Parameter System)

LAXFAC is used only in the steady state routine CINDSM and represents the number of iterations to be performed on a linear lumped parameter system with no updating of elements during a set of LAXFAC iterations. The system elements are re-evaluated for the new set of temperatures and in turn temperatures are recalculated for another set of LAXFAC iterations with a more severe relaxation criterion. The number of iterations will not exceed control constant NL00P which represents the total number of iterations. NL00P will not be met only if relaxation criteria are met during an iterative loop and between iterative loops and if the system energy balance as stored in BALENG is satisfied (refer to Section 6.5.3 for details).

NL00P (Number of Iteration Loops)

This control constant must be specified for the implicit and the steady state routines; if not specified, the "run" terminates with an error message printout. NL00P may be optionally specified for the explicit routines since it is used for the arithmetic nodes; if not specified, NL00P is set to 1. The value of NL00P to be used depends upon the problem.

to be solved. For a steady state problem it is not unusual to have NLØØP equal to several hundred, whereas for a transient problem the implicit routines NLØØP should be specified as discussed for control constant DTIMEI. In general, a trial and error procedure is required to arrive at a suitable value of NLØØP.

OUTPUT (Time Interval for Activating ØUTPUT CALLS)

This control constant must be specified for all numerical solution routines except steady state routines since the first time-step used is generally set to ØUTPUT. The input value is left to the judgment of the user. Normally, the output interval is gauged by the length of the run and the expected temperature response characteristics. As a "rule-of-thumb" the output interval lies between CSGMIN and CSGMAX, with ØUTPUT being several times larger than CSGMIN. The values of CSGMIN and CSGMAX can be obtained from the output subroutines CSGDMP and RCDUMP.^{3,4} Subroutines CSGDMP and RCDUMP are designed to aid in the checkout of thermal problem data decks and should be used before making a transient computer run.

TIMEND (Problem Stop Time)

The use of this control constant is self-explanatory. For the subroutines as they are presently coded, TIMEND must be specified as larger than TIMEØ, otherwise an error message is printed and the "run" terminated. For the explicit routines, if TIMEND is not larger than TIMEØ a time-step of zero will result and the "TIME STEP TOO SMALL" error message will be printed. The implicit routines will print the error message, "TRANSIENT TIME NOT SPECIFIED." If a solution is to be terminated by the use of a criteria, but the run is not to be terminated, this can be accommodated by setting TIMEND=TIMEØ when the criteria is met.

TIMEØ ("Old" Time or Problem Start Time)

This control constant represents the "old" time or the problem start time for the transient routines. If not specified, TIMEØ is set to 0.0. An important consideration in the use of TIMEØ is that TIMEØ may be set to negative.

6.2.4 Time-Step Calculations

Each numerical solution routine requires the use of a time-step that depends upon many considerations, such as the output interval, the end

of the problem time, the stability criterion for explicit routines, etc. In spite of the unique solution procedure of each of the numerical solution routines, the overall time-step calculation procedure for the transient routines is essentially identical. The numerous time-step checks, as well as the selection of the time-step, are indicated below (for definition of control constants refer to Section 6.2.3):

- (1) Check that elapsed time, t , does not exceed problem end time.

If: $\text{TIME}\phi + \phi\text{OUTPUT} > \text{TIMEND}$

Set: $\phi\text{OUTPUT} = \text{TIMEND} - \text{TIME}\phi$

$\text{TIME}\phi$ is the old time

ϕOUTPUT is the output time interval

TIMEND is the problem stop time

- (2) Set initial time-step, Δt , which is stored in DTIMEU (control constant for time-step). The initial time step for the SINDA numerical routines is as follows:

<u>Numerical Routines</u>	<u>Initial Time-Step</u>
EXPLICIT CNFRWD	ϕOUTPUT
EXPLICIT CNFRDL	ϕOUTPUT
EXPLICIT CNEXPN	ϕOUTPUT
EXPLICIT CNDUFR	ϕOUTPUT
EXPLICIT CNQUIK	ϕOUTPUT
EXPLICIT CNFAST	DTIMEL (minimum time-step allowed)
IMPLICIT CNBACK	DTIMEI (specified time-step)
IMPLICIT CNFWBK	DTIMEI
IMPLICIT CNVARB	DTIMEI

- (3) Check Δt (stored in DTIMEU) against maximum allowable time-step.

If: $\text{DTIMEU} > \text{DTIMEH}$

Set: $\text{DTIMEU} = \text{DTIMEH}$

- (4) Check sum of elapsed time since last printout, TSUM , and time-step, DTIMEU , against ϕOUTPUT .

If: $\text{TSUM} + \text{DTIMEU} > \phi\text{OUTPUT}$

Set: $\Delta t = \phi\text{OUTPUT} - \text{TSUM}$

If: $\text{TSUM} + \Delta t < \phi\text{OUTPUT}$

and if: $TSUM + 2(\Delta t) > \emptyset OUTPUT$

Set: $\Delta t = 1/2 (OUTPUT - TSUM)$

(5) Store

Set: $DTIMEU = \Delta t$

(6) Check DTIMEU against minimum allowable time-step.

If: $DTIMEU < DTIMEL$

Result: An error message is printed and the "run" terminated
except for CNFAST, CNBACK, CNFWBK and CNVARB.

(7) Set new time (TIMEN)

Set: $TIMEN = TPRINT + TSUM + \Delta t$

$TPRINT$ is the time of the last printout.

$TSUM$ is the time from the last printout.

(8) Set mean time (TIMEM)

Set: $TIMEM = 1/2 (TIMEN + TIME\emptyset)$

(9) Calculate (or specify) time-step.

The calculated (or specified) time-step for the SINDA numerical routines is as follows:

<u>Numerical Routines</u>	<u>Calculated Time-Step</u>
EXPLICIT CNFRWD	$0.95 * CSGMIN/CSGFAC$
EXPLICIT CNFRDL	$0.95 * CSGMIN/CSGFAC$
EXPLICIT CNEXPN	$0.95 * CSGMIN * CSGFAC$
EXPLICIT CNDUFR	$0.95 * CSGMIN * CSGFAC$
EXPLICIT CNQUIK	$0.95 * CSGMIN * CSGFAC$
EXPLICIT CNFAST	larger of CSGMIN or DTIMEL
IMPLICIT CNBACK	DTIMEI
IMPLICIT CNFWBK	DTIMEI
IMPLICIT CNVARB	DTIMEI

$CSGMIN = C_i / \sum G_{ij}$ (minimum value, $i = 1, 2, \dots, NND$)

where: C_i is the capacitance of the i th node

G_{ij} is the conductance from node i to node j

$CSGFAC$ is the time-step factor (see above).

- (10) It should be recognized that individual routines may have slight variations to the time-step calculations.

6.2.5 Computation of Temperatures

The actual calculation of temperatures, be it for diffusion nodes or for arithmetic nodes, represents the end result of a long computational procedure with many checks and criteria. Nevertheless, if one confines the discussion to the $D\emptyset$ loops of nodal types, a rather compact but general computational pattern becomes apparent. More details are presented in the individual sections describing each numerical solution routine. (Sections 6.3 - 6.5)

6.2.5.1 Transient Explicit Routines

For the explicit routines the diffusion and arithmetic nodes are treated separately. Diffusion-node temperatures are calculated explicitly, whereas the arithmetic-node temperatures are computed implicitly. This means that at each time-step an iterative loop is set-up for the arithmetic nodes; none is required for the diffusion nodes.

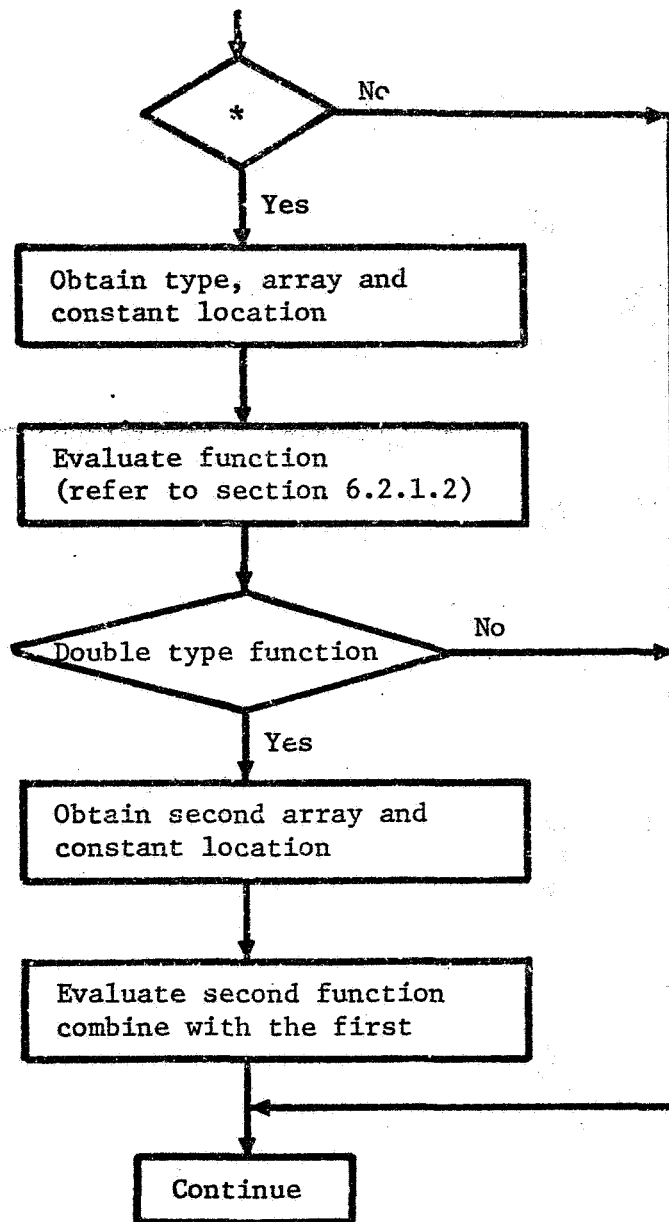
Diffusion-Node Temperatures

Calculation of the diffusion-node temperatures follows the VARIABLES 1 call; the computational pattern is:

$D\emptyset-L\emptyset\emptyset P$ ($I = 1, NND$) on the diffusion nodes is established.

The functions associated with the variable capacitance C_i , the variable impressed source q_i , and the variable coefficients G_k (a_{ij} for conduction and σ_{ij} for radiation), between diffusion-diffusion and diffusion-arithmetic nodes are updated at the beginning of each time-step. These functional types are described in Section 6.2.1.2 and the computational pattern is indicated in the flow chart of Figure 6.2-3.

Using the updated C_i , q_i and G_k , the branch heat flow sum, Q_{si} , and conductance sum X_i , are calculated (refer for example to flow chart of Figure 6.3-1).



* Variable capacitance (c_i),
impressed source (q_i), or
variable coefficient (G_k).

Figure 6.2-3. Evaluation of Nonlinear Capacitance,
Source or Conductance

$$Q_{si} = \sum_{j=1}^P G_{ij,n} (T_{j,n} - T_{i,n}) + q_{i,n} \quad (6.2-1)$$

$$X_i = \sum_{j=1}^P G_{ij,n} \quad (6.2-2)$$

where, p = total number of nodes; n = time-step old
 C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Table 6.2-1 - 6.2-4)

$$G_{ij,n} = a_{ij,n} + ob_{ij,n} (T_{j,n}^2 + T_{i,n}^2)(T_{j,n} + T_{i,n})$$

Stability criterion $C_i / \sum_{j=1}^P G_{ij,n}$ is computed and the smallest value is stored in control constant CSGMIN. If $CSGMIN \leq 0.0$, an error message is printed and the "run" terminated.

Diffusion-node temperatures are calculated by using the appropriate finite difference expression associated with each routine. These routines and algorithms are identified as:

CNFRWD and CNFRDL (Section 6.3.1), uses standard forward-difference algorithm.

CNFAST (Section 6.3.2), uses a modified CNFRWD computational procedure to decrease the computational time.

CNEXPN (Section 6.3.3), uses the exponential prediction method.

CNDUFR (section 6.3.4), uses DuFort-Frankel method.

CNQUIK (Section 6.3.5), uses half DuFort-Frankel and half exponential prediction method.

Symbolically, the expression for the diffusion-node temperatures may be written as,

$$T_{i,n+1} = T_{i,n} + \frac{\Delta t Q_{si}}{C_i} \quad (6.2-3)$$

Except for CNFAST the maximum diffusion-node temperature change which is stored in DTMPCC is checked against the allowable diffusion node temperature change which may be specified by the user via the control constant DTMPCA (if not specified $DTMPCA = 1.0E+8$). If DTMPCA is not satisfied, the time-step is decreased to,

$$\Delta t = .95 * \Delta t (DTMPCA/DTMPCC)$$

and all temperatures re-set to former values. The computational procedure is repeated with the smaller time-step. CNFAST does not allow for the recalculation of diffusion-node temperatures.

Arithmetic-Node Temperatures

Calculation of the arithmetic-node temperatures always follows the computation of the diffusion-node temperatures and uses "successive point" iteration. The computational pattern is as follows:

Arithmetic-node damping factors DN and DD are established.

DN = DAMPA (optionally specified user constant, if not specified DAMPA = 1.0; factor for the current time-step temperature change)

DD = 1.0 - DN (factor that allows a certain fraction of the "old" temperature to be included as part of the temperature change for the current time-step)

Iterative D ϕ -L $\phi\phi$ P (K=1, NL $\phi\phi$ P) is established (NL $\phi\phi$ P is the number of iterations specified by the user, if not specified, NL $\phi\phi$ P = 1).

D ϕ -L $\phi\phi$ P (I=NND, NND + NNA) for the arithmetic nodes is established.

Impressed source q_i and coefficient G_k (a_{ij} for conduction and ob_{ij} for radiation) are updated once for each time-step.

Using the updated G_k and q_i , the branch heat flow sum Q_{si} and the conductance sum X_i are calculated (refer to flow chart of Figure 6.3-2).

$$Q_{si} = \sum_{j=1}^P G_{ij,n} (T_{j,k} - T_{i,k}) \quad (6.2-4)$$

$$X_i = \sum_{j=1}^P G_{ij,n} \quad (6.2-5)$$

Arithmetic node temperatures are calculated for each iterative loop by using the following "successive point" expression, which is employed in all of the routines,

$$T_{i,k+1} = DD * T_{i,k} + DN * \left(\frac{q_{i,n} + \sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^p G_{ij,n} T_{j,k}}{\sum_{j=1}^p G_{ij,n}} \right) \quad (6.2-6)$$

where, $i = (NND+1), (NND+2), \dots, (NND + NNA)$

$T_{j,k} = \text{constant}, (NND + NNA) < j \leq p$

$p = \text{total number of nodes}$

$T_{i,k} = \text{temperature at } k\text{th iteration}$

$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$

$(\ell = k \text{ if } j \geq i \text{ and } \ell = k+1 \text{ if } j < i)$

$(a_{ij,n} \text{ and } b_{ij,n} \text{ mean updating at time-step, } n)$

$q_i, a_{ij}, b_{ij} = \text{optionally specified (refer to Tables 6.2-1 - 6.2-4)}$

$DN \equiv \text{DAMPA (arithmetic node damping factor)}$

$DD = 1.0 - DN$

The maximum arithmetic-node relaxation temperature change is calculated and checked against the allowable arithmetic-node relaxation temperature change which may be specified via the control constant ARLXCA. This relaxation convergence check is made during each iterative step calculation and is used in conjunction with control constant NLØØP. Satisfaction of either ARLXCA or NLØØP during any iterative step terminates the arithmetic-node temperature calculation.

For each time step, except for CNFAST, the maximum arithmetic-node temperature change which is stored in control constant ATMPCC is checked against the allowable arithmetic-node temperature change which may be specified via the control constant ATMPCA (if not specified, ATMPCA = 1.0E+8). If ATMPCA is not satisfied, the time-step is decreased to,

$$\Delta t = .95 * \Delta t (ATMPCA/ATMPCC)$$

and all temperatures re-set to former values. The computational procedure is repeated with the smaller time-step. CNFAST does not allow for recalculation of arithmetic-node temperatures.

6.2.5.2 Transient Implicit Routines

Both diffusion-node and arithmetic-node temperatures are calculated by "successive point" iteration. Although these calculations are performed

on the same iterative pass, diffusion node temperatures are evaluated on its own computational loop using a specified algorithm associated with a particular implicit routine. Calculation of the arithmetic-node temperatures is also done on its own computational loop and is identical in all the implicit routines. As a matter of fact, arithmetic-node temperatures are calculated in the same manner in all the SINDA numerical solution routines. Use of a separate computational loop for the diffusion nodes permits the extrapolation of diffusion-node temperatures provided acceleration of convergence criterion is met (refer to Section 6.2.7).

Diffusion-Node Temperatures

In order to facilitate the discussion to follow on the computational procedure, it is convenient to examine the forward-backward finite difference expression.¹³

$$C_i \frac{(T_{i,k+1} - T_{i,n})}{\Delta t} = \beta T_{\text{forward}} + (1 - \beta) T_{\text{backward}} \quad (6.2-7)$$

where: $\beta =$ factor with range $0 \leq \beta \leq 1/2$

$$T_{\text{forward}} = q_{i,n} + \sum_{j=1}^P a_{ij,n} (T_{j,n} - T_{i,n}) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \quad (6.2-8)$$

$$T_{\text{backward}} = q_{i,n} + \sum_{j=1}^P a_{ij,n} (T_{j,k+1} - T_{i,k+1}) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,k+1}^4 - T_{i,k+1}^4) \quad (6.2-9)$$

$$i = 1, 2, \dots, N$$

$$T_{j,n}; T_{j,k+1} = \text{constant}, N < j \leq P$$

$n =$ nth time-step; $k =$ kth iteration within a given time-step.

$C_i, q_i, a_{ij}, b_{ij} =$ optionally specified (refer to Tables 6.2-1 -- 6.2-4)

Any value of β less than one yields an implicit set of equations which must be solved simultaneously. For values of β less than or equal to one-half equation (6.2-7) represents an unconditionally stable set of equations, whereas values of β greater than one-half yields a set of equations with conditional stability.

The standard implicit algorithm used in subroutine CNBACK follows directly from equation (6.2-7) by letting $\beta = 0$, whereas the Crank-Nicolson method used in subroutine CNFWBK follows by letting $\beta = 1/2$. Subroutine

CNVARB uses a variable factor which is based upon the ratio of CSGMIN/DTIMEU; this ratio is internally calculated in CNVARB (refer to Section 6.4.3.2). In order to simplify the presentation, the following notation is used.

For CNBACK ($\beta = 0$):

$$Q_i = q_{i,n} + \bar{C}_{i,n} T_{i,n} \quad (6.2-10)$$

$$Q_{\text{sum}} = Q_i + \sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,n} T_{j,k} \quad (6.2-11)$$

$$G_{\text{sum}} = \bar{C}_{i,n} + \sum_{j=1}^P a_{ij,n} \quad (6.2-12)$$

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} T_{j,\ell}^3 \quad (6.2-13)$$

($\ell = k$, if $j \geq i$ and $\ell = k+1$, if $j < i$)

$$(q_i)_{\text{ave}} = \frac{1}{2} \sum_{j=1}^P \sigma b_{ij,n} [(T_{i,k}^4) + (T_{i,k}^4)_2], \text{ average heat loss} \quad (6.2-14)$$

from i th node, called radiation damping (refer to Section 6.2.6 for details)

= 0, if radiation is not present

For CNFWBK ($\beta = \frac{1}{2}$) (note equation (6.2-7) is multiplied by 2):

$$Q_i = 2q_{i,n} + 2\bar{C}_{i,n} T_{i,n} + \sum_{j=1}^P a_{ij,n} (T_{j,n} - T_{i,n}) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \quad (6.2-15)$$

$$Q_{\text{sum}} = \text{same as equation (6.2-11)}$$

$$G_{\text{sum}} = 2\bar{C}_{i,n} + \sum_{j=1}^P a_{ij,n} \quad (6.2-16)$$

$$G_{ij,n} = \text{same as equation (6.2-13)}$$

$$(q_i)_{\text{ave}} = \text{same as equation (6.2-14)}$$

For CNVARB (variable β') (note that equation (6.2-7) is multiplied by 2, so that $\beta' = 2\beta$ now ranges, $0 \leq \beta' \leq 1.0$):

$$Q_i = 2q_{i,n} + 2\bar{C}_{i,n} T_{i,n} + \beta' \left(\sum_{j=1}^P a_{ij,n} (T_{j,n} - T_{i,n}) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \right) \quad (6.2-17)$$

$$Q_{sum} = Q_i + (2.0 - \beta') \left(\sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,n} T_{j,k} \right) \quad (6.2-18)$$

$$G_{sum} = 2\bar{C}_{i,n} + (2.0 - \beta') \sum_{j=1}^P a_{ij,n} \quad (6.2-19)$$

$G_{ij,n}$ = same as equation (6.2-13)

$$(q_i)_{ave} = \frac{2.0 - \beta'}{2} \sum_{j=1}^P \sigma b_{ij,n} [(T_{i,k}^4) + (T_{i,k}^4)_2], \text{ average heat} \quad (6.2-20)$$

loss from ith node, called radiation damping (refer to Section 6.2.6 for details)

= 0, if radiation is not present

$i = 1, 2, \dots, N$

$\beta = 2.0 * CSGMIN / DTIMEU$ (range allowed, $0 \leq \beta' \leq 1.0$)

$T_{j,n}; T_{j,k} = \text{constant}, N < j \leq p$ (p is the total number of nodes)

n = nth time-step; k = kth iteration

C_i, q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

$$\bar{C}_{i,n} = C_{i,n} / \Delta t$$

Calculation of the diffusion-node temperatures follows VARIABLES 1 call; the computational pattern is:

Iterative DO-LOOP (kl=1, NLOOP) for the total nodal system is established.

First Iterative Loop:

DO-LOOP (I=1, NND) on diffusion nodes is established.

The functions associated with the variable capacitance C_i , the variable impressed source q_i , and the variable coefficients G_k (a_{ij} for conduction and σb_{ij} for radiation) between diffusion-diffusion and diffusion-arithmetic nodes are updated once for each time-step. These functional types are described in Section 6.2.1.2 and the computational pattern is indicated in the flow chart of Figure 6.2-3.

All known quantities (those evaluated at time-step n) are summed and are identified by the symbol Q_i (equations 6.2-10, 6.2-15 and 6.2-17). CSGMIN is evaluated.

Radiation damping is used; average radiation heat loss, $(q_i)_{ave}$, from the ith node is evaluated (refer to Section 6.2.6).

For CNVARB, $\beta' = 2.0 * CSGMIN / DTIMEU$ is calculated.

The diffusion-node temperatures are calculated by "successive point" iteration (actually CNBACK and CNFWBK have slightly different first iterative pattern than CNVARB but the difference is not significant).

$$T_{i,k+1} = DD * T_{i,k} + DN * [Q_{sum} - (q_i)_{ave}] / G_{sum} \quad (6.2-21)$$

DN = DAMPD (use specified diffusion node damping factor, if not specified, DAMPD = 1.0)

DD = 1.0 - DN

For CNVARB, the diffusion-node relaxation temperature change is calculated; maximum value is stored in DRLXCC.

Second and Succeeding Iterative Loops:

With the iterative loops after the first, those quantities C_i , q_i , and G_k which were updated during the first iteration are held constant.

Diffusion-node temperatures are found by using equation (6.2-21).

The diffusion-node relaxation temperature change is calculated and the maximum value stored in DRLXCC.

Check of DRLXCC against DRLXCA (allowable maximum diffusion-node relaxation temperature change) is made after the arithmetic-node temperature calculations.

Each third iteration, a check on solution convergence is made; if convergence is occurring linear extrapolation to accelerate convergence is made (refer to Section 6.2.7).

Arithmetic-Node Temperatures (if any)

During the first iterative loop the impressed source q_i and coefficient G_k (a_{ij} for conduction and σb_{ij} for radiation) between arithmetic-arithmetic nodes are updated once each time-step. On every loop, arithmetic-

node temperatures are calculated using "successive point" iteration. The finite difference algorithm is presented in Section 6.2.5.1 (equation 6.2-6).

The arithmetic-node relaxation temperature change is calculated and the maximum is stored in ARLXCC.

During Each Iterative Loop After the First

Both DRLXCC and ARLXCC are checked against DRLXCA and ARLXCA, respectively. If both DRLXCA and ARLXCA are satisfied, the iteration ceases.

If LØØPCT equals NLØØP the message "RELAXATION CRITERIA NOT MET" is printed.

Both the calculated maximum diffusion-node and arithmetic-node temperature change (stored in DTMPCC and ATMPCC, respectively) are checked against the corresponding allowable temperature change stored in DTMPCA and ATMPCA. If DTMPCA is not satisfied, the time-step is decreased to,

$$\Delta t = .95 * \Delta t (DTMPCA / DTMPCC)$$

and all temperatures re-set to former values. The computational procedure is repeated with the smaller time-step.

If ATMPCA is not satisfied, the time-step is decreased to,

$$\Delta t = .95 * \Delta t (ATMPCA / ATMPCC)$$

and all temperatures re-set to former values. The computational procedure is repeated with the smaller time-step.

6.2.5.3 Steady State Routines

Diffusion nodes and arithmetic nodes are treated separately in CINDSS and CINDSL even though from a physical standpoint a distinction between diffusion nodes (nodes with capacitance) and arithmetic nodes (nodes with no capacitance) doesn't exist. Thus, the set of control constants for the diffusion nodes and another set of control constants for arithmetic nodes are similar to those used in the transient routines. No distinction in the type of nodes is made in CINDSM.

The computational procedure to be discussed applies only to CINDSS and CINDSL: CINDSM is considerably different (refer to Section 6.5.3).

Diffusion-Node Temperatures (nodes specified with capacitance even though the problem is steady state)

An iterative DØ-LØØP (K1=1,NLØØP) is established.

Within this iterative loop a DØ-LØØP (I=1,NND) on the diffusion nodes is made. The functions associated with the impressed source q_i and the variable coefficients G_k (a_{ij} for conduction and σ_{ij} for radiation) between diffusion-diffusion and diffusion-arithmetic nodes are updated each iteration.

Diffusion-node temperatures are calculated using "block" iteration for CINDSS and "successive point" iteration for CINDSL.

"Block" iteration (CINDSS):

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{\left(q_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k} \right)}{\sum_{j=1}^P G_{ij,k}} \quad (6.2-22)$$

$$G_{ij,k} = a_{ij,k} + \sigma_{ij,k} (T_{j,k}^2 + T_{i,k}^2) (T_{j,k} + T_{i,k})$$

DN = DAMPD (diffusion-node damping factor)

DD = 1.0 - DN

i = 1,2,...,NND (number of diffusion nodes)

k = kth iteration; p = total number of nodes

q_i, a_{ij}, b_{ij} = optionally specified to Tables (6.2-1 - 6.2-4)

$T_{j,k}$ = constant, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes)

"Successive point" iteration (CINDSL):

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{\left(q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k} \right)}{\sum_{j=1}^P G_{ij,k}} \quad (6.2-23)$$

$$G_{ij,k} = a_{ij,k} + \sigma_{ij,k} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$$

($\ell = k$ if $j \geq i$ and $\ell = k+1$ if $j < i$)

DN = DAMPD

DD = 1.0 - DN

$i = 1, 2, \dots, (NND + NNA)$
 $k = \text{kth iteration}; p = \text{total number of nodes}$
 $q_i, a_{ij}, b_{ij} = \text{optionally specified to Tables (6.2-1 - 6.2-4)}$
 $T_{j,k} = \text{constant, } (NND + NNA) < j \leq p \text{ (NNA is the total number of arithmetic nodes)}$

Diffusion-node relaxation temperature change is calculated and the maximum is stored in DRLXCC.

Arithmetic-Node Temperatures (nodes specified with no capacitance)

Within this iterative D ϕ -L $\phi\phi$ P a D ϕ -L $\phi\phi$ P (I=NND+1, NND + NNA) is established.

The functions associated with impressed source q_i and variable coefficients G_k (a_{ij} for conduction and b_{ij} for radiation) between arithmetic-arithmetic nodes are updated each iteration.

Arithmetic-node temperatures are calculated using "successive point" iteration.

$$T_{i,k+1} = AD * T_{i,k} + AN * \frac{\left(q_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k} \right)}{\sum_{j=1}^P G_{ij,k}} \quad (6.2-24)$$

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,\ell})$$

$$(\ell = k \text{ if } j \geq i \text{ and } \ell = k+1 \text{ if } j < i)$$

AN = DAMPA (arithmetic-node damping factor)

AD = 1.0 - AN

$i = (NND+1), (NND+2), \dots, (NND + NNA)$ (number of arithmetic nodes)

$k = \text{kth iteration}$

$p = \text{total number of nodes}$

$T_{j,k} = \text{constant, } (NND + NNA) < j \leq p$

The arithmetic-node relaxation temperature change is calculated and the maximum value is stored in ARLXCC.

During Each Iterative Loop

Both DRLXCC and ARLXCC are checked against DRLXCA and ARLXCA, respectively. If both relaxation criteria, DRLXCA and ARLXCA, are satisfied, the iteration ceases.

If both relaxation criteria, DRLXCA and ARLXCA, are not met with NLØØP iterations, the message "ITERATION COUNT EXCEEDED, NLØØP = " is printed.

Energy balance of the system is calculated and is stored in control constant ENGBAL.

6.2.6 Radiation Damping

Radiation damping denotes an averaging of radiation heat loss technique used to prevent or minimize large temperature oscillations. This method is currently employed in only the implicit routines. The technique which is original with J. D. Gaski is based upon practical and computational considerations. Solution of numerous problems without large temperature oscillations indicates the effectiveness of the approach.

The radiation averaging technique is relatively simple conceptually and rather easily incorporated in the numerical solution routines. The computational pattern is such that the diffusion nodes are encountered sequentially. Let the encountered node be the i th node. A check is made for the presence of a radiation coefficient, $G_k = \sigma b_{ij}$, to the i th node. If one or more radiation connections is present, the radiation heat loss, $(q_i)_{rl}$, from the i th node is calculated based upon the previous temperature $T_{i,k}$.

$$(q_i)_{rl} = \sum_j \sigma b_{ij,n} T_{i,k}^4 \quad (6.2-24)$$

where, j = all radiation connections to node i
 n = n th time-step
 k = k th iteration

Using $(q_i)_{rl}$, a second temperature $(T_{i,k})_2$, is found as follows:

$$(T_{i,k})_2 = [Q_{sum} - (q_i)_{rl}] / G_{sum} \quad (6.2-25)$$

where,

$$Q_{\text{sum}} = \bar{C}_i T_{i,n} + q_{i,n} + \sum_{j=1}^i a_{ij,n} T_{j,k+1} + \sum_{j=i+1}^p a_{ij,n} T_{j,k} \\ + \sum_{j=1}^i \sigma b_{ij,n} T_{j,k+1}^4 + \sum_{j=i+1}^p \sigma b_{ij,n} T_{j,k}^4 \quad (6.2-26)$$

$$G_{\text{sum}} = \bar{C}_i + \sum_{j=1}^p a_{ij,n} \quad (6.2-27)$$

Note that in the evaluation of $(T_{i,k})_2$, the damping factor DAMPD is not used. Note further that G_{sum} does not contain $\sum_j \sigma b_{ij,n} T_{i,k}^3$ since it is accounted for in the radiation loss term, $(q_i)_{r1}$.

Now a second radiation heat loss based on $(T_{i,k})_2$ is found,

$$(q_i)_{r2} = \sum_j \sigma b_{ij,n} (T_{i,k})_2^4 \quad (6.2-28)$$

Equations (6.2-24) and (6.2-28) are then averaged,

$$(q_i)_{\text{ave}} = [(q_i)_{r1} + (q_i)_{r2}] / 2.0 \quad (6.2-29)$$

This average radiation heat loss from an ith node is used in the diffusion-node finite difference algorithm as follows,

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{(Q_{\text{sum}} - (q_i)_{\text{ave}})}{G_{\text{sum}}} \quad (6.2-30)$$

where,

$$DN = \text{DAMPD}$$

$$DD = 1.0 - DN$$

$$(q_i)_{\text{ave}} = \text{average radiation heat loss (equation 6.2-29)}$$

$$G_{\text{sum}} = \bar{C}_i + \sum_j a_{ij,n}$$

Q_{sum} = of the form shown by equation (6.2-26). The actual expression depends upon algorithm. Equation (6.2-26) is for the standard implicit method.

The reason behind the use of $(q_i)_{\text{ave}}$ is that if the initial temperature $T_{i,k}$ is too large, the heat loss from the ith node, $(q_i)_{r1}$ would then be too large. As a result the evaluation of $(T_{i,k})_2$ with $(q_i)_{r1}$ would yield a temperature that is too low. Thus, the averaging of $(q_i)_{r1}$ and $(q_i)_{r2}$ would be much closer to the true heat loss from the

ith node. If $T_{i,k}$ is too small then $(T_{i,k})_2$ would be too large; the averaging scheme still holds true.

6.2.7 Acceleration of Convergence by Extrapolation Technique

Several of the SINDA numerical solution routines use an extrapolation technique to accelerate convergence of the iterative procedure. The extrapolation technique is used in the implicit routines CNBACK, CNFWBK, and CNVARB for the iterative temperature solution of the diffusion nodes, but is not used for the iterative temperature solutions of the arithmetic nodes. The extrapolation method is also used in the steady state routines CINDSL and CINDSM for the iterative temperature solution of both the diffusion and the arithmetic nodes.

6.2.7.1 Extrapolation Technique

The extrapolation is based on a zero temperature difference condition which is defined to be a point where the temperature change of a particular node over two successive iterations is zero. The governing equations are developed as follows:

Consider the temperatures of an ith node at three successive iterations as shown in Figure 6.2-4a. Let these temperatures, which are assumed to be successively decreasing (or increasing), be denoted as,

$$T_{i,k-2}, T_{i,k-1} \text{ and } T_{i,k}$$

where, k is the present iteration
 $k-1$ is the previous iteration
 $k-2$ is two iterations before the k th iteration

By taking the differences,

$$\Delta T_{i,k-1} = T_{i,k-2} - T_{i,k-1}$$

$$\Delta T_{i,k} = T_{i,k-1} - T_{i,k}$$

and plotting these temperature differences as a function of iterations, the iterative point of zero temperature difference can be found by linear extrapolation as shown in Figure 6.2-4b. The corresponding expression for the line is found by using the point, $\Delta T_{i,k}$ at $I = k$ and the slope, $(\Delta T_{i,k} - \Delta T_{i,k-1}) / (k - (k-1))$, to yield,

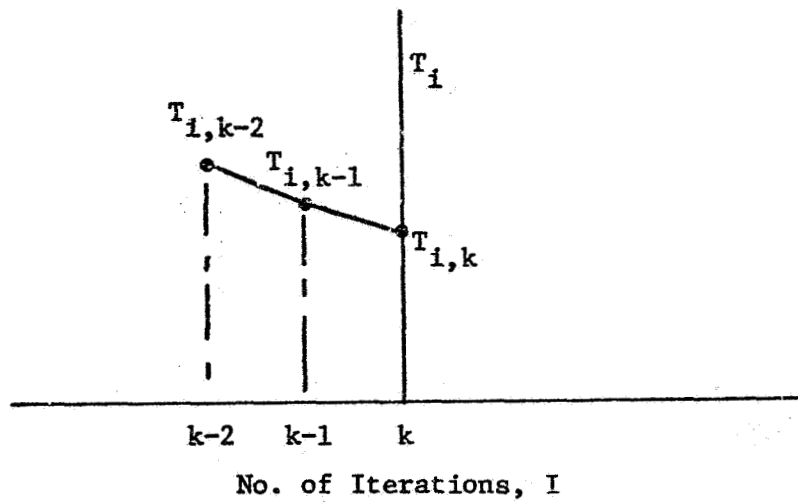


Figure 6.2-4a. Temperature (ith) vs. No. of Iterations

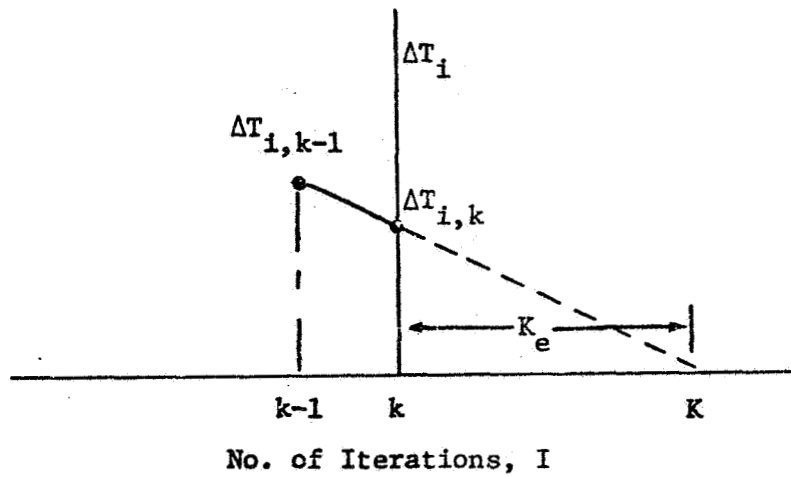


Figure 6.2-4b. Temperature Difference vs. No. of Iterations

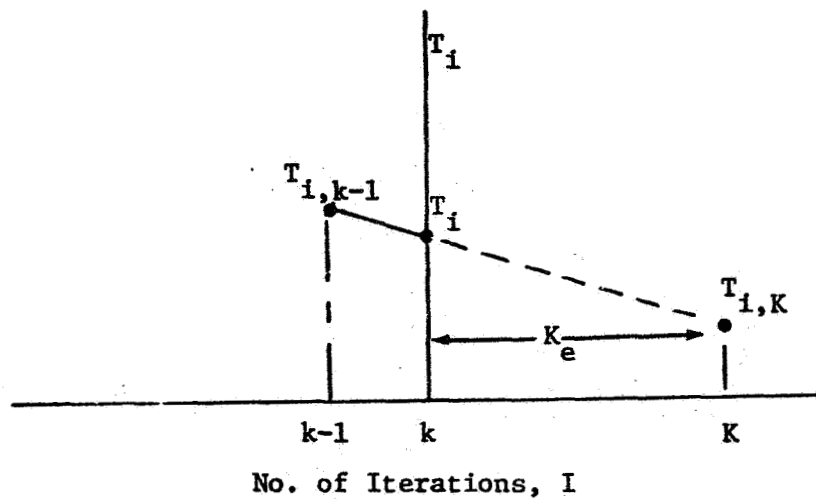


Figure 6.2-4c. Extrapolation of Temperature (ith) to New Value

Figure 6.2-4. Method of Extrapolation to Accelerate Convergence

$$\Delta T_{i,I} = \Delta T_{i,k} + (\Delta T_{i,k} - \Delta T_{i,k-1})(I - k) \quad (6.2-31)$$

where, $I = \text{iterations}$

Since at the zero temperature difference condition, $\Delta T_{i,I} = 0$, the expression for the extrapolated iterations, $K_e = (K - k)$, is found to be,

$$K_e = - \Delta T_{i,k} / (\Delta T_{i,k} - \Delta T_{i,k-1}) \quad (6.2-32)$$

Now, by extrapolating the line established by the temperatures, $T_{i,k-1}$ and $T_{i,k}$ to the line $I = K$, as shown in Figure 6.2-4c, the extrapolated temperature $T_{i,K}$ is found. The expression is readily found to be,

$$T_{i,I} = T_{i,k} + (T_{i,k} - T_{i,k-1})(I - k) \quad (6.2-33)$$

Since $I = K$ and $K - k = K_e$, equation (6.2-33) becomes,

$$T_{i,K} = T_{i,k} (1 + K_e) - K_e T_{i,k-1} \quad (6.2-34)$$

6.2.7.2 Programming Considerations

Each applicable node is tested at the completion of each third iteration to determine if the extrapolation method should be applied. If K_e is calculated to be less than or equal to zero, extrapolation is neglected since the error function is diverging. If K_e is calculated to be greater than zero, a new temperature is calculated based on equation (6.2-34); however, to avoid problems associated with a nearly-zero slope of the line representing the temperature difference vs. number of iterations relationship (Figure 6.2-4b), K_e is set to a number K_m ; otherwise, K_e could be a very large number. For the implicit routines, CNBACK, CNFWBK, and CNVARB, $K_m = 10$. For the steady state routine CINDSL $K_m = 8$ and for steady state routine CINDSM a criterion based upon the maximum temperature is used.

6.2.7.3 Routines Using Acceleration of Convergence

SINDA numerical solution routines that employ the acceleration of convergence features are:

CINDSL, CINDSM	Steady state routines
CNBACK, CNFWBK, CNVARB	Transient implicit routines

6.2.7.4 Comment on Acceleration of Convergence

Neither an extensive study on the value of the acceleration convergence feature has been made, nor has one been reported, but the

limited results presently available indicate that for the steady state routine CINDSL, the number of iterations is reduced approximately 20%. Results are not available for the implicit routines.

A study of the acceleration of convergence feature is made difficult because the method is not a user option in the applicable SINDA numerical solution routines. Thus, the user must be sufficiently versed with the routines in order to delete the acceleration of convergence feature.

6.2.8 Other Characteristics of the SINDA Numerical Solution Routines

6.2.8.1 Units

SINDA, as presently coded, requires that the temperatures must be specified in degrees Fahrenheit ($^{\circ}\text{F}$) since the conversion factor to obtain degrees absolute is internally set at 460.0. This means that the units must be consistent with $^{\circ}\text{F}$ (or $^{\circ}\text{R}$). The execution routines as presently coded do not permit the use of other units.

6.2.8.2 General Comments on Computational Features

Many of the computational features such as radiation damping are original with J. D. Gaski. No theoretical proofs are offered since a practical "gut-feel" development was often used in lieu of a sophisticated mathematical approach; the features, in general, appear to meet the intended objectives. It should be particularly noted that the numerical solution routines are computationally similar; within a particular numerical solution class explicit, implicit or steady state, the computational similarity is even more pronounced. Yet on the other hand, similarity of patterns are broken for no particular reason other than the programmer's whim.

6.3 Transient Explicit Solution Routines

SINDA explicit solution routine number six. These are identified as follows:

- CNFRWD Conditionally stable explicit forward difference.
Requires short pseudo-compute sequence (SPCS).
- CNFRDL Identical to CNFRWD except that the long pseudo-compute sequence (LPCS) is required.
- CNFAST Modified CNFRWD for accelerated forward differencing.
Requires short pseudo-compute sequence (SPCS).
- CNEXPN Unconditionally stable explicit differencing using exponential prediction.
Requires short pseudo-compute sequence (SPCS).
- CNDUFR Unconditionally stable explicit differencing using DuFort-Frankel method.
Requires short pseudo-compute sequence (SPCS).
- CNQUIK Unconditionally stable explicit differencing using a combination of half CNEXPN and half CNDUFR.
Requires short pseudo-compute sequence (SPCS).

A detailed description of each explicit routine is presented on the pages to follow with heavy reliance upon the general description of Section 6.2. A brief description of these routines is summarized first.

CNFRWD uses an explicit forward differencing algorithm and requires the short pseudo-compute sequence (SPCS). The explicit method is characterized by computational simplicity and stability limitations. Since the allowable time-step is governed by the smallest time constant of the network, care must be given in reducing the physical system to a reasonable lumped-parameter model. Arithmetic-node temperatures are calculated by "successive point" iteration.

CNFRDL is identical to CNFRWD except that CNFRDL requires the long pseudo-compute sequence instead of the short pseudo compute sequence. CNFRDL requires slightly less solution time than CNFRWD but the difference is not significant; CNFRDL does require more core storage, however.

CNFAST represents a modified CNFRWD with the modifications intended to decrease the computational time. A user specified control constant DTIMEL which contains the minimum time-step allowed is used as a criterion for isolating those diffusion nodes that are to receive the steady state calculations. A large pocket of internally converted diffusion nodes can present considerable accuracy problems.

CNEXPN uses an unconditionally stable explicit method with the intent to reduce computational time at the expense of temperature accuracy. If accuracy is an important consideration, another routine such as CNFRWD would be a better choice. As a note of interest, CNEXPN solutions tend to lag in time the true solutions.

CNDUFR uses the unconditionally stable DuFort-Frankel method with the intent to reduce computational time by using time-steps greater than those allowed with the conditionally stable explicit methods. Again accuracy may be compromised. CNDUFR solutions tend to lead in time the true solutions.

CNQUIK uses half CNEXPN and half CNDUFR. Why? Since CNEXPN solutions tend to lag in time and CNDUFR solutions tend to lead in time, a combination may yield better solutions. Preliminary results indicate that CNQUIK solutions are more accurate than either CNEXPN or CNDUFR for the same computational time.

6.3.1 Subroutines: CNFRWD and CNFRDL

6.3.1.1 General Comments

Subroutines CNFRWD and CNFRDL are numerical solution routines that use the forward finite difference explicit approximation^{12, 14} of the parabolic differential equation. CNFRWD and CNFRDL are identical except that CNFRDL requires the short pseudo compute sequence (SPCS) whereas CNFRDL requires the long pseudo compute sequence (LPCS). The need for both routines becomes apparent when it is understood that if a steady state numerical solution routine is followed by a transient numerical solution routine, both routines must have consistent PCS (LPCS or SPCS). As a note of interest, each arithmetic node receives the long pseudo compute sequence (LPCS) but this is done internally by the program.

The forward finite difference explicit method as used in CNFRWD and CNFRDL is the conventional Euler method that neither provides a check on the accuracy nor does it provide any scheme of correction once the temperature values are calculated except for the arithmetic nodes which are reiterated $NL\emptyset\emptyset P$ -times. The explicit method is characterized by computational simplicity and stability limitations with the temperature error at any time point being of the order Δt , $O(\Delta t)$, provided the stability criterion is satisfied. For a rapidly changing boundary condition, such as a heat source, there is no assurance that the calculated temperatures are accurate during the transient period, particularly near the start of the transient, even though the stability criterion is satisfied. Since the allowable time step is governed by the smallest time constant of the network, care must be given in reducing the physical system to a lumped-parameter model. Nonlinearity due to the presence of thermal radiation exchange or temperature-time varying coefficients can lead to numerical solution difficulties; the presence of arithmetic nodes can also present difficulties. These routines offer a number of control constants many of which can be optionally specified by the user to affect the numerical results.

Even with the experience gained through the use of these routines, no realistic criteria can be stated except for the qualitative guidelines indicated above. It is thus recommended that the user becomes familiar with various control constants and their role. The presentation

to follow is intended to provide the instructional information.

6.3.1.2 Finite Difference Approximation and Computational Algorithm

The forward finite difference explicit formulation of the lumped parameter heat balance equations was presented in Section 5.2.1. For convenience, the expression is repeated here.

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_{i,n} - \sum_{j=1}^p a_{ij} (T_{j,n} - T_{i,n}) + \sum_{j=1}^p \sigma b_{ij} (T_{j,n}^4 - T_{i,n}^4)$$

(From equation 5.2-1 of Section 5.2.1)

where, $i = 1, 2, \dots, N$

$T_{j,n}$ = constant, $N < j \leq p$

p = total number of nodes

Δt = time-step

n = nth time-step

By letting $G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{j,n}^2 + T_{i,n}^2)(T_{j,n} + T_{i,n})$, equation (5.2-1) becomes,

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_{i,n} + \sum_{j=1}^p G_{ij,n} (T_{j,n} - T_{i,n}) \quad (6.3-1)$$

The algorithm as used in the subroutines for the diffusion nodes and for the arithmetic nodes may be expressed as follows.

Diffusion Nodes

$$T_{i,n+1} = T_{i,n} + \frac{\Delta t}{C_i} \left[q_{i,n} + \sum_{j=1}^p G_{ij,n} (T_{j,n} - T_{i,n}) \right] \quad (6.3-2)$$

where, n = nth time-step

Δt = time-step (refer to Section 6.2.4)

$i = 1, 2, \dots, NND$ (number of diffusion nodes)

$T_{j,n}$ = constant, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

C_i, q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 through 6.2-4).

Arithmetic Nodes (if any)

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{\left(q_{i,n} + \sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,n} T_{j,k} \right)}{\sum_{j=1}^P G_{ij,n}} \quad (6.3-3)$$

where, $k = k$ th iteration loop; $i = (NND + 1), (NND + 2), \dots, (NND + NNA)$

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

$T_{j,k}$ = constant, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$DN \equiv$ DAMPA (arithmetic node damping factor, refer to Section 6.2.3.2)

$DD = 1.0 - DN$

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$$

($\ell = k$, if $j \geq i$ and $\ell = k+1$, if $j < i$)

6.3.1.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutines CNFRWD and CNFRDL are indicated in Table 6.3-1. For a detailed step-by-step computational description, the user must examine the computer listings for CNFRWD and CNFRDL presented in Appendix A, but some general computational details are given in Section 6.2.5.1. Both CNFRWD and CNFRDL use essentially the same computational steps with the difference occurring in the calculation of the diffusion-node temperatures as shown in the flow chart of Figure 6.3-1; a flow chart for the calculation of the arithmetic-node temperatures is shown in Figure 6.3-2. A functional flow chart of CNFRWD and CNFRDL is shown in Figure 6.3-3. The difference between CNFRWD and CNFRDL is due to the use of the short pseudo-compute sequence (SPCS) by CNFRWD and the use of the long pseudo-compute sequence (LPCS) by CNFRDL.

All diffusion-node temperatures are calculated by a two-pass operation prior to the calculation of the arithmetic node temperatures. On the first pass the pseudo-compute sequence for the diffusion nodes is addressed and the heat flow is calculated and the direction determined for each conductor encountered; the appropriate heat flow and conductance summations are performed. Refer to Section 6.2.5.1 for more details on the computational procedure.

The stability criterion of each diffusion node is calculated and the minimum value is placed in control constant CSGMIN. The time-step used (stored in control constant DTIMEU) is calculated as 95% of CSGMIN divided by control constant CSGFAC which is set at 1.0 unless specified larger by the user. A "look ahead" feature is used when DTIMEU is calculated. If one time-step will pass the output time point the time-step is set to lie on the output time point; if two time-steps will pass the output time point, the time-step is set so that the end of the two time-steps will lie on the output time point. DTIMEU is checked against both DTIMEH and DTIMEL. If DTIMEU exceeds DTIMEH, DTIMEU is set equal to DTIMEH, and if DTIMEU is less than DTIMEL, the "run" is terminated. DTIMEL is internally set to zero if not specified and DTIMEH is set to 1.0E+8 if not specified. The maximum diffusion node temperature change over a time-step is placed in control constant DTMPC and is checked against the allowable diffusion node temperature change stored in the optionally user specified control constant DTMPCA which is not specified is set to 1.0E+8. If DTMPC is larger than DTMPCA, DTIMEU is shortened and the calculations repeated. Refer to Section 6.2.4 for detailed procedure on time-step calculation.

The user may iterate the arithmetic node calculations during a time-step by specifying control constant NLØØP and adjust the solution by the use of ARLXCA. The maximum arithmetic node temperature change over an iteration is placed in control constant ARLXCC and is checked against the arithmetic node temperature change criterion stored in ARLXCA. Satisfaction of either NLØØP or ARLXCA terminates the iterative process for that time-step. If the arithmetic node iteration count exceeds NLØØP the results are retained and computation proceeds without user notification. The maximum arithmetic node temperature change over the time-step is stored in control constant ATMPCC and is checked against the allowable temperature change stored in ATMPCA. If larger, the time-step is shortened and the calculation repeated. The user may also specify the control constant DAMPA in order to dampen possible oscillation due to nonlinearities.

6.3.1.4 Control Constants

Control constants ØOUTPUT and TIMEND (> TIMEØ) must be specified as indicated in Table 6.2-5 and described in Section 6.2.3.2; otherwise the "run" will terminate with an error message. The function of optionally

specified control constants ARLXCA, ATMPCA, BACKUP, CSGFAC, DAMPA, DTIMEH, DTIMEL, DTMPCA, NLØØP, and TIMEØ is described in Section 6.2.3.2. Note particularly that TIMEØ may be set negative and that NLØØP is set to one if not specified.

6.3.1.5 Error and Other Messages

If control constants ØOUTPUT and TIMEND are not specified, the following error message will be printed for each,

ØOUTPUT	"NØ ØOUTPUT INTERVAL"
TIMEND	"TIME STEP TØØ SMALL"

The reason for the TIMEND error message is that a direct check on TIMEND is not made; the resultant error message just happens to be a quirk in the coding.

If the short pseudo-compute sequence SPCS is not specified, the error message will be,

"CNFRWD REQUIRES SHORT PSEUDØ-COMPUTE SEQUENCE"

If the long pseudo-compute LPCS is not specified, the error message will be,

"CNFRDL REQUIRES LONG PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < (NND + NNA)$), the message will be,

"_____ LØCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If the time-step used is less than the time-step allowed (DTIMEL) which may be optionally specified by the user, the message will be,

"TIME STEP TØØ SMALL"

If $CSGMIN \leq 0$, the message printed will be,

"CSGMIN ZERØ or NEGATIVE"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the "run" terminating if a single check is not satisfied,

OUTPUT, pseudo-compute sequence, dynamic storage locations

It should be particularly noted that no message is printed if ARLXCA is not satisfied with NLØØP iterations; ARLXCA and NLØØP are optionally specified control constants.

Table 6.3-1. Basic Computational Steps for CNFRWD and CNFRDL

1. Specification of control constants (all control constants are pre-set to zero). Control constants ØUTPUT and TIMEND must be specified. SPCS is required for CNFRWD and LPCS for CNFRDL. (Refer to Table 6.2-4 for nominal values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = NND + NNA (NND = diffusion nodes and NNA = arithmetic nodes).
3. Setting and/or calculation of time-step, Δt. (Refer to Section 6.2.4 for detailed procedure.)
4. Setting of source and diffusion node dynamic storage locations at zero.
5. Calling of VARIABLES 1. (Refer to Section 6.2.2.2.)
6. Checking of BACKUP. (Refer to Section 6.2.3.2.)
7. Calculation of diffusion-node temperatures. (Refer to Section 6.2.5.1 for description and to flow chart of Figure 6.3-1.)

Diffusion-node temperatures are calculated by using: (refer to Section 6.3.1.2.)

$$T_{i,n+1} = T_{i,n} + \Delta T_{i,n}$$

$$\text{where, } \Delta T_{i,n} = \frac{\Delta t}{C_{i,n}} \left[q_{i,n} + \sum_{j=1}^P G_{ij,n} (T_{j,n} - T_{i,n}) \right]$$

8. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion DTMPCA is not satisfied and recalculation of temperatures with reduced time-step.
9. Calculation of arithmetic-node temperatures; if the number of iterations equals NLØØP the temperatures are retained without user modification. (Refer to Section 6.2.5.1 for description and to flow chart of Figure 6.3.2)
10. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion ATMPCA is not satisfied and recalculation of temperatures with reduced time-step.
11. Setting of BACKUP to 0.0 and the calling of VARIABLES 2.
If BACKUP is nonzero, temperatures are re-set to former values and the computational procedure repeated.
12. Advancing of time, checking of time to print, and the printing at the output interval.
13. Calling of ØUTPUT CALLS.
14. Checking for problem end-time stored in user specified control constant TIMEND.

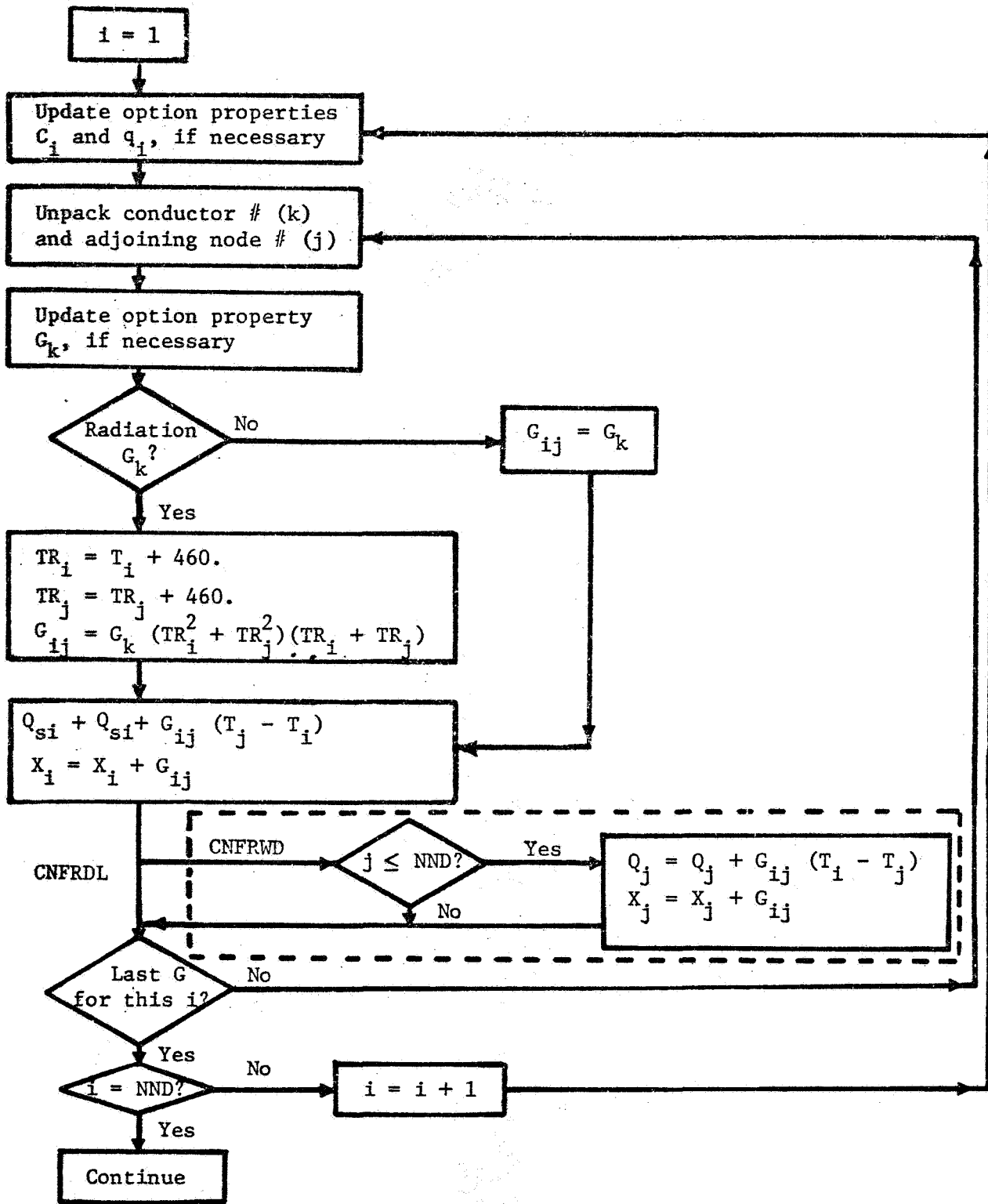


Figure 6.3-1. QSUM and GSUM for "Block" Diffusion-Node Temperature Calculation, CNFRWD and CNFRDL

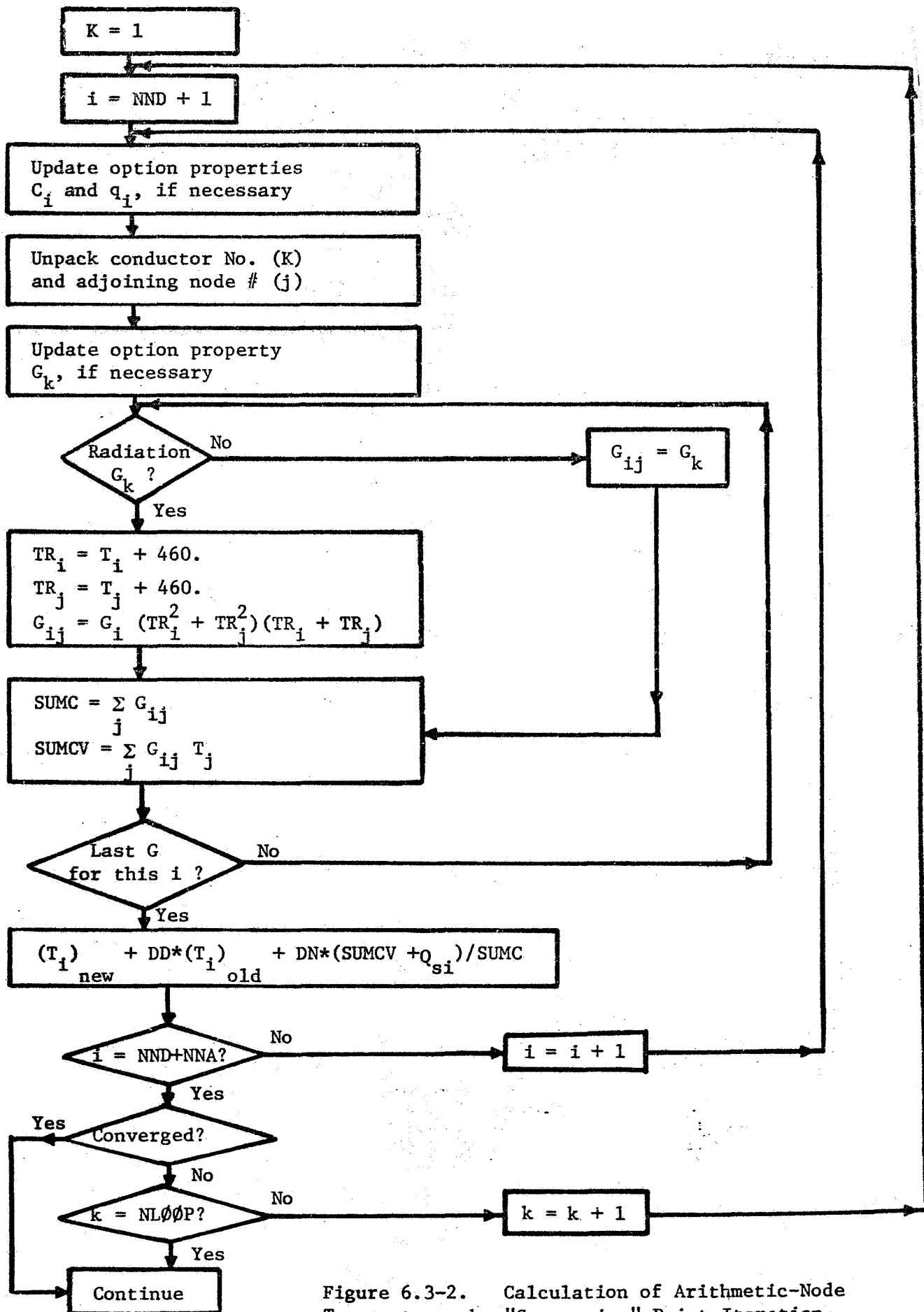


Figure 6.3-2. Calculation of Arithmetic-Node Temperatures by "Successive" Point Iteration

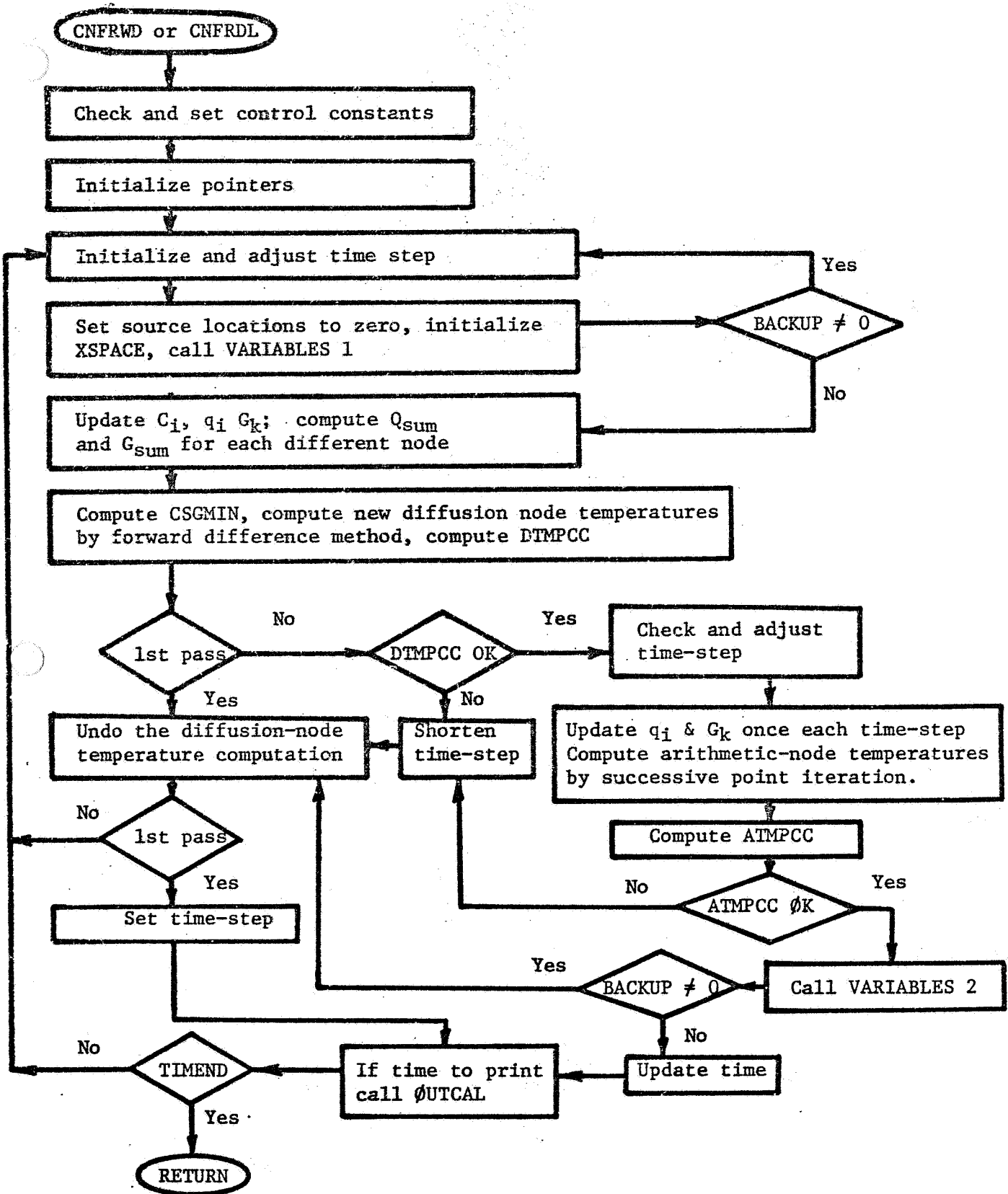


Figure 6.3-3. Functional Flow Chart for CNFRWD and CNFRDL

6.3.2 Subroutine: CNFAST

6.3.2.1 General Comments

Subroutine CNFAST, which requires the short pseudo compute sequence (SPCS) represents a modified CNFRWD with the modifications intended to decrease the computational time. Use of CNFAST requires a user specification of control constant DTIMEL which represents the minimum time-step allowed in addition to control constant ØUTPUT. With minimum computational time and adequate temperature values as the objective, the computational procedure is simplified. A number of checks on control constants are eliminated and temperature nodes with CSGMIN less than the allowable time-step, DTIMEL, are calculated using the steady state equations.

Although experience on the use of CNFAST is rather limited at this time, it is clear that the user specified DTIMEL should be sufficiently small that only a small number of the diffusion nodes should receive the steady state equations. These steady state equations are computed only once during a time-step and thus are not treated computationally the same as the other user-specified arithmetic nodes. A large pocket of internally converted diffusion nodes would lead to large temperature inaccuracies.

6.3.2.2 Finite Difference Approximation and Computational Algorithm

The finite difference expressions for CNFAST are the same as those indicated in Section 6.3.1.2 for subroutines CNFRWD and CNFRDL, but the application of these equations in the computation procedure is different.

Diffusion Nodes

If the user specified control constant, DTIMEL, which represents the maximum time-step allowed as specified by the user is less than or equal to CSGMIN, the diffusion node temperature is calculated as,

$$T_{i,n+1} = T_{i,n} + \frac{\Delta t}{C_i} \left[q_{i,n} + \sum_{j=1}^p G_{ij,n} (T_{j,n} - T_{i,n}) \right] \quad 6.3-4$$

where, Δt = time-step (refer to Section 6.2.4); n = nth time-step

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

i = 1,2,...,NND (of diffusion nodes with DTIMEL \leq CSGMIN)

$T_{j,n}$ = constant, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$G_{ij,n} = a_{ij,n} + ob_{ij,n} (T_{j,n}^2 + T_{i,n}^2)(T_{j,n} + T_{i,n})$$

If DTIMEL > CSGMIN, the time-step is set at DTIMEL and the diffusion node temperature calculated with no iterations as,

$$T_{i,n+1} = \left(\frac{q_{i,n} + \sum_{j=1}^P G_{ij,n} (T_{j,n} - T_{i,n})}{\sum_{j=1}^P G_{ij,n}} \right) \quad (6.3-5)$$

where, n means the nth time-step

i = 1,2,...,NND (number of diffusion nodes with DTIMEL > CSGMIN)

Arithmetic Nodes (if any)

The arithmetic-node temperatures are calculated in the same manner as in CNFRWD (Section 6.3.1.2) or refer to Section 5.2.3 for finite difference algorithm.

6.3.2.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutine CNFAST are indicated in Table 6.3-2 and a functional flow chart is shown in Figure 6.3-4. For a detailed computational description, the user should examine the computer listing for CNFAST in Appendix A, but some general computational details are presented in Section 6.2.5.1. The computational procedure is similar to the one used in CNFRWD with the major difference being the use of DTIMEL which represents the user specified minimum time-step allowed. The time-step calculations stored in DTIMEU proceed exactly as in CNFRWD until the check with DTIMEL is made. If DTIMEU (CSGMIN of a node) \geq DTIMEL, the diffusion node temperature calculation is identical to CNFRWD. If DTIMEU (CSGMIN of a node) < DTIMEL, the diffusion node receives the steady state calculation.

Control constants DTMPCA which contains the allowable diffusion-node temperature change and ATMPCA which contains the arithmetic-node temperature change are not checked in CNFAST. Thus time-steps are not shortened and temperature calculations repeated. The remainder of the computational procedure follows those of CNFRWD (Section 6.3.1.3).

6.3.2.4 Control Constants

Control constants DTIMEL, \emptyset OUTPUT and TIMEND (\emptyset TIME \emptyset) must be specified as indicated in Table 6.2-5 and described in Section 6.2.3.2;

otherwise the "run" will terminate with an error message. The function of optionally specified control constants ARLXCA, BACKUP, DAMPA, DTIMEH, NLØØP and TIMEØ is described in Section 6.2.3.2. As mentioned before in a previous paragraph, the user should take considerable amount of caution in specifying DTIMEL in order to prevent large pockets of nodes that receive the steady state equation without reiteration. Note also that TIMEØ may be set negative and that NLØØP is set to one if not specified.

6.3.2.5 Error and Other Messages

If control constants DTIMEL, ØUTPUT and TIMEND are not specified, the following error message will be printed for each,

DTIMEL	"NØ DTIMEL"
ØUTPUT	"NØ ØUTPUT INTERVAL"
TIMEND	no message

A direct check on TIMEND is not made; an indirect message is printed for the other explicit routines but is not output for CNFAST.

If the short pseudo-compute sequence SPCS is not specified, the error message will be,

"CNFAST REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < NND$), the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If $CSGMIN \leq 0$, the message printed will be,

"C/SK ZERØ or NEGATIVE"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence, with the run terminating if a single check is not satisfied,

ØUTPUT, DTIMEL, pseudo-compute sequence, and dynamic storage locations.

It should be particularly noted that no message is printed if ARLXCA is not satisfied with NLØØP iterations; ARLXCA and NLØØP are optionally specified control constants.

Table 6.3-2. Basic Computational Steps for CNFAST

1. Specification of control constants. Control constants DTIMEL, ØOUTPUT and TIMEND must be specified. SPCS is required for CNFAST. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for descriptions.)
2. Sufficiency check on dynamic storage. Requirements = NND (NND = diffusion nodes).
3. Setting and/or calculation of time-step, Δt. (Refer to Section 6.2.4 for detailed procedure.)

Note that initial time-step equal DTIMEL and subsequent time-step is the larger of CSGMIN or DTIMEL.

4. Setting of source and diffusion node dynamic storage locations to zero.
5. Calling of VARIABLES 1. (Refer to Section 6.2.2.2 for description.)
6. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
7. Calculation of diffusion-node temperatures. (Refer to Section 6.2.5.1 for description.) Calculation differs from the other explicit routines, since diffusion nodes with CSGMIN less than DTIMEL receive steady state calculation (refer to Section 6.3.2.2.)

If $DTIMEL \leq CSGMIN$, the node temperature is calculated as,

$$T_{i,n+1} = T_{i,n} + \frac{\Delta t}{C_i} \left[\sum_{j=1}^P G_{ij,n} (T_{j,n} - T_{i,n}) + q_{i,n} \right]$$

If $DTIMEL > CSGMIN$, the node temperature is calculated using the steady state expression,

$$T_{i,n+1} = \left(\frac{q_{i,n} + \sum_{j=1}^P G_{ij,n} (T_{j,n} - T_{i,n})}{\sum_{j=1}^P G_{ij,n}} \right)$$

8. Calculation of arithmetic-node temperatures if the number of iterations equals NLØØP the temperatures are retained without user notification. (Refer to Section 6.2.5.1 for details.)
9. Calling of VARIABLES 2. (Refer to Section 6.2.2.3 for description.)
10. Advancing of time, checking of time to print, and the printing at the the output interval.
11. Calling of ØOUTPUT CALLS.
12. Checking for problem end-time stored in user specified control constant TIMEND

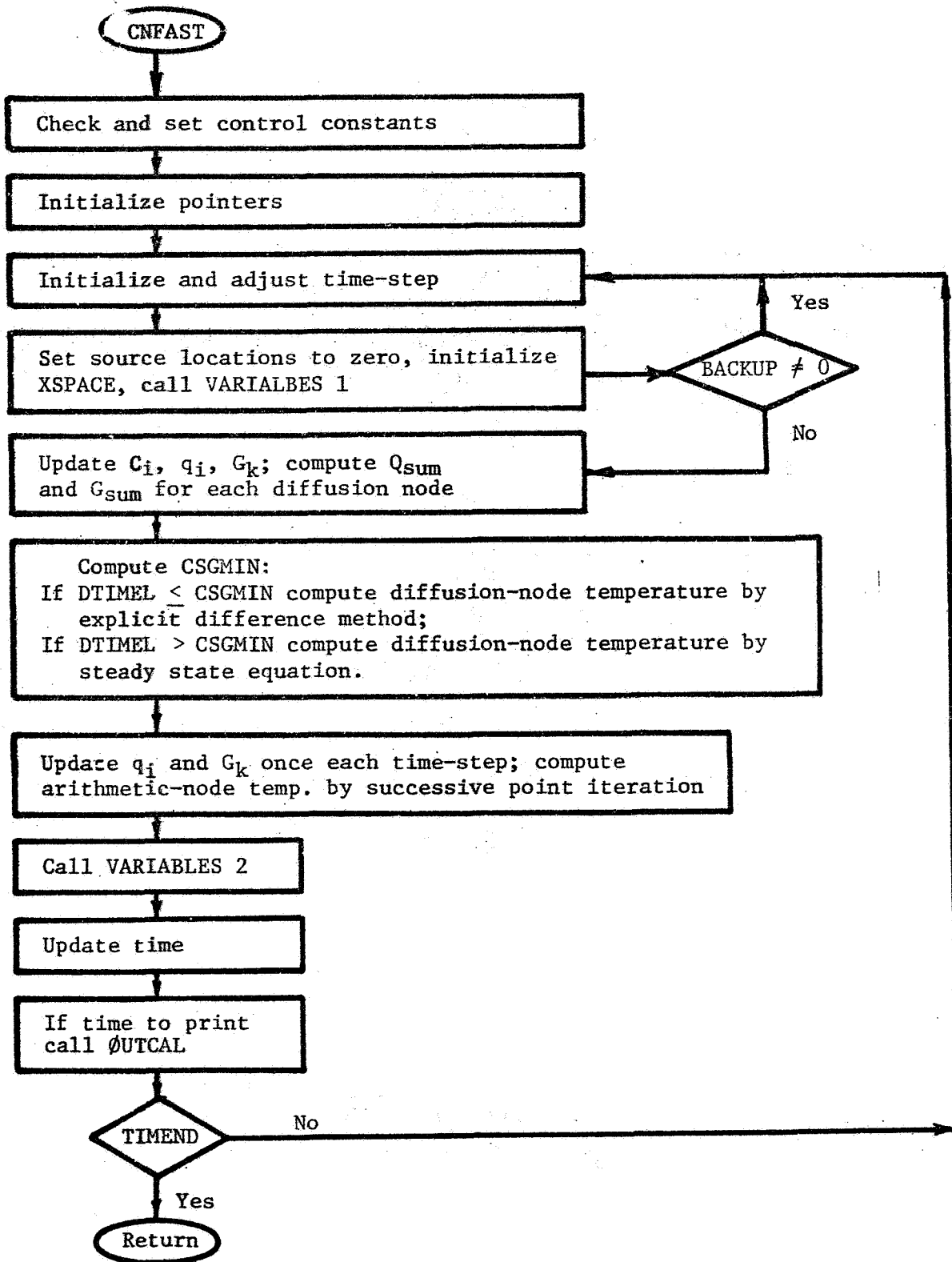


Figure 6.3-4. Functional Flow Chart for CNFAST

6.3.3 Subroutine: CNEXPN

6.3.3.1 General Comments

Subroutine CNEXPN is an explicit routine based upon the exponential prediction method;^{1, 17} the method being unconditionally stable permits any size time-steps and requires the short pseudo-compute sequence (SPCS). An infinite time-step reduces the transient equation to a steady state one. Although the method is unconditionally stable, stability should not be confused with accuracy. Comparison of several numerical methods, including the exponential approximation, is given in Reference 17.

If accuracy is an important consideration, time-steps should not be larger than those taken with the standard explicit method such as used in CNFRWD. If high accuracy is not an important consideration, considerable savings in computational time can be affected with the use of a large time-step. It should be noted that the same savings in computational time may be possible with the implicit routines. As another note of interest, CNEXPN solutions have a tendency to lag in time the true temperatures.

6.3.3.2 Finite Difference Approximation and Computational Algorithm

Diffusion Nodes

The expression for the numerical method used in subroutine CNEXPN for solving the diffusion-node temperatures may be derived from the heat balance equation (5.1-6).

$$\frac{dT_i}{dt} = \frac{1}{C_i} \left[q_i + \sum_{j=1}^p a_{ij} (T_j - T_i) + \sum_{j=1}^p \sigma b_{ij} (T_j^4 - T_i^4) \right] \quad \begin{array}{l} \text{(equation} \\ \text{5.1-6 of} \\ \text{Section 5)} \end{array}$$

$$i = 1, 2, \dots, N$$

$$T_j = \text{constant}, N < j \leq p$$

If $G_{ij} = a_{ij} + \sigma b_{ij} (T_j^2 + T_i^2)$ equation (5.1-6) becomes,

$$\frac{dT_i}{dt} = \frac{1}{C_i} \left[q_i + \sum_{j=1}^p G_{ij} (T_j - T_i) \right] \quad (6.3-6)$$

$$i = 1, 2, \dots, N$$

$$T_j = \text{constant}, N < j \leq p$$

If we further let G_{ij} , q_i and T_j be invariant with time and temperature, equation (6.3-6) may be integrated rather easily to yield,

$$T_{i,n+1} = T_{i,n} e^{-\alpha_n \Delta t} + \frac{q_{i,n} + \sum_{j=1}^p G_{ij,n} T_{j,n}}{\sum_{j=1}^p G_{ij,n}} \left(1 - e^{-\alpha_n \Delta t}\right) \quad (6.3-7)$$

where, $n = \text{nth time-step}$; $i = 1, 2, \dots, \text{NND}$ (number of diffusion nodes)
 C_i, q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)
 $T_{j,n} = \text{constant}$, $(\text{NND} + \text{NNA}) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$\alpha_n = \frac{\sum_{j=1}^p G_{ij,n}}{C_{i,n}}$$

$\Delta t = \text{time-step}$ (refer to Section 6.2.4)

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{i,n}^2 + T_{j,n}^2)(T_{i,n} + T_{j,n})$$

Computationally equation (6.3-7) is applied to the diffusion nodes. It should be noted that the form of equation (6.3-7) represents a "block" change in temperatures since the evaluation of $T_{i,n+1}$ is based upon $T_{i,n}$.

Arithmetic Nodes (if any)

Arithmetic-node temperatures are calculated in the same manner as in CNFRWD (Section 6.3.1.2) or refer to Section 5.2.3 for finite difference algorithm.

6.3.3.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutine CNEXPXN are indicated in Table 6.3-3 and a functional flow chart is shown in Figure 6.3-5. A detailed computational procedure requires the examination of the CNEXPXN computer listing which is presented in Appendix A but some general computational details are given in Section 6.2.5.1. The computational process of subroutine CNEXPXN is essentially identical to CNFRWD with the difference being the finite difference expression used for the calculation of the diffusion nodes and the time-step which is calculated as $\text{CSGMIN} * \text{CSGFAC}$ in lieu of $\text{CSGMIN} / \text{CSGFAC}$. The "look ahead" feature for

time-step calculation as well as a check with DTIMEH, DTIMEL and DTMPCA is identical to CNFRWD. Temperatures of arithmetic nodes are calculated after the diffusion nodes and utilize NLØØP, ARLXCA, and DAMPA in exactly the same way as CNFRWD. The verbal flow description of CNFRWD (Section 6.3.1.3) applies here except for the differences indicated above.

6.3.3.4 Control Constants

Control constants ØUTPUT and TIMEND (> TIMEØ) must be specified as indicated in Table 6.2-5 and described in Section 6.2.3.2; otherwise the "run" will terminate with an error message. The function of optionally specified control constants ARLXCA, ATMPCA, BACKUP, CSGFAC, DAMPA, DTIMEH, DTIMEL, DTMPCA, NLØØP, and TIMEØ is described in Section 6.2.3.2. The user should take particular care in the selection of CSGFAC since too large of a time-step would lead to grossly inaccurate temperatures even though the solution is stable. Note also that TIMEØ may be set negative and that NLØØP is set to one if not specified.

6.3.3.5 Error and Other Messages

If control constants ØUTPUT and TIMEND are not specified, the following error message will be printed for each,

ØUTPUT	"NØ ØUTPUT INTERVAL"
TIMEND	"TIME STEP TØØ SMALL"

The reason for the TIMEND error message is that a direct check on TIMEND is not made; the resultant error message just happens to be a quirk in the coding.

If the short pseudo-compute sequence SPCS is not specified, the error message will be,

"CNEXPN REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient $NDIM < (NND + NNA)$, the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If the time-step used is less than the time-step allowed (DTIMEL) which may be optionally specified by the user, the message will be,

"TIME STEP TOO SMALL"

If CSGMIN \leq 0, the message printed will be,

"CSGMIN ZERO or NEGATIVE"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the run terminating if a single check is not satisfied,

OUTPUT, pseudo-compute sequence, dynamic storage locations.

It should be particularly noted that no message is printed if ARLXCA is not satisfied with NLPPP iterations; ARLXCA and NLPPP are optionally specified control constants.

Table 6.3-3. Basic Computational Steps for CNEXPAN

1. Specification of control constants. Control constants \emptyset UTPUT and TIMEND must be specified. SPCS is required for CNEXPAN. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = NND + NNA (NND = diffusion nodes and NNA = arithmetic nodes).
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Calculated time-step = $0.95 * CSGMIN * CSGFAC$.
4. Setting of source and diffusion node dynamic storage locations to zero.
5. Calling of VARIABLES 1. (Refer to Section 6.2.2.2 for description.)
6. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
7. Calculation of diffusion-node temperatures. (Refer to Section 6.2.5.1 for description.)

Diffusion-node temperatures are calculated by using (refer to Section 6.3.3.2),

$$T_{i,n+1} = T_{i,n} e^{-\alpha_n \Delta t} + \frac{q_{i,n} + \sum_{j=1}^P G_{ij,n} T_{j,n}}{\sum_{j=1}^P G_{ij,n}} \left(1 - e^{-\alpha_n \Delta t} \right)$$

where,

$$\alpha_n = \frac{\sum_{j=1}^P G_{ij,n}}{C_{i,n}}$$

8. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion DTMPA is not satisfied and recalculation of temperatures with reduced time-step.
9. Calculation of arithmetic-node temperatures. If the number of iterations equal NL $\emptyset\emptyset$ P, the temperatures are retained without user notification
10. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion ATMPA is not satisfied and recalculation of temperatures with reduced time-step.
11. Calling of VARIABLES 2 and checking of BACKUP. (Refer to Section 6.2.2.3 and 6.2.3.2 for description.)
12. Advancing of time, checking of time to print, and the printing at the output interval.
13. Calling of \emptyset UTPUT CALLS.
14. Checking for problem end time stored in user specified control constant TIMEND.

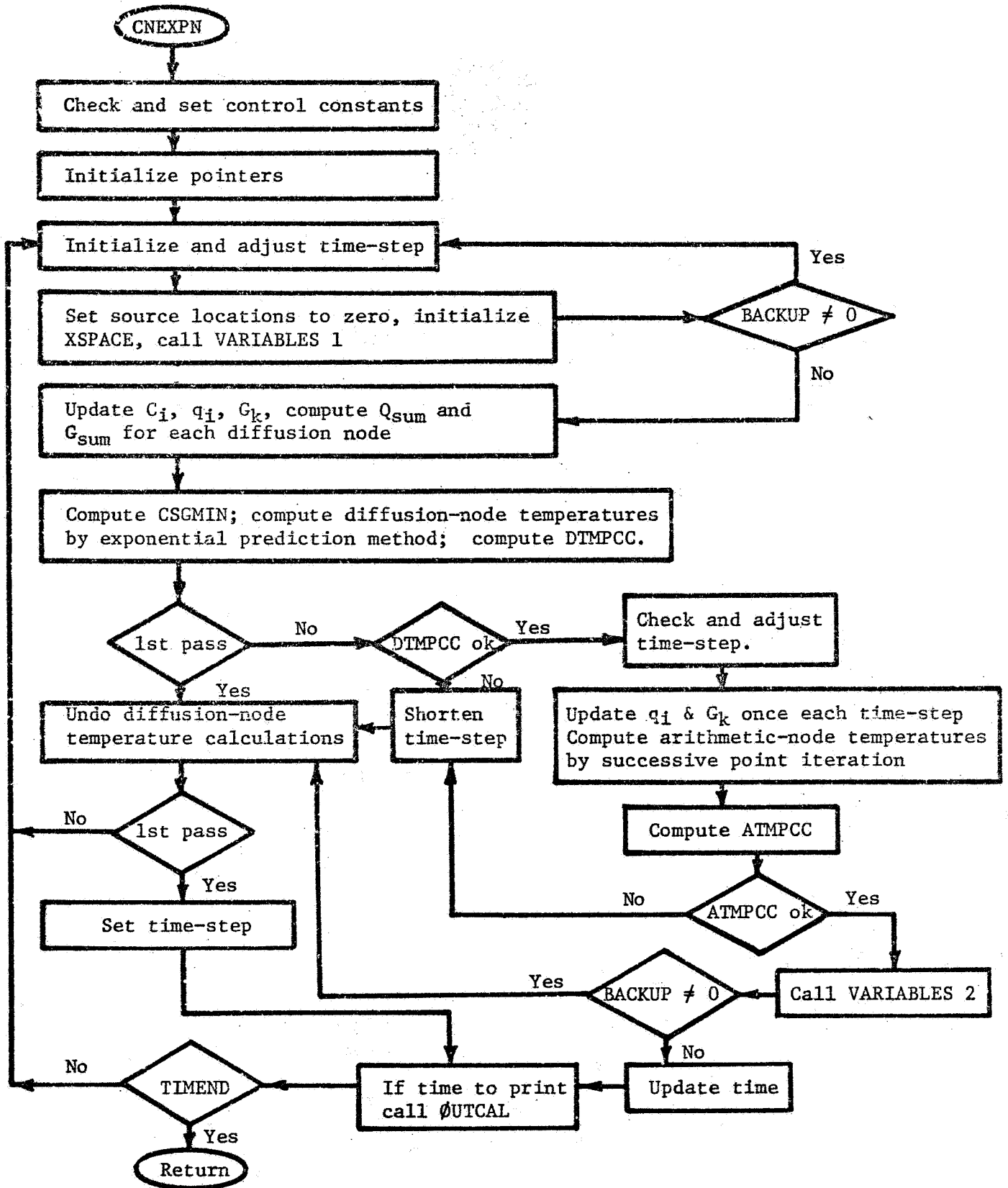


Figure 6.3-5. Functional Flow Chart for CNEXPN

6.3.4 Subroutine: CNDUFR

6.3.4.1 General Comments

Subroutine CNDUFR is an explicit numerical solution routine that uses an unconditionally stable DuFort-Frankel method.^{9, 12, 17} The DuFort-Frankel method replaces the present temperature of the node being operated on by the average of future and past temperatures in the forward differencing equation. In subroutine CNDUFR the present temperature of the node being operated on is replaced by a time-weighted average of future and past temperatures. CNDUFR requires the short pseudo-compute sequence (SPCS).

The intent of an unconditionally stable routine such as CNDUFR is the reduction of computational time by using time-steps greater than those allowed with the conditionally stable explicit methods as constrained by the stability criterion. However, less accuracy can be expected with a lengthened time-step. The time-step controlled with control constant CSGFAC represents a user decision that is difficult and must be aided by a trial and error procedure.

Examination of several CNDUFR solutions reveals a tendency to lead in time the true temperatures.

6.3.4.2 Finite Difference Approximation and Computational Algorithm

Diffusion Nodes

The DuFort-Frankel explicit finite difference expression^{9, 12, 17} for calculating the diffusion-node temperatures may be readily determined as follows:

Using the standard explicit finite difference expression,

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_{i,n} + \sum_{j=1}^p G_{ij,n} (T_{j,n} - T_{i,n}) \quad (6.3-8)$$

$$i = 1, 2, \dots, N$$

$$T_{j,n} = \text{constant}, N < j \leq p$$

letting the present temperature, $T_{i,n}$, be replaced by the average of future temperature, $T_{i,n+1}$, and past temperature, $T_{i,n-1}$,

$$T_{i,n} = \frac{T_{i,n+1} + T_{i,n-1}}{2} \quad (6.3-9)$$

where, $\Delta t_i = \Delta t_{i+1}$, $i = 1, 2, \dots, M$ (equal time-steps)

and defining,

$$\bar{C}_i = C_i / \Delta t \quad (\text{refer to Section 6.2.4 for discussion on } \Delta t) \quad (6.3-10)$$

equation (6.3-8) can be expressed as,

$$T_{i,n+1} = \frac{\bar{C}_{i,n} T_{i,n-1} + 2 q_{i,n} + \sum_{j=1}^P G_{ij,n} (2 T_{j,n} - T_{i,n-1})}{\bar{C}_{i,n} + \sum_{j=1}^P G_{ij,n}} \quad (6.3-11)$$

$i = 1, 2, \dots, N$

In CNDUFR the present temperature, $T_{i,n}$, of equation (6.3-8) is replaced by a weighted average of future temperature, $T_{i,n+1}$, and past temperature, $T_{i,n-1}$. The weighting is based on unequal time-steps.

$$T_{i,n} = \frac{(\Delta t_{n-1} T_{i,n+1} + \Delta t_n T_{i,n-1})}{\Delta t_{n-1} + \Delta t_n} \quad (6.3-12)$$

where, $\Delta t_{n-1} = t_n - t_{n-1}$ (past time-step)

$$\Delta t_n = t_{n+1} - t_n \quad (\text{present time-step})$$

Let

$$\tau_{n-1} = \frac{\Delta t_{n-1}}{\Delta t_{n-1} + \Delta t_n} \quad (6.3-13)$$

$$\tau_n = \frac{\Delta t_n}{\Delta t_{n-1} + \Delta t_n} \quad (6.3-14)$$

Equation (6.3-8) becomes,

$$T_{i,n+1} = \frac{\tau_n T_{i,n-1} \left(\bar{C}_{i,n} - \sum_{j=1}^P G_{ij,n} \right) + \sum_{j=1}^P G_{ij,n} T_{j,n} + q_{i,n}}{\bar{C}_{i,n} - \tau_{n-1} \left(\bar{C}_{i,n} - \sum_{j=1}^P G_{ij,n} \right)} \quad (6.3-15)$$

where, $i = 1, 2, \dots, NND$ (number of diffusion nodes)

$T_{j,n} = \text{constant}$, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{j,n}^2 + T_{i,n}^2) (T_{j,n} + T_{i,n})$$

C_i, q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

In CNDUFR, equation (6.3-15) is applied to the diffusion nodes with the computational procedure being a "block" change in temperature from one time-step to another.

Arithmetic Nodes (if any)

Arithmetic-node temperatures are calculated in the same manner as in CNFRWD (Section 6.3.1.2) or refer to Section 5.2.3 for the finite difference algorithm.

6.3.4.3 Comments on the Computational Procedure

The important steps of the computation procedure used in subroutine CNDUFR are indicated in Table 6.3-4 and a functional flow chart is shown in Figure 6.3-6. A computer listing of CNDUFR is found in Appendix A but some general computational details are given in Section 6.2.5.1. The computational procedure for CNDUFR follows the CNEXPN computational pattern, but with the temperatures of the diffusion nodes calculated by the DuFort-Frankel method of the exponential prediction method. Another significant difference is that CNDUFR must provide for two sets of past temperatures which are required for DuFort-Frankel algorithm; two time-steps for consecutive time-step calculations are also required. Otherwise, checks and control constant use are identical to CNEXPN. Thus, the verbal flow description of Section 6.3.1.3 applies directly except for the differences indicated above.

6.3.4.4 Control Constants

Control constants \emptyset UTPUT and TIMEND ($>$ TIME \emptyset) must be specified as indicated in Table 6.2-5 and described in Section 6.2.3.2; otherwise the "run" will terminate with an error message. The function of optionally specified control constants ARLXCA, ATMPCA, BACKUP, CSGFAC, DAMPA, DTIMEH, DTIMEL, DTMPCA, NL $\emptyset\emptyset$ P, and TIME \emptyset is described in Section 6.2.3.2. The user should take particular care in the selection of CSGFAC since too large of a time-step would lead to grossly inaccurate temperatures even though the solution is stable. Note also that TIME \emptyset may be set negative and that NL $\emptyset\emptyset$ P is set to one if not specified.

6.3.4.5 Error and Other Messages

If control constants \emptyset UTPUT and TIMEND are not specified, the following error message will be printed for each,

\emptyset UTPUT	"N \emptyset \emptyset UTPUT INTERVAL"
TIMEND	"TIME STEP T $\emptyset\emptyset$ SMALL"

The reason for the TIMEND error message is that a direct check on TIMEND is not made; the resultant error message just happens to be a quirk in the coding.

If the short pseudo-compute sequence SPCS is not specified, the error message will be,

"CNDUFR REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < (2*NND + NNA)$), the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If the time-step used is less than the time-step allowed (DTIMEL), which may be optionally specified by the user, the message will be,

"TIME STEP T $\emptyset\emptyset$ SMALL"

If $CSGMIN \leq 0$, the message printed will be,

"CSGMIN ZER \emptyset or NEGATIVE"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the run terminating if a single check is not satisfied,

\emptyset UTPUT, pseudo-compute sequence, dynamic storage locations

It should be particularly noted that no message is printed if ARLXCA is not satisfied with NL $\emptyset\emptyset$ P iterations; ARLXCA and NL $\emptyset\emptyset$ P are optionally specified control constants.

Table 6.3-4. Basic Computational Steps for CNDUFR

1. Setting of control constants to nominal values. Control constants \emptyset UTPUT and TIMEND must be specified. SPCS is required for CNDUFR. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = $2*NND + NNA$ (NND = diffusion nodes and NNA = arithmetic nodes).
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Calculated time-step = $0.95*CSGMIN*CSGFAC$.
4. Setting of source and diffusion node dynamic storage locations to zero.
5. Calling of VARIABLES 1. (Refer to Section 6.2.2.2 for description.)
6. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
7. Calculation of diffusion-node temperatures. (Refer to Section 6.2.5.1 for description.)

Diffusion-node temperatures are calculated by using (refer to Section 6.3.4.2),

$$T_{i,n+1} = \frac{\tau_n T_{i,n-1} (\bar{C}_{i,n} - \sum_{j=1}^P G_{ij,n}) + \sum_{j=1}^P G_{ij,n} T_{j,n} + q_{i,n}}{\bar{C}_{i,n} - \tau_{n-1} (\bar{C}_{i,n} - \sum_{j=1}^P G_{ij,n})}$$

where,

$$\tau_{n-1} = \frac{\Delta t_{n-1}}{\Delta t_{n-1} + \Delta t_n}$$

$$\tau_n = \frac{\Delta t_n}{\Delta t_{n-1} + \Delta t_n}$$

8. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion DTMPCA is not satisfied and temperature recalculation with reduced time-step.
9. Calculation of arithmetic-node temperatures; if the number of iterations equal NL $\emptyset\emptyset$ P, the temperatures are retained without user notification (refer to Section 6.2.5.1 for details).
10. Erasure of arithmetic-node temperatures for latest time-step if allowable temperature change criterion ATMPCA is not satisfied and temperature recalculation with reduced time-step.
11. Calling of VARIABLES 2 and checking of BACKUP. (Refer to Section 6.2.2.3 and 6.2.3.2 for description.)
12. Advancing of time, checking of time to print, and the printing at the output interval.
13. Calling of \emptyset UTPUT CALLS.
14. Checking for problem end time stored in user specified control constant TIMEND.

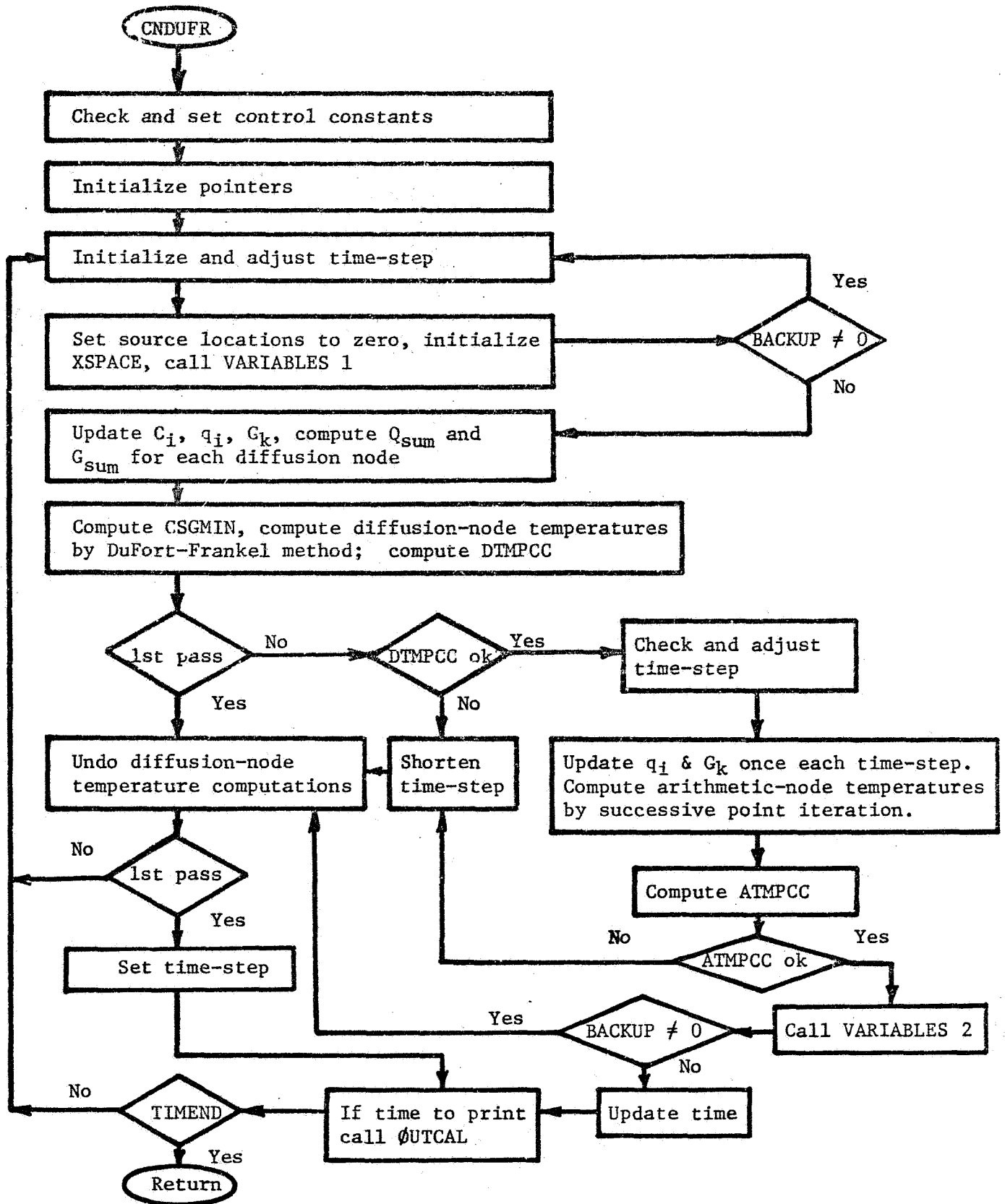


Figure 6.3-6. Functional Flow Chart for CNDUFR

6.3.5 Subroutine: CNQUIK

6.3.5.1 General Comments

Subroutine CNQUIK is a numerical solution routine that uses an algorithm composed of half DuFort-Frankel method^{9, 12, 17} and half exponential prediction method.^{1, 17} CNQUIK requires the short pseudo-compute sequence (SPCS); characteristics of subroutines CNDUFR and CNEXPN, as described in Section 6.3.3 and 6.3.4, also apply to CNQUIK.

Why CNQUIK? Examination of CNDUFR and CNEXPN solutions reveals that CNDUFR has a tendency to yield temperatures which lead the true temperatures, whereas CNEXPN has a tendency to lag the true temperatures. Thus, it was theorized that a combination of CNDUFR and CNEXPN should yield a more accurate solution than either one. Preliminary results indicate that CNQUIK is more accurate than either CNDUFR or CNEXPN with approximately the same solution time. It can also be theorized that a more accurate combination of the DuFort-Frankel and exponential prediction is probably possible than the half and half used in CNQUIK. However, a detailed study will be required before a realistic evaluation of CNQUIK can be made.

6.3.5.2 Finite Difference Approximation and Computational Algorithm

Diffusion Nodes

Subroutine CNQUIK uses a numerical solution algorithm composed of half DuFort-Frankel and half exponential prediction. That is the temperature of the diffusion nodes is calculated by using,

$$T_{i,n+1} = \left(T_{\text{CNDUFR}} + T_{\text{CNEXPN}} \right) / 2.0 \quad (6.3-16)$$

$$T_{\text{CNDUFR}} = \frac{\tau_n T_{i,n-1} \left(\bar{C}_{i,n} - \sum_{j=1}^P G_{ij,n} \right) + \sum_{j=1}^P G_{ij,n} T_{j,n} + q_{i,n}}{\bar{C}_{i,n} (1 - \tau_{n-1}) + \sum_{j=1}^P G_{ij,n}}$$

(equation 6.3-15 of Section 6.3.4.2)

$$T_{\text{CNEXPN}} = T_{i,n} e^{-\alpha_n \Delta t} + \frac{q_{i,n} + \sum_{j=1}^P G_{ij,n} T_{j,n}}{\sum_{j=1}^P G_{ij,n}} \left(1 - e^{-\alpha_n \Delta t} \right)$$

(equation 6.3-7 of Section 6.3.3.2)

n = nth time-step

i = 1,2,...,NND (number of diffusion nodes)

$T_{j,n}$ = constant, (NND + NNA) < j ≤ p (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$\alpha_n = \frac{\sum_{j=1}^p G_{ij,n}}{C_{i,n}}$$

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} (T_{i,n}^2 + T_{j,n}^2)(T_{i,n} + T_{j,n})$$

$$\tau_n = \frac{\Delta t_n}{\Delta t_{n-1} + \Delta t_n}; \quad \tau_{n-1} = \frac{\Delta t_{n-1}}{\Delta t_{n-1} + \Delta t_n}$$

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

$$\bar{C}_i = C_i/\Delta t \text{ (refer to Section 6.2.4 for discussion of } \Delta t)$$

Arithmetic Nodes (if any)

Temperatures of arithmetic nodes are calculated in the same manner as in CNFRWD (Section 6.3.1.2) or refer to Section 5.2.3 for the finite difference algorithm.

6.3.5.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutine CNQUIK are indicated in Table 6.3-5 and a functional flow chart is shown in Figure 6.3-7. A computer listing of CNQUIK is found in Appendix A. General computational details are given in Section 6.2. The computational procedure for CNQUIK follows CNEXP or CNDUFR with the diffusion-node temperatures calculated with the half DuFort-Frankel and half exponential prediction algorithm being the only difference. Arithmetic-node temperatures are calculated in the same manner as the other SINDA explicit routines. Note that the time-step is calculated as CSGMIN*CSGFAC and checks are the same as CNEXP or CNDUFR. Thus, the verbal flow description of Section 6.3.1.3 applies directly except for the differences indicated above.

6.3.5.4 Control Constants

Control constants \emptyset UTPUT and TIMEND ($>$ TIME \emptyset) must be specified as indicated in Table 6.2-5 and described in Section 6.2.3.2; otherwise the "run" will terminate with an error message. The function of optionally specified control constants ARLXCA, ATMPCA, BACKUP, CSGFAC, DAMPA, DTIMEH, DTIMEL, DTMPCA, NL $\emptyset\emptyset$ P, and TIME \emptyset is described in Section 6.2.3.2. Again, caution must be exercised in the selection of CSGFAC since too large of a time-step would lead to grossly inaccurate temperatures even though the solution is stable. Note also that TIME \emptyset may be set negative and that NL $\emptyset\emptyset$ P is set to one if not specified.

6.3.5.5 Error and Other Messages

If control constants \emptyset UTPUT and TIMEND are not specified, the following error message will be printed for each,

\emptyset UTPUT	"N \emptyset \emptyset UTPUT INTERVAL"
TIMEND	"TIME STEP T $\emptyset\emptyset$ SMALL"

The reason for the TIMEND error message is that a direct check on TIMEND is not made; the resultant error message just happens to be a quirk in the coding.

If the short pseudo-compute SPCS is not specified, the error message will be,

"CNQUIK REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient (NDIM $<$ (2*NND + NNA), the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If the time-step used is less than the time-step allowed (DTIMEL), which may be optionally specified by the user, the message will be,

"TIME STEP T $\emptyset\emptyset$ SMALL"

If CSGMIN \leq 0, the message printed will be,

"CSGMIN ZER \emptyset or NEGATIVE"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the run terminating if a single check is not satisfied,

ØOUTPUT, pseudo-compute sequence, dynamic storage locations.

It should be particularly noted that no message is printed if ARLXCA is not satisfied with NLØØP iterations; ARLXCA and NLØØP are optionally specified control constants.

Table 6.3-5. Basic Computational Steps for CNQUIK

1. Specification of control constants. Control constants \emptyset UTPUT and TIMEND must be specified. SPCS is required for CNEXP. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = $2(NND) + NNA$ (NND = diffusion nodes and NNA = arithmetic nodes).
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Calculated time-step = $0.95 * CSGMIN * CSGFAC$.
4. Setting of source and diffusion node dynamic storage locations to zero.
5. Calling of VARIABLES 1. (Refer to Section 6.2.2.2 for description.)
6. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
7. Calculation of diffusion-node temperatures. (Refer to Section 6.2.5.1 for description.)

Diffusion-node temperatures are calculated by using (refer to Section 6.3.5.2),

$$T_{i,n+1} = (T_{CNDUFR} + T_{CNEXP})/2.0$$

(Refer to equation 6.3-17, Section 6.3.5.2.)

8. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion DTMPCA is not satisfied and temperature recalculation with reduced time-step.
9. Calculation of arithmetic-node temperatures; if the number of iterations equal NL $\emptyset\emptyset$ P, the temperatures are retained without user notification. (Refer to Section 6.2.5.1 for details.)
10. Erasure of all temperature calculations for latest time-step if allowable temperature change criterion ATMPCA is not satisfied and temperature recalculation with reduced time-step.
11. Calling of VARIABLES 2 and checking of BACKUP. (Refer to Section 6.2.2.3 and 6.2.3.2 for description.)
12. Advancing of time, checking of time to print, and the printing at the output interval.
13. Calling of \emptyset UTPUT CALLS.
14. Checking for problem end time stored in user specified control constant TIMEND.

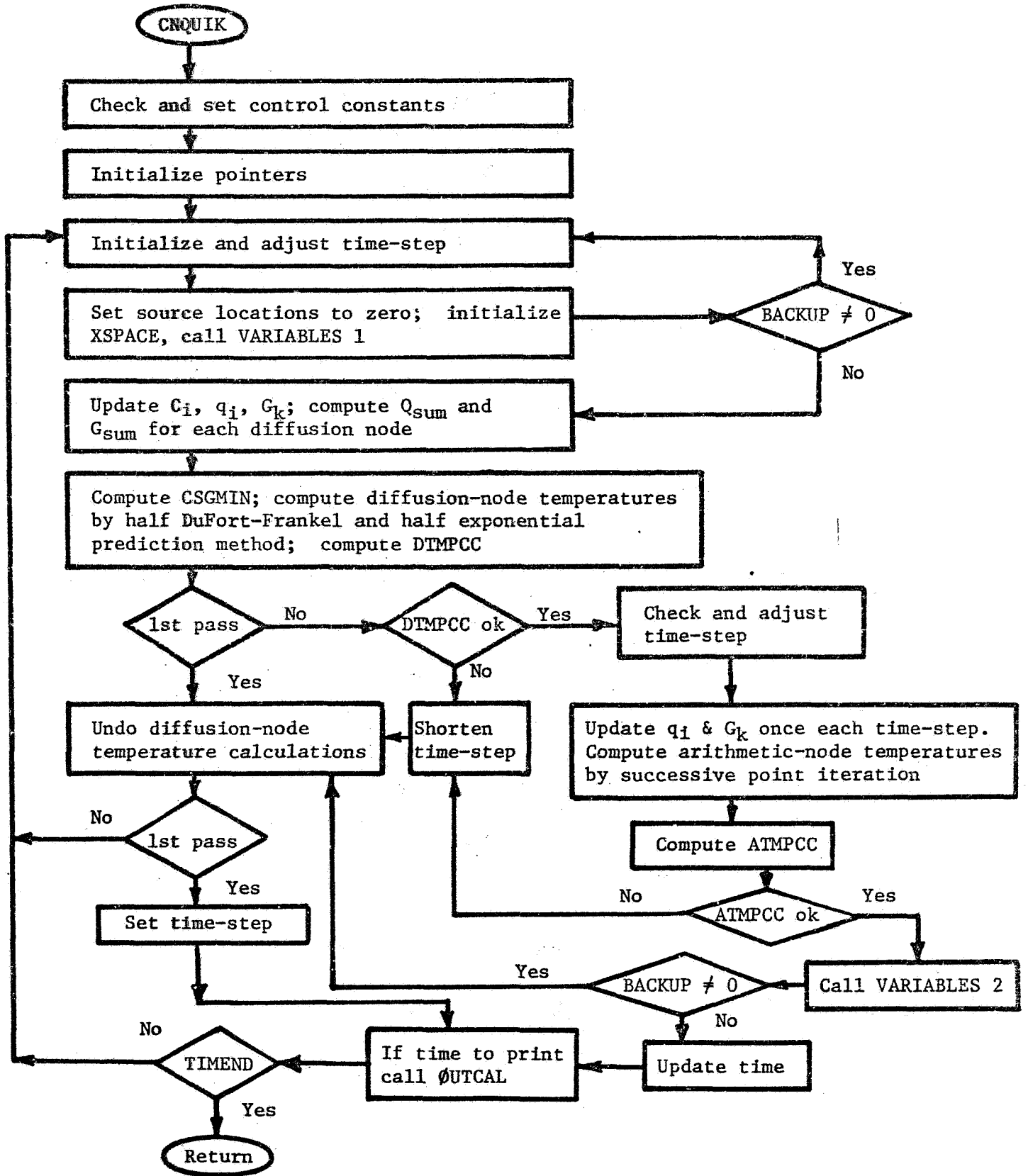


Figure 6.3-7. Functional Flow Chart for CNQUIK

6.4 Transient Implicit Solution Routines

SINDA implicit solution routines number three; these routines are identified as follows:

- CNBACK Implicit backward difference method.
Requires long pseudo-compute sequence (LPCS).
- CNFWBK Implicit forward-backward differencing, using Crank-Nicolson method.
Requires long pseudo-compute sequence (LPCS).
- CNVARB Combination of CNBACK and CNFWBK.
Requires long pseudo-compute sequence (LPCS).

Implicit methods generally tend to be more accurate than explicit methods and are unconditionally stable as are some explicit methods. With implicit methods the time-step is specified in contrast to the calculated time-steps of explicit methods with their stability criterion. An important consideration in the use of implicit methods is that the time-step DTIMEI should be specified in conjunction with control constant NLØØP which represents the maximum number of computational iterations during each time-step. Since each iterative calculation is essentially equivalent to a time-step calculation for an explicit method, the combination of DTIMEI and NLØØP for a given time period should be set less than the total number of time-steps used by the explicit method for the same time period, if computational time is to be reduced; this of course assumes that during each time-step the maximum number of iterations is required. If the NLØØP iterations are required during a time-step, the temperature accuracy is affected but the magnitude would depend upon the value used for the maximum allowable relaxation temperature change criteria, ARLXCA and DRLXCA. It should be noted if NLØØP iterations are required during a time-step, the message "RELAXATION CRITERIA NOT MET" is printed.

A detailed description of each implicit routine, as presented on the pages to follow, relies on the general description of Section 6.2. A brief description of these routines is summarized first.

CNBACK uses the standard backward differencing algorithm and requires the long pseudo-compute sequence (LPCS). The time-step must be

specified via control constant DTIMEI and used in conjunction with the control constant NL00P. CNBACK uses the acceleration of convergence feature.

CNFWBK uses the Crank-Nicolson algorithm which is composed of half forward differencing and half backward differencing. CNFWBK solutions tend to be more accurate than CNBACK solutions with approximately 25% less iterations; however CNFWBK solutions have "blown" on occasions.

CNVARB uses a combination of forward differencing and backward differencing. Unlike CNFWBK which is half and half, CNVARB uses a variable beta factor which ranges from 0 to 1. Thus CNVARB uses a method that is somewhere between forward differencing and backward differencing.

6.4.1 Subroutine: CNEACK

6.4.1.1 General Comments

Subroutine CNEACK is an implicit routine that uses the standard backward difference expression and requires the long pseudo-compute (LPCS). Time-step must be specified via DTIMEI otherwise the "run" will terminate with an error message printout. The time-step value is arbitrary but the user should consider DTIMEI in conjunction with the control constant NLØØP which represents the maximum number of computational iterations during each time-step (refer to Section 6.2.3.2 for description).

Implicit methods tend to be more accurate than explicit methods and are unconditionally stable, but implicit solutions often oscillate at start up or boundary step changes when heat transfer by radiation is present. CNEACK internally controls sudden radiation heat transfer changes by an averaging technique which is termed "radiation damping" (refer to Section 6.2.6 for details). This automatic damping has been very effective in many solutions that have been examined and lessens the need for the use of DAMPD and DAMPA.

6.4.1.2 Finite Difference Approximation and Computational Algorithm

The numerical solution algorithm used in subroutine CNEACK is the standard backward-difference expression^{12, 13, 17} which may be expressed as:

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = q_{i,n} + \sum_{j=1}^p a_{ij} (T_{j,n+1} - T_{i,n+1}) + \sum_{j=1}^p \sigma b_{ij} (T_{j,n+1}^4 - T_{i,n+1}^4)$$

(equation 5.2-5 of Section 5.2.2)

$$i = 1, 2, \dots, N$$

$$T_{j,n+1} = \text{constant}, N < j \leq p$$

$$T_{i,n} \equiv T_i(n\Delta t)$$

The computational procedure for the backward difference formulation must necessarily be re-iterative because of the need to solve a set of simultaneous nonlinear equations.

Diffusion Nodes

Diffusion node temperatures are solved by "successive point" iteration but differs from the arithmetic-node temperature calculation because of the capacitance term and the use of "radiation damping" (refer to Section 6.2.5.2).

$$T_{i,k+1} = DD * T_{i,k}$$

$$T_{i,k+1} = DD * \frac{\bar{C}_{i,n} T_{i,n} + q_{i,n} + \sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,n} T_{j,k} - (q_i)_{ave}}{\bar{C}_{i,n} + \sum_{j=1}^P a_{ij,n}} \quad (6.4-1)$$

where, $i = 1, 2, \dots, NND$

$n =$ nth time-step

$k =$ kth iteration

DN = DAMPD (diffusion-node damping factor)

DD = 1.0 - DN

$$G_{ij,n} = a_{ij,n} + \sigma b_{ij,n} T_{j,\ell}^3 \quad (\ell = k \text{ if } j \geq i \text{ and } \ell = k+1 \text{ if } j < i)$$

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

$$\bar{C}_{i,n} = C_{i,n} / \Delta t \quad (\Delta t = \text{time step, refer to Section 6.2.4})$$

$$(q_i)_{ave} = \sum_{j=1}^P \sigma b_{ij,n} [(T_{i,k}^4) + (T_{i,k}^4)_2] / 2.0, \text{ average heat loss from}$$

the i th node (refer to Section 6.2.6 on radiation damping for details)

Details on the computational procedure for implicit routines are presented in Sections 5.2.2 and 6.2.5.2.

Arithmetic Nodes

Arithmetic-node temperatures are calculated identically the same in all the SINDA numerical solution routines. Thus, refer to either Section 6.3.1.2 or Section 6.2.5.2 for the finite difference algorithm.

6.4.1.3 Comments on the Computational Procedure

The important steps of the computational procedure used in sub-routine CNBACK are indicated in Table 6.4-1. For a detailed step-by-step computational description, the user must examine the computer listing for

CNBACK in Appendix B, but some general computational details are given in Section 6.2.5.2. A functional flow chart of CNBACK is shown in Figure 6.4-1.

Both diffusion-node temperatures and arithmetic-node temperatures are calculated by "successive point" iteration. Each third iteration, diffusion-node temperatures which are decreasing over two time-steps are extrapolated in an attempt to accelerate convergence (refer to Section 6.2.7). Temperature convergence is examined during each time-step by checking DRLXCC and ARLXCC against the user control constants DRLXCA (for diffusion nodes) and ARLXCA (for arithmetic nodes), respectively. If temperatures have not converged with NLØØP iterations, the message "RELAXATION CRITERIA NOT MET" is printed. Control constant NLØØP is used to specify the maximum number of iterations allowed during each time-step.

VARIABLES 1 and VARIABLES 2 are performed only once for each time-step. Since this subroutine is implicit, the user must specify the time-step to be used through the control constant DTIMEI in addition to control constant TIMEND and ØUTPUT. The look ahead feature for the time-step calculation used in CNFRWD is also employed in CNBACK as are checks for maximum allowable time-step DTIMEH, maximum allowable temperature change between time-steps, DTMPCA (diffusion nodes) and ATMPCA (arithmetic nodes). The minimum time-step DTIMEL is not checked however. Damping of solutions can be achieved through the use of the control constants DAMPD and DAMPA but "radiation damping" (refer to Section 6.2.6) used by CNBACK lessens the need for the damping factors DAMPD and DAMPA.

6.4.1.4 Control Constants

Control constants ARLXCA, DRLXCA, DTIMEI, NLØØP, ØUTPUT, and TIMEND must be specified as indicated in Table 6.2-5 and as described in Section 6.2.3.2; otherwise "run" will terminate with an appropriate error message. The function of optionally specified control constants ATMPCA, BACKUP, DAMPA, DAMPD, DTIMEH, DTMPCA, and TIMEØ is described in Section 6.2.3.2.

Specification of time-step DTIMEI should be done in conjunction with control constant NLØØP which represents the maximum number of computational iterations during each time-step. Since each iterative calculation is essentially equivalent to a time-step calculation for an explicit

method, the combination of DTIMEI and NLØØP for a given time period should be less than the total number of time-steps by the explicit method for the same period. Note also that TIMEØ may be set negative. Specification of ARLXCA and DRLXCA depends upon the problem but a typical value is 0.1.

6.4.1.5 Error and Other Messages

If control constants ARLXCA, DRLXCA, DTIMEI, NLØØP, ØUTPUT and TIMEND are not specified, the following error message will be printed for each,

ARLXCA	"NØ ARLXCA"
DRLXCA	"NØ DRLXCA"
DTIMEI	"NØ DTIMEI"
NLØØP	"NØ NLØØP"
ØUTPUT	"NØ ØUTPUT INTERVAL"
TIMEND	"TRANSIENT TIME NØT SPECIFIED"

If the long pseudo-compute sequence LPCS is not specified, the error message will be,

"CNBACK REQUIRES LONG PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < (3*NND + NNA + NNB)$), the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If $CSGMIN \leq 0$, the following message will be printed,

"CSGMIN ZERØ or NEGATIVE"

If either ARLXCA or DRLXCA is not satisfied with NLØØP iterations, the following message will be printed,

"RELAXATIØN CRITERIA NØT MET"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the "run" terminating if a single check is not satisfied,

NLØØP, TIMEND, ØUTPUT, ARLXCA, DTIMEI, DRLXCA, LPCS and dynamic storage allocation.

Table 6.4-1. Basic Computational Steps for CNBACK

1. Specification of control constants. Control constants ARLXCA (if $NNA > 0$), DRLXCA (if $NND > 0$), DTIMEI, NLØØP, ØUTPUT and TIMEND ($TIMEND \geq TIMEØ$) must be specified. LPCS is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = $3*NND + NNA + NNB$ (NND = diffusion nodes, NNA = arithmetic nodes and NNB = boundary nodes).
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Time-step = DTIMEI.
4. Setting of iterative DØ loop, 1 to NLØØP.
5. Setting of source locations to zero.
6. Calling of Variables 1. (Refer to Section 6.2.2.2 for description.)
7. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
8. Diffusion-node temperature calculations, first iteration only.
 - Evaluation of q_i , C_i and G_k .
 - Damping of radiation heat transfer. (Refer to Section 6.2.5.2.)
 - Calculation of diffusion-node temperature.
 - The computational algorithm depends upon the presence of radiation heat transfer, but the method of solution is the standard implicit algorithm (refer to Section 6.2.5.2).
9. Conversion of $T_{i,k+1}$ to degrees Rankine.
10. Diffusion-node temperature calculations, successive iterations after first.
 - Repeating of step 8, except that q_i , C_i and G_k are not updated.
 - Calculation of DRLXCC.
11. Acceleration of convergence every third iteration if linear extrapolation is met (refer to Section 6.2.7).
12. Conversion of $T_{i,k+1}$ to degrees Fahrenheit.
13. Calculation of arithmetic-node temperatures, second and succeeding iterations; arithmetic-node temperatures are not calculated on the first iteration (refer to Section 6.2.5.2 for details).
14. Conversion of temperatures to degrees Rankine.
15. Checking of ARLXCA and DRLXCA for convergence and ØPEITR for output. If both ARLXCA and DRLXCA are satisfied, iterations during a time-step ceases, otherwise NLØØP iterations are performed.
16. Checking of ATMPCA and DTMPCA. If either one is not satisfied time-step is shortened, previous temperatures erased, and temperatures recalculated for shortened time-steps (refer to Section 6.2.5.2).
17. Conversion of temperatures back to degrees Fahrenheit.
18. Calling of VARIABLES 2 and checking of BACKUP (refer to Section 6.2.2.3 and 6.2.3.2).
19. Advancing of time, checking of time to print, and the printing of the output interval.
20. Calling of ØUTPUT CALLS.
21. Checking for problem end time stored in control constant TIMEND.

6.4.2 Subroutine: CNFWBK

6.4.2.1 General Comments

Subroutine CNFWBK is an implicit numerical solution routine that uses the Crank-Nicolson algorithm.^{7, 8, 12} The long pseudo-compute sequence (LPCS) is required and the nodal temperatures (both diffusion and arithmetic) are solved by "successive point" iterations. The iteration looping, convergence criteria and other control constant checks are identical to CNBACK. Time-step must be specified via control constant DTIMEI. Diffusion and arithmetic temperature calculations may be damped through the use of DAMPD and DAMPA, respectively. Thermal radiation heat transfer is uniquely "handled" via a so-called "radiation damping" (refer to Section 6.2.6), and acceleration of convergence (refer to Section 6.2.7) is also available in CNFWBK.

CNFWBK solutions which are based on a half forward differencing and a half backward differencing method tend to be more accurate than CNBACK solutions with approximately the same solution time.

6.4.2.2 Finite Difference Approximation and Computational Algorithm

The numerical solution algorithm used in subroutine CNFWBK is the Crank-Nicolson method, which is half forward differencing and half backward differencing, and may be expressed as:

$$C_i \frac{(T_{i,n+1} - T_{i,n})}{\Delta t} = \frac{1}{2} (T_{\text{forward}} + T_{\text{backward}}) \quad (6.4-2)$$

$$T_{\text{forward}} = q_{i,n} + \sum_{j=1}^p a_{ij,n} (T_{j,n} - T_{i,n}) + \sum_{j=1}^p \sigma b_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \quad (6.4-3)$$

$$T_{\text{backward}} = q_{i,n} + \sum_{j=1}^p a_{ij,n} (T_{j,n+1} - T_{i,n+1}) + \sum_{j=1}^p \sigma b_{ij,n} (T_{j,n+1}^4 - T_{i,n+1}^4) \quad (6.4-4)$$

n = nth time-step

i = 1, 2, ..., N

p = total number of nodes

$T_{j,n}; T_{j,n+1}$ = constant, $N < j \leq p$

The computational procedure for the forward-backward difference formulation must be re-iterative because of the need to solve a set of simultaneous nonlinear equations. The pattern of computation is very similar to that used in CNBACK.

Diffusion-Nodes

Diffusion node temperatures are solved by "successive point" iteration but the algorithm differs from the algorithm used in CNBACK because of the additional terms arising from the forward difference portion of the expression.

$$T_{i,k+1} = DD * T_{i,k} + DN * [Q_{sum} - (q_i)_{ave}] / G_{sum} \quad (6.4-5)$$

where,

$$Q_{sum} = Q_i + \sum_{j=1}^i a_{ij,n} T_{j,k+1} + \sum_{j=i+1}^p a_{ij,n} T_{j,k} + \sum_{j=1}^i \sigma b_{ij,n} T_{j,k+1}^4 + \sum_{j=i+1}^p \sigma b_{ij,n} T_{j,k}^4 \quad (6.4-6)$$

$$Q_i = 2 q_{i,n} + 2 \bar{C}_{i,n} T_{i,n} + \sum_{j=1}^p a_{ij,n} (T_{j,n} - T_{i,n}) \quad (6.4-7)$$

$$G_{sum} = 2 \bar{C}_{i,n} + \sum_{j=1}^p a_{ij,n} \quad (6.4-8)$$

n = nth time-step; k = kth iteration

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

DN = DAMPD (diffusion-node damping factor)

DD = 1.0 - DN

$\bar{C}_{i,n} = C_{i,n} / \Delta t$ (Δt = time-step)

$(q_i)_{ave} = \sum \sigma b_{ij,n} [(T_{i,k}^4) + (T_{i,k}^4)_2] / 2.0$, average heat loss from ith node (refer to Section 6.2.6 on radiation damping for details)

(Note that the known quantities at time-step, n, are indicated by Q_i , equation 6.4-7.)

Arithmetic Nodes

Arithmetic-node temperatures are calculated identically the same in all the SINDA numerical solution routines. Thus, refer to Section 6.3.1.2 or Section 6.2.5.2 for the finite difference algorithm.

6.4.2.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutine CNFWBK are indicated in Table 6.4-2. For a detailed step-by-step computational description, the user must examine the computer listing for CNFWBK in Appendix B, but some general computational details are given in Section 6.2.5.2. A functional flow chart of CNFWBK is shown in Figure 6.4-2.

The computational flow pattern for CNFWBK is identical to CNBACK with the only difference between the routines being the diffusion-node temperature finite-difference algorithm. On the first iteration only the source locations zeroed out and the present temperatures stored, VARIABLES 1 is called and variable C_i , impressed source q_i and variable coefficients G_i (diffusion-diffusion and diffusion-arithmetic) evaluated. All quantities which are evaluated at time, t_n , are summed in accordance with equations (6.4-6) and (6.4-8). CSGMIN is evaluated and the diffusion-node temperatures calculated; note the arithmetic-node temperatures are not calculated on the first iteration.

On the second and succeeding iterations the quantities C_i , q_i and G_k (diffusion-diffusion and diffusion-arithmetic) are not updated. Diffusion-node temperatures are calculated and DRLXCC determined. Every third iteration, if a diffusion-node temperature is converging, a linear extrapolation to accelerate convergence is performed (refer to Section 6.2.7). If arithmetic nodes are encountered, the appropriate q_i and G_k (for arithmetic nodes) are evaluated once per time-step. Arithmetic-node temperatures are calculated and ARLXCC determined.

Control constants DRLXCC and ARLXCC are checked against DRLXCA and ARLXCA, respectively each time-step; if both criteria are satisfied the iterations cease, otherwise the iterations continue NLØØP times and the message "RELAXATION CRITERIA NOT MET" is printed.

Diffusion-node and arithmetic-node temperature changes between time-steps are calculated and stored in DTMPCC and ATMPCC, respectively.

If DTMPCC > DTMPCA or if ATPCC > ATPCA, the just completed calculations are erased and the time-step shortened (refer to Section 6.2.5.2).

6.4.2.4 Control Constants

The control constants for CNFWBK are used in exactly the same way as used in CNBACK. Control constants ARLXCA, DRLXCA, DTIMEI, NLØØP, ØOUTPUT, and TIMEND must be specified as indicated in Table 6.2-5 and as described in Section 6.2.3.2; otherwise "run" will terminate with an appropriate error message. The function of optionally specified control constants ATPCA, BACKUP, DAMPA, DAMPD, DTIMEH, DTMPCA, and TIMEØ is described in Section 6.2.3.2.

Specification of time-step DTIMEI should be done in conjunction with control constant NLØØP which represents the maximum number of computational iterations during each time-step. Since each iterative calculation is essentially equivalent to a time-step calculation for an explicit method, the combination of DTIMEI and NLØØP for a given time period should be less than the total number of time-steps by the explicit method for the same time period. Note also that TIMEØ may be set negative. Specification of ARLXCA and DRLXCA depends upon the problem but a typical value is 0.1.

6.4.2.5 Error and Other Messages

If control constants ARLXCA, DRLXCA, DTIMEI, NLØØP, ØOUTPUT and TIMEND are not specified the following error message will be printed for each,

ARLXCA	"NØ ARLXCA"
DRLXCA	"NØ DRLXCA"
DTIMEI	"NØ DTIMEI"
NLØØP	"NØ NLØØP"
ØOUTPUT	"NØ ØOUTPUT INTERVAL"
TIMEND	"TRANSIENT TIME NØT SPECIFIED"

If the long pseudo-compute sequence LPCS is not specified, the error message will be,

"CNFWBK REQUIRES LØNG PSEUDØ-CØMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient (NDIM < (3* NND + NNA + NNB)), the message will be,

" _____ LOCATIONS AVAILABLE"

Note that the number presented will be negative indicating the additional storage locations required.

If $CSGMIN \leq 0$, the following message will be printed,

"CSGMIN ZERO OR NEGATIVE"

If either ARLXCA or DRLXCA is not satisfied with NLOOP iterations, the following message will be printed,

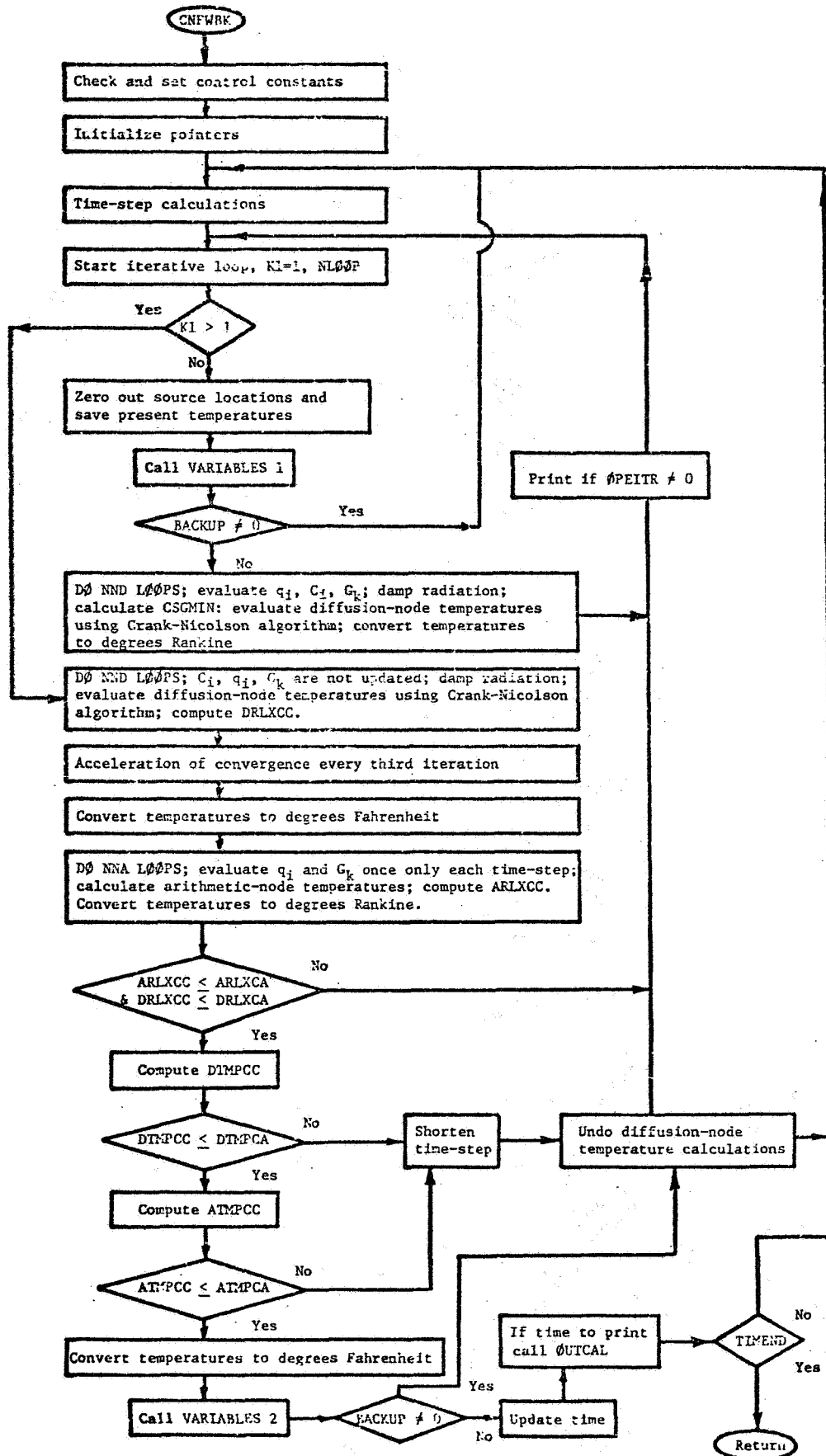
"RELAXATION CRITERIA NOT MET"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the "run" terminating if a single check is not satisfied,

NLOOP, TIMEND, OUTPUT, ARLXCA, DTIMEI, DRLXCA, LPCS and dynamic storage allocation.

Table 6.4-2. Basic Computational Steps for CNFWBK

1. Specification of control constants. Control constants ARLXCA (if $NNA > 0$), DRLXCA (if $NND > 0$), DTIMEI, NLØØP, ØOUTPUT and TIMEND ($TIMEND > TIMEO$) must be specified. LPCS is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage.
Requirements = $3*NND + NNA + NNB$ (NND = diffusion nodes, NNA = arithmetic nodes and NNB = boundary nodes)
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Time-step = DTIMEI.
4. Setting of iterative DØ loop, 1 to NLØØP.
5. Setting of source locations to zero.
6. Calling of Variables 1. (Refer to Section 6.2.2.2 for description.)
7. Checking of BACKUP. (Refer to Section 6.2.3.2 for description.)
8. Diffusion-node temperature calculations, first iteration only. Evaluation of q_i , C_i and G_k . Damping of radiation heat transfer. (Refer to Section 6.2.5.2.) Calculation of diffusion-node temperature. The computational algorithm depends upon the presence of radiation heat transfer, but the method of solution is the Crank-Nicolson algorithm (half forward and half backward, refer to Section 6.2.5.2).
9. Conversion of $T_{i,k+1}$ to °R (Rankine).
10. Diffusion-node temperature calculation, successive iterations after first. Repeating of step 8 except that q_i, C_i and G_k are not updated. Calculation of DRLXCC.
11. Acceleration of convergence every third iteration if linear extrapolation is met (refer to Section 6.2.7).
12. Conversion of $T_{i,k+1}$ to degrees Fahrenheit.
13. Calculation of arithmetic-node temperatures, second and succeeding iterations; arithmetic-node temperatures are not calculated on the first iteration (refer to Section 6.2.5.2 for details).
14. Conversion of temperatures to degrees Rankine.
15. Checking of ARLXCA and DRLXCA for convergence and ØPEITR for output. If both ARLXCA and DRLXCA are satisfied, iterations during a time-step cease, otherwise NLØØP iterations are performed.
16. Checking of ATMPCA and DTMPCA. If either one is not satisfied time-step is shortened, previous temperatures erased, and temperatures recalculated for shortened time-steps (refer to Section 6.2.5.2).
17. Conversion of temperatures back to degrees Fahrenheit.
18. Calling of VARIABLES 2 and checking of BACKUP (refer to Section 6.2.2.3 and 6.2.3.2).
19. Advancing of time, checking of time to print, and the printing of the output interval.
20. Calling of ØOUTPUT CALLS.
21. Checking for problem end-time stored in user specified control constant TIMEND.



6.4-2. Functional Flow Chart for CNFWBK

6.4.3 Subroutine: CNVARB

6.4.3.1 General Comments

Subroutine CNVARB uses an implicit finite difference algorithm that is a composition of forward-differencing and backward-differencing. The proportion of forward to backward to be used is calculated internally by using a weighting factor, β , that is dependent upon the ratio of the explicit stability criterion as stored in the control constant CSGMIN divided by the computational time-step stored in DTIMEU. The weighting factor can vary each time-step but is constrained to range, $0 \leq \beta \leq 1/2$ (refer to Section 6.2.5.2 or Section 6.4.3.2). A β of one-half yields the Crank-Nicolson half-forward and half-backward expression, whereas a β of zero yields the standard backward-difference expression.

Except for the weighting factor, β , the computational procedure and the use of the various control constants in CNVARB is essentially identical to subroutine CNFWBK.

Solution characteristics should be very similar to CNFWBK solutions with expectation that CNVARB solutions would be more optimum in terms of accuracy and solution time. Solutions are not presently available to verify or refute the expected advantages of CNVARB solutions.

6.4.3.2 Finite Difference Approximation and Computational Algorithm

The numerical solution algorithm used in subroutine CNVARB is a combination of forward-differencing and backward-differencing with the weighting of each determined by the ratio of control constants CSGMIN/DTIMEU.

The combination forward-backward differencing with weighting can be expressed as:

$$\frac{C_i}{\Delta t} (T_{i,n+1} - T_{i,n}) = \beta \left(q_{i,n} + \sum_{j=1}^P a_{ij,n} (T_{j,n} - T_{i,n}) \right) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \\ + (1.0 - \beta) \left(q_{i,n} + \sum_{j=1}^P a_{ij,n} (T_{j,n+1} - T_{i,n+1}) \right) + \sum_{j=1}^P \sigma b_{ij,n} (T_{j,n+1}^4 - T_{i,n+1}^4) \quad (6.4-9)$$

$$i = 1, 2, \dots, N$$

$$n = \text{nth time-step}$$

$$\beta = \text{weighting factor } (0 < \beta \leq 1/2)$$

$$T_{j,n}; T_{j,n+1} = \text{constant, } N < j \leq P$$

If equation (6.4-9) is multiplied by 2.0 and the known quantities (at time-step, n) and the unknown quantities (at time-step, n+1) separated, the algorithm used in CNVARB may be obtained readily.

Diffusion Nodes

Diffusion-node temperatures are solved by "successive point" iteration. The finite difference iterative form as used in CNVARB can be found by multiplying equation (6.4-9) by 2.0 and by using appropriate time-step, n, and iteration, k subscripts.

$$T_{i,k+1} = DD* T_{i,k} + DN* [Q_{sum} - (q_i)_{ave}] / G_{sum} \quad (6.4-10)$$

where, $Q_i = 2 q_{i,n} + 2 \bar{C}_{i,n} T_{i,n} + \beta' \left(\sum_{j=1}^P a_{ij,n} (T_{j,n} - T_{i,n}) + \sum_{j=1}^P ob_{ij,n} (T_{j,n}^4 - T_{i,n}^4) \right)$ (6.4-11)

$$Q_{sum} = Q_i + (2.0 - \beta') \left(\sum_{j=1}^i G_{ij,n} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,n} T_{j,k} \right) \quad (6.4-12)$$

$$G_{sum} = 2 \bar{C}_{i,n} + (2.0 - \beta') \sum_{j=1}^P a_{ij,n} \quad (6.4-13)$$

$$G_{ij,n} = a_{ij,n} + ob_{ij,n} T_{j,\ell}^3 \quad (6.4-14)$$

($\ell = k$, if $j \geq i$ and $\ell = k+1$, if $j < i$)

$$(q_i)_{ave} = \frac{(2.0 - \beta')}{2} \sum_{j=1}^P ob_{ij,n} [(T_{i,k}^4)_1 + (T_{i,k}^4)_2] \quad (6.4-15)$$

average heat loss from the ith node, called radiation damping (refer to Section 6.2.6 for details)

= 0, if radiation is not present

$\beta' = 2.0 * CSGMIN / DTIMEU$ (range allowed, $0 \leq \beta' \leq 1.0$, note $\beta' = 2\beta$)

n = nth time-step; k = kth iteration

C_i, q_i, a_{ij}, b_{ij} = optionally specified (refer to Tables 6.2-1 - 6.2-4)

$\bar{C}_{i,n} = C_{i,n} / \Delta t$

i = 1, 2, ..., NND

$T_{j,n}; T_{j,k}$ = constant, $(NND + NNA) < j \leq p$ (p is the total number of nodes and NNA is the number of arithmetic nodes)

Arithmetic Nodes

Arithmetic nodes are calculated in the same manner in all the SINDA numerical solution routines. For the finite difference algorithm refer to either Section 6.3.1.2 or Section 6.2.5.2.

6.4.3.3 Comments on the Computational Procedure

The important steps of the computational procedure used in subroutine CNVARB are indicated in Table 6.4-3. For a detailed step-by-step computational description, the user must examine the computer listing for CNVARB in Appendix B, but some general computational details are given in Section 6.2.5.2. A functional flow chart of CNVARB is shown in Figure 6.4-3.

The computational flow pattern for CNVARB is very similar to CNFWBK or CNBACK; the slight difference is shown in the flow chart of Figure 6.4-3. The basic difference between CNVARB and the other two implicit routines is the use of a variable beta, β' , which is calculated internally by the routine. Thus, the updating of the variable capacitance C_i , the impressed source q_i and the variable coefficients (a_{ij} for conduction and σ_{ij} for radiation) during the first iteration and the subsequent calculation of diffusion-node temperatures in subsequent iterations are identical to CNFWBK except for the finite difference algorithm. Use of the various control constants and checks are identical to CNFWBK.

6.4.3.4 Control Constants

Control constants for CNVARB are used in exactly the same way as used in CNFWBK. Control constant ARLXCA, DRLXCA, DTIMEI, NL $\emptyset\emptyset$ P, \emptyset UTPUT, and TIMEND must be specified as indicated in Table 6.2-5 and as described in Section 6.2.3.2; otherwise "run" will terminate with an appropriate error message. The function of optionally specified control constants ATMPCA, BACKUP DAMPA, DAMPD, DTIMEH, DTMPCA and TIME \emptyset is described in Section 6.2.3.2.

6.4.3.5 Error and Other Messages

If control constants ARLXCA, DRLXCA, DTIMEI, NL $\emptyset\emptyset$ P, \emptyset UTPUT and TIMEND are not specified, the following error message will be printed for each,

APLXCA	"NO ARLXCA"
DRLXCA	"NO DRLXCA"
DTIMEI	"NO DTIMEI"
NLØØP	"NO NLØØP"
ØUTPUT	"NO OUTPUT INTERVAL"
TIMEND	"TRANSIENT TIME NOT SPECIFIED"

If the long pseudo-compute sequence LPCS is not specified, the error message will be,

"CNVARB REQUIRES LONG PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < (3 * NND + NNA + NNB)$), the error message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number presented will be negative indicating the additional storage locations required.

If $CSGMIN \leq 0$, the following message will be printed,

"CSGMIN ZERO or NEGATIVE"

If either ARLXCA or DRLXCA is not satisfied with NLØØP iterations, the following message will be printed,

"RELAXATION CRITERIA NOT MET"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following sequence with the "run" terminating if a single check is not satisfied,

NLØØP, TIMEND, ØUTPUT, ARLXCA, LPCS and dynamic storage allocation.

Table 6.4-3. Basic Computational Steps for CNVARB

1. Specification of control constants. Control constants ARLXCA (if $NNA > 0$), DRLXCA (if $NND > 0$), DTIMEI, NLØØP, ØUTPUT and TIMEND ($TIMEND > TIMEØ$) must be specified. LPCS is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = $3*NND + NNA + NNB$ (NND = diffusion nodes, NNA = arithmetic nodes and NNB = boundary nodes).
3. Setting and/or calculation of time-step, Δt . (Refer to Section 6.2.4 for detailed procedure.) Time-step = DTIMEI.
4. Setting of iterative DØ loop, 1 to NLØØP.
5. Setting of source locations to zero.
6. Calling of Variables 1 (refer to Section 6.2.2.2 for description).
7. Checking of BACKUP (refer to Section 6.2.3.2 for description).
8. Diffusion-node temperature calculations, first iteration only.
 - Checking of stable stability criteria.
 - Calculation of weighting factor $\beta' = 2.0*CSGMIN/DTIMEU$. ($0 \leq \beta' \leq 1.0$)
 - Conversion of temperatures to degrees Rankine.
 - Damping of radiation heat transfer (refer to Section 6.2.5.2).
 - Calculation of diffusion-node temperatures using forward-backward algorithm with variable beta (β').
 - Calculation of DRLXCC.
9. Diffusion-node temperature calculations, successive iterations after first. Repeating of step 8 except that q_i , C_i and G_k are not updated. Calculation of DRLXCC.
10. Acceleration of convergence every third iteration if linear extrapolation criterion is met (refer to Section 6.2.7).
11. Conversion of $T_{i,k+1}$ to degrees Fahrenheit.
12. Calculation of arithmetic-node temperatures every iteration (refer to Section 6.2.5.2 for details).
13. Conversion of temperatures to degrees Rankine.
14. Checking of ARLXCA and DRLXCA for convergence and ØPEITER for output. If both ARLXCA and DRLXCA are satisfied, iterations during a time-step cease, otherwise NLØØP iterations are performed.
15. Checking of ATMPCA and DTMPCA. If either one is not satisfied time-step is shortened, previous temperatures erased, and temperatures recalculated for shortened time-steps (refer to Section 6.2.5.2).
16. Conversion of temperatures back to degrees Fahrenheit.
17. Calling of VARIABLES 2 and checking of BACKUP (refer to Section 6.2.2.3 and 6.2.3.2).
18. Advancing of time, checking of time to print, and the printing of the output interval.
19. Calling of ØUTPUT CALLS.
20. Checking for problem end time stored in user specified control constant TIMEND.

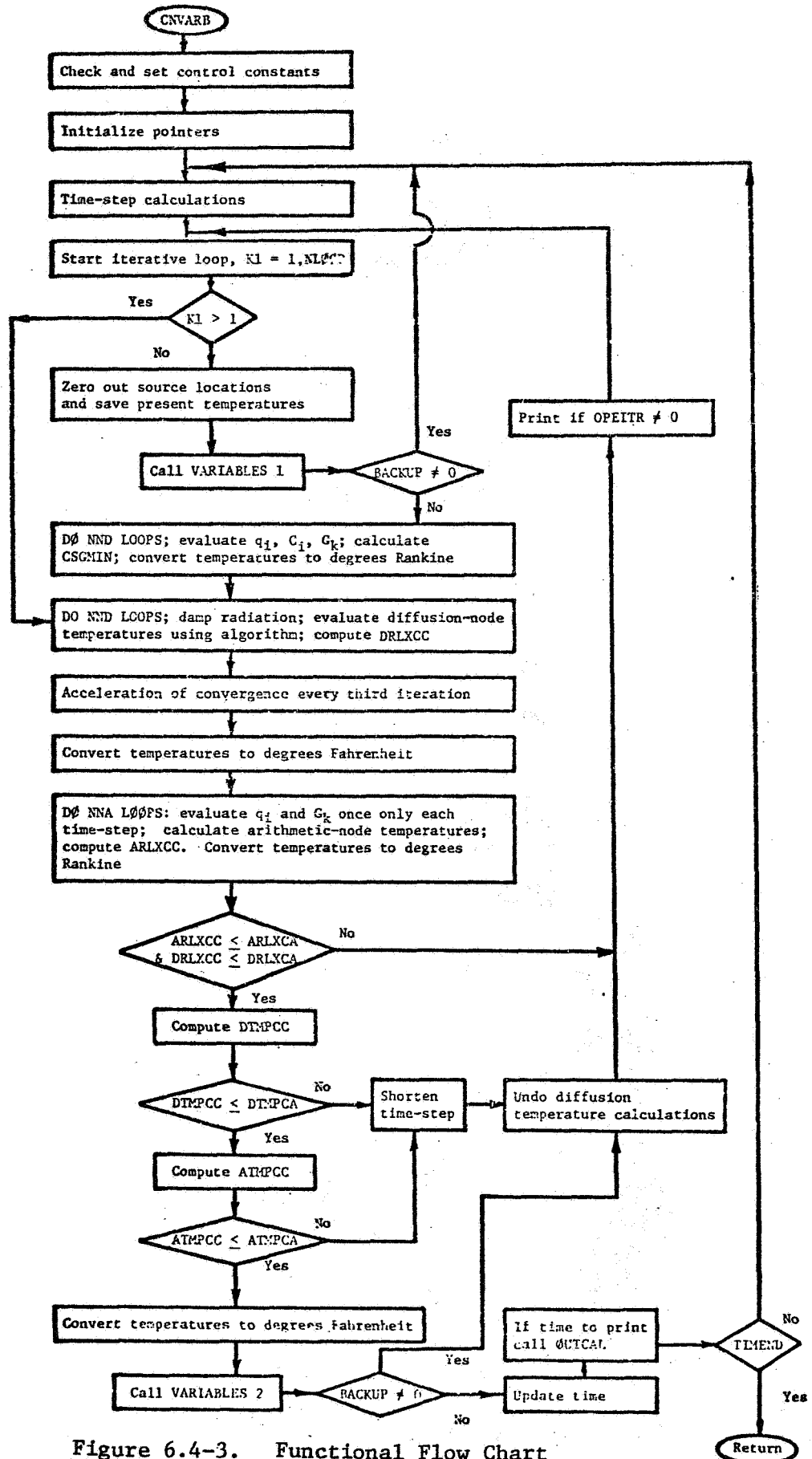


Figure 6.4-3. Functional Flow Chart for CNVARB

6.5 Steady State Numerical Solution Routines

SINDA steady state numerical solution routines number three. These steady state routines are identified as follows:

- CINDSS Block iterative method
Requires short pseudo-compute sequence (SPCS)
- CINDSL Successive point iterative method
Requires long pseudo-compute sequence (LPCS)
- CINDSM Modified CINDSL for radiation-dominated problems
Requires long pseudo-compute sequence (LPCS)

A detailed description of steady state routines is presented in the pages to follow with liberal reference to materials presented in Section 6.2. A brief description of these routines follows.

CINDSS which uses the short pseudo-compute sequence (SPCS) was the first steady state routine developed for SINDA (via CINDA and CINDA-3G); as a result, some of the features contained in subsequent steady state routines are not used in CINDSS. If a transient analysis is to be performed following a steady state analysis, CINDSS must be used with a transient routine that also requires SPCS. The "block" iterative method (refer to Section 5.2.3) used by CINDSS should lend itself to some types of problems which are highly nonlinear with terms such as $G_{ij} (T_j^4 - T_i^4)$. With "block" iteration, both T_j and T_i are changed simultaneously. Solution convergence is based upon a temperature relaxation criterion stored in DRLXCA for diffusion nodes and ARLXCA for arithmetic nodes.

CINDSL requires the long pseudo-compute sequence (LPCS) and uses the "successive point" iteration method (refer to Section 5.2.3). Any transient analysis routine coupled with CINDSL must require LPCS. Solution time for CINDSL is less than CINDSS; as a result, it is used more often than CINDSS. A major problem with CINDSL is that a highly nonlinear problem can present convergence difficulties unless considerable amount of damping is used. For example, a radiation-dominated problem contains many $\sigma_{ij} (T_j^4 - T_i^4)$. With "successive point" iteration, T_j may be updated and T_i not for a given conductor; as a result, the resultant heat flow calculation could present difficulties because of large change in values. CINDSL has the acceleration of convergence feature, whereas CINDSS does not.

Solution convergence is based upon temperature relaxation criterion stored in DRLXCA for diffusion nodes and ARLXCA for arithmetic nodes.

CINDSM is the latest addition to the SINDA library of steady state routines. CINDSM requires the long pseudo-compute sequence and uses "successive point" iteration. The routine was specifically developed to solve radiation-dominated problems. Solution convergence is based upon system energy criterion stored in BALENG.

6.5.1 Subroutine: CINDSS

6.5.1.1 General Comments

Subroutine CINDSS is a steady state routine that requires the short pseudo-compute sequence (SPCS) and ignores the capacitance values of diffusion nodes to calculate steady state temperatures. Diffusion nodes are solved by a "block" iterative method as discussed in Section 6.5.2.3, whereas arithmetic nodes are solved by a "successive point" iterative method also discussed in Section 6.5.2.3. For steady state solutions diffusion nodes are not necessary; as a matter of fact, solutions will be achieved more quickly if all diffusion nodes are specified as arithmetic. The use of diffusion nodes in a steady state solution allows for the direct use of the transient model.

A series of steady state solutions at various points in a time period can be accomplished by specifying control constants TIMEN and ØUTPUT. ØUTPUT is used both as the output interval and the computational interval. The instructions with the appropriate call are made in VARIABLES 1 to modify boundary conditions with time.

The CINDSS call can be followed by a call to one of the transient solution subroutines which has the same short pseudo-compute sequence requirements such as CNFRWD. In this manner the steady state solution becomes the initial conditions for the transient analysis. It is important to remember that control constants specified for the steady state routine will be used by the transient routine unless initialized to the desired values. Since CINDSS utilizes control constants TIMEND and ØUTPUT for the steady state-transient problem, the user must specify their values in the execution block after the steady state call and prior to the transient analysis call. CINDSS does not utilize the acceleration of convergence feature as discussed in Section 6.2.7.

Solution convergence is based upon a temperature relaxation criterion stored in control constants DRLXCA for diffusion nodes and ARLXCA for arithmetic nodes. Normally, identical values are specified for both DRLXCA and ARLXCA. Sufficient information is not presently available to indicate different values for DRLXCA and ARLXCA. A method to indicate the accuracy of the "converged" temperatures is not presently available. It

should also be noted that "converged" temperatures could have large system energy unbalance.

6.5.1.2 Finite Difference Approximation and Computational Algorithm

The steady state heat balance equation at the i th node may be readily expressed as,

$$q_i + \sum_{j=1}^P a_{ij} (T_j - T_i) + \sum_{j=1}^P \sigma b_{ij} (T_j^4 - T_i^4) = 0 \quad (6.5-1)$$

$$i = 1, 2, \dots, N$$

$$T_j = \text{constant}, N < j \leq p$$

Equation (6.5-1) represents a set of nonlinear algebraic equations to be solved simultaneously. Since CINDSS solves temperature of nodes specified as diffusion (nodes with capacitance even though a steady state solution is desired) by the "block" iteration method and temperatures of nodes specified as arithmetic (no capacitance) by the "successive point" iteration method, two successive approximation algorithms are used.

Diffusion Nodes (if any)

$$T_{i,k+1} = DD * T_{i,k} + \frac{DN * (q_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}} \quad (6.5-2)$$

where, $k = k$ th iteration; $i = 1, 2, \dots, NND$ (number of diffusion nodes)

q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

$T_{j,k} = \text{constant}, (NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,k}^2 + T_{i,k}^2) (T_{j,k} + T_{i,k})$$

$DN \equiv DAMPD$ (diffusion node damping factor)

$$DD = 1.0 - DN$$

Arithmetic Nodes (if any)

$$T_{i,k+1} = AD * T_{i,k} + \frac{AN * (q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}}$$

where, $k = k$ th iteration; $i = (NND + 1), (NND + 2), \dots, (NND + NNA)$

$q_i, a_{ij}, b_{ij} =$ optionally specified (refer to Tables 6.2-1 - 6.2-4)

$T_{j,k} =$ constant, $(NND + NNA) < j \leq p$ (NNA is the number of arithmetic nodes and p is the total number of nodes)

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,\ell}^2 + T_{i,\ell}^2) (T_{j,\ell} + T_{i,\ell})$$

($\ell = k$, if $j \geq i$ and $\ell = k+1$, if $j < i$)

AN \equiv DAMPA (arithmetic node damping factor)

AD = 1.0 - AN

6.5.1.3 Comments on the Computational Procedure

The important steps of the computational procedure used in the steady state subroutine CINDSS are indicated in Table 6.5-1. For a detailed procedural description, the user must examine the computer listing for CINDSS in Appendix C, but some general computational details are given in Section 6.2.5.3. A functional flow chart of CINDSS is shown in Figure 6.5-1. The user is required to specify the maximum number of iterations to be performed via control constant NLØØP and the diffusion-node temperature change relaxation criteria DRLXCA and the arithmetic-node temperature change criteria ARLXCA. The iterations continue until either NLØØP is satisfied or both DRLXCA and ARLXCA are satisfied. If DRLXCA and ARLXCA are not satisfied with NLØØP iterations, an appropriate message is printed. VARIABLES 1 and ØUTPUT CALLS are performed at the start and VARIABLES 2 and ØUTPUT CALLS are performed upon completion. Control constants DAMPD for diffusion nodes and DAMPA for arithmetic nodes are so-called damping factors which are multipliers of the "new" temperatures; the factor 1.0 - DAMPD (or 1.0 - DAMPA) is a multiplier for the "old" temperatures. This weighting of "old" and "new" temperatures is useful for damping oscillations due to nonlinearities. For nonlinear systems, the damping factors are specified to be less than one. If not specified, the damping factor is set to 1.0. As a point of interest, it appears that if a linear system is to be solved, the convergence could be accelerated by using the damping factor greater than one. The diffusion nodes receive a "block" iteration, whereas the arithmetic nodes receive a "successive point" iteration; acceleration features are not utilized.

6.5.1.4 Control Constants

Control constant NLØØP must be specified and control constants ARLXCA and DRLXCA must be specified if $NNA > 0$ and $NND > 0$, respectively; otherwise "run" will terminate with an appropriate error message. Control constants DAMPA and DAMPD may be optionally specified among others. Control constant characteristics are tabulated in Table 6.2-5 and description of these control constant is presented in Section 6.2.3.2. Specification of NLØØP is dependent upon the values of ARLXCA and DRLXCA and thus the accuracy of solution. Since the type of problem will influence accuracy, it appears that a trial and error procedure is the only practical way of determining realistic control constant values.

6.5.1.5 Error and Other Messages

If control constants ARLXCA, DRLXCA and NLØØP are not specified, the following error message will be printed for each,

ARLXCA	"NØ ARLXCA"
DRLXCA	"NØ DRLXCA"
NLØØP	"NØ NLØØP"

If the short pseudo-compute sequence SPCS is not specified, the error message will be,

"CINDSS REQUIRES SHØRT PSEUDO-CØMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient ($NDIM < NND$) will be,

"_____ LØCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If both temperature change relaxation criteria ARLXCA and DRLXCA are not met with NLØØP iterations, the message will be,

"ITERATION CØUNT EXCEEDED, LØØPCT = _____"

Checks on the control constants, the pseudo-compute sequence, and the dynamic storage allocation are made in the following order with the "run" terminating if a single check is not satisfied.

NLØØP, ARLXCA, DRLXCA, SPCS, and dynamic storage allocation.

Table 6.5.1. Basic Computational Steps for CINDSS

1. Specification of control constants. Control constants ARLXCA (if NNA > 0), DRLXCA (if NND > 0) and NLØØP must be specified. SPCS is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = NND (NND = diffusion nodes).
3. Setting of TIMEN for first iteration and succeeding iterations.
 $TIMEN = TIMEØ$, first iteration
 $TIMEN = TIMEØ + ØUTPUT$, succeeding iterations
4. Setting of iterative loop for all nodes, $k1 = 1$, NLØØP
5. Setting of source locations to zero.
6. Calling of VARIABLES 1 (refer to Section 6.2.2.2 for description).
7. Calculation of diffusion-node temperatures by "block" iteration if NND > 0 (refer to sections 6.2.5.3 and 6.5.1.2).

$$T_{i,k+1} = DD * T_{i,k} + \frac{DN * (q_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}}$$

$$DN = DAMPD \text{ and } DD = 1.0 - DN$$

8. Calculation of DRLXCC.
9. Calculation of arithmetic-node temperatures by "successive point" iteration if NNA > 0 (refer to Sections 6.2.5.3 and 6.5.1.2).

$$T_{i,k+1} = AD * T_{i,k} + \frac{AN * (q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}}$$

$$AN = DAMPA$$

$$AD = 1.0 - DAMPA$$

10. Calculation of ARLXCC.
11. Checking of DRLXCC and ARLXCC against the relaxation criteria DRLXCA and ARLXCA, respectively, for convergence. If both ARLXCA and DRLXCA are satisfied, iterations cease, otherwise NLØØP iterations are performed.
12. Calculation of system energy balance which is stored in ENGBAL.
13. Call VARIABLES 2 and ØUTCAL, print ENGBAL and LØØPCT.
14. Check if TIMEND = TIMEN.

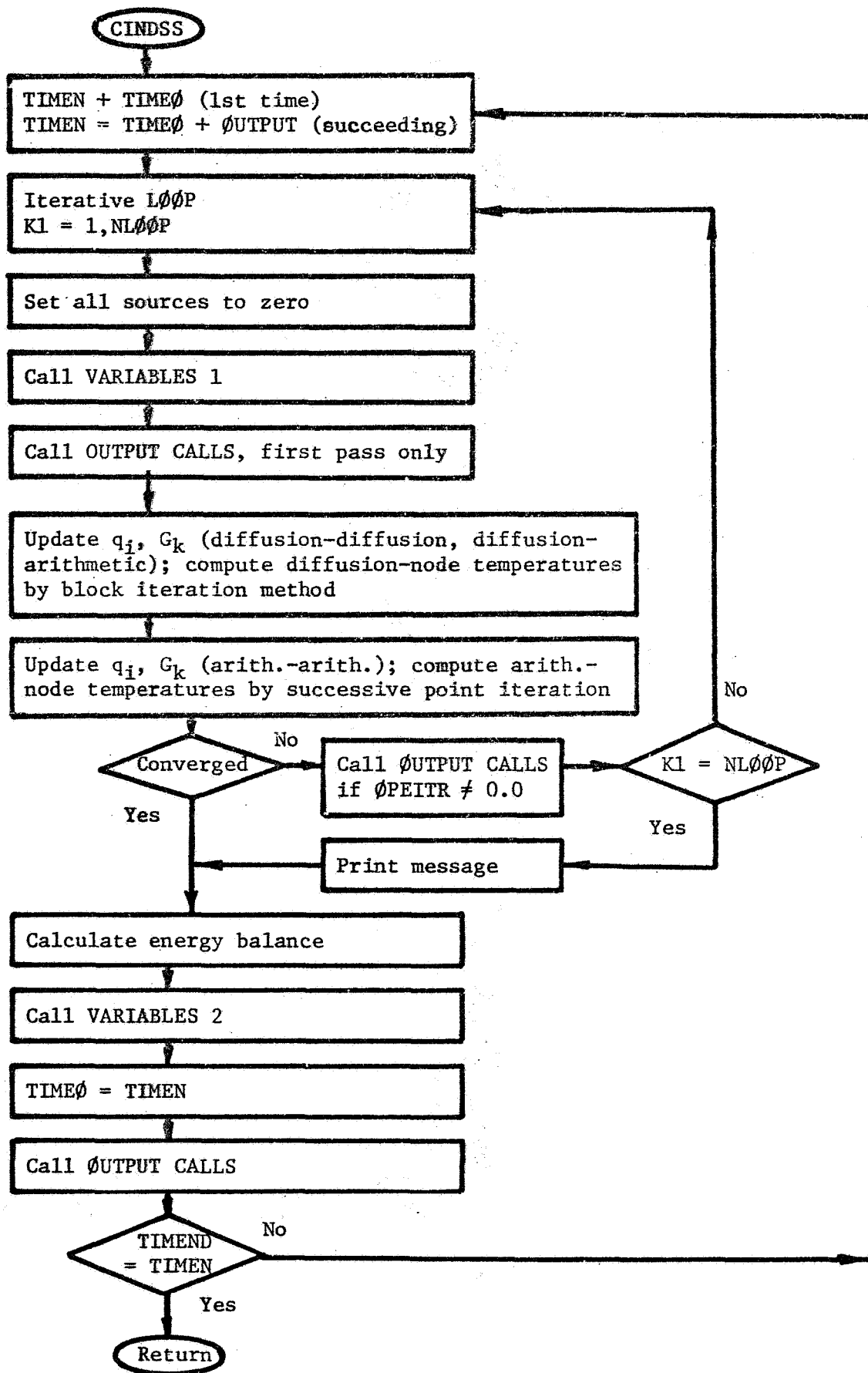


Figure 6.5-1. Functional Flow Chart for CINDSS

6.5.2 Subroutine: CINDSL

6.5.2.1 General Comments

Subroutine CINDSL is a steady state routine that requires the long pseudo-compute sequence (LPCS). Both diffusion- and arithmetic-node temperatures are calculated by a "successive point" iteration computational technique. Every third iteration a linear extrapolation is performed to accelerate convergence. CINDSL generally yields significantly faster solutions than CINDSS, but nonlinear problems such as those with radiation heat transfer can pose considerable convergence difficulties unless a large amount of damping (low values of DAMPA and DAMPD) is imposed.

A series of steady state solutions at various points in time can be generated by specifying control constants TIMEND and ØUTPUT. ØUTPUT is used both as the output interval and the computation interval; this requires appropriate calls in VARIABLES 1 to modify boundary conditions with time.

CINDSL can be followed by a call to one of the transient numerical solution routines which have the same LPCS requirements. Used in this manner the steady state solutions become the initial conditions for the transient analysis. Note that since CINDSL utilizes control constants TIMEND and ØUTPUT for the coupled steady state-transient problem, the user must specify the values of TIMEND and ØUTPUT in the execution block after the steady state call and prior to the transient analysis call.

Solution convergence is based upon a temperature relaxation criterion stored in control constants DRLXCA for diffusion nodes and ARLXCA for arithmetic nodes. Normally, identical values are specified for both DRLXCA and ARLXCA for lack of anything better. The damping factors DAMPD for diffusion nodes and DAMPA for arithmetic nodes are merely multipliers of "new" temperatures and the factor $1.0 - \text{DAMPD}$ (or $1.0 - \text{DAMPA}$) is a multiplier of the "old" temperatures. Normally, these damping factors are specified to be less than 1.0, but for a linear system the convergence probably could be accelerated by using a damping factor greater than one.

6.5.2.2 Finite Difference Approximation and Computational Algorithm

The set of steady state heat balance equations,

$$q_i + \sum_{j=1}^P a_{ij} (T_j - T_i) + \sum_{j=1}^P \sigma b_{ij} (T_j^4 - T_i^4) = 0$$

$$i = 1, 2, \dots, N$$

$$T_j = \text{constant } N < j \leq P$$

is solved by a re-iterative scheme called a "successive point" iterative method here. Both diffusion-node and arithmetic-node temperatures are solved in this manner. The only difference between the two algorithms is that control constant DAMPD is used with diffusion nodes and control constant DAMPA is used with arithmetic nodes.

Diffusion Nodes (if any)

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{(q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}} \quad (6.5-4)$$

where, $i = 1, 2, \dots, NND$; $k = k$ th iteration

q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

$T_{j,k}$ = constant, $(NND + NNA) < j \leq P$ (NNA is the number of arithmetic nodes and P is the total number of nodes)

DN = DAMPD (diffusion-node damping factor)

DD = 1.0 - DN

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$$

($\ell = k$ if $j \geq i$ and $\ell = k+1$ if $j < i$)

Arithmetic Nodes (if any)

$$T_{i,k+1} = AD * T_{i,k} + AN * \frac{(q_{i,k} + \sum_{j=1}^P G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}} \quad (6.5-5)$$

where, $i = (NND + 1), (NND + 2), \dots, (NND + NNA)$

q_i, a_{ij}, b_{ij} = may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

$T_{j,k}$ = constant $(NND + NNA) < j \leq P$ (NNA is the number of arithmetic nodes and P is the total number of nodes)

AN = DAMPA (arithmetic-node damping factor)

AD = 1.0 - AN

$$G_{ij,k} = a_{ij,k} + \sigma b_{ij,k} (T_{j,\ell}^2 + T_{i,k}^2) (T_{j,\ell} + T_{i,k})$$

($\ell = k$ if $j \geq i$ and $\ell = k+1$ if $j < i$)

6.5.2.3 Comments on the Computational Procedure

The important steps of the computational procedure used in the steady state subroutine CINDSL are indicated in Table 6.5-2. For a detailed procedural description, the user must examine the computer listing for CINDSL in Appendix C, but some general computational details are given in Section 6.2.5.3. A functional flow chart of CINDSL is shown in Figure 6.5-2.

The computational pattern of CINDSL is very similar to CINDSS with the differences being that CINDSL uses the long pseudo-compute sequence, whereas CINDSS uses the short pseudo-compute sequence, and that CINDSL contains the acceleration convergence feature, whereas CINDSS does not. The user is required to specify the maximum number of iterations to be performed via control constant NLØØP and the diffusion-node temperature change relaxation criteria DRLXCA and the arithmetic-node temperature change relaxation criteria ARLXCA. The iterations continue until either NLØØP is satisfied or both DRLXCA and ARLXCA are satisfied. If DRLXCA and ARLXCA are not satisfied with NLØØP, an appropriate message is printed. Acceleration of convergence is performed every third iteration if a temperature is converging over two time-steps.

6.5.2.4 Control Constants

Control constant NLØØP must be specified and control constants ARLXCA and DRLXCA must be specified if NNA > 0 and NND > 0, respectively; otherwise "run" will terminate with an appropriate error message. Control constants DAMPA and DAMPD may be optionally specified among others. Control constant characteristics are tabulated in Table 6.2-5 and description of these control constants is presented in Section 6.2.3.2. Specification of NLØØP is dependent upon the values of ARLXCA and DRLXCA and thus the accuracy of the solution. Since the type of problem will influence accuracy, it appears that a trial and error procedure is the only practical way of determining realistic control constant values.

6.5.2.5 Error and Other Messages

If control constants ARLXCA, DRLXCA and NLØØP are not specified, the following error message will be printed for each,

ARLXCA	"NØ ARLXCA"
DRLXCA	"NØ DRLXCA"
NLØØP	"NØ NLØØP"

If the long pseudo-compute sequence LPCS is not specified, the error message will be,

"CINDSL REQUIRES LONG PSEUDO-COMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient, (NDIM < 2* (NNA + NND)), the message will be,

"_____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

"LØØPCT = _____ and ENGBAL = _____"

If both temperature change relaxation criteria, ARLXCA and DRLXCA, are not met with NLØØP iterations, the message will be,

"ITERATION COUNT EXCEEDED, LØØPCT = _____"

Checks on the control constants, the pseudo-compute sequence, and the dynamic storage allocation are made in the following order with the "run" terminating if a single check is not satisfied.

NLØØP, ARLXCA, DRLXCA, LPCS, and dynamic storage allocation.

Table 6.5.2 Basic Computational Steps for CINDSL

1. Specification of control constants. Control constants ARLXCA (if $NNA > 0$), DRLXCA (if $NND > 0$) and NLØØP must be specified. LPCS is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirements = $2 * (NND + NNA)$ (NND = diffusion nodes and NNA = arithmetic nodes).
3. Setting of TIMEN for first and succeeding iterations.

$$\text{TIMEN} = \text{TIMEØ}, \text{ first iteration}$$

$$\text{TIMEN} = \text{TIMEØ} + \text{ØUTPUT}, \text{ succeeding iterations}$$
4. Setting of iterative loop for all nodes, $k1 = 1$, NLØØP.
5. Setting of source locations to zero.
6. Calling of VARIABLES 1 (refer to Section 6.2.2.2 for description).
7. Calculation of diffusion-node temperatures by "block" iteration if $NND > 0$ (refer to Section 6.2.5.2 and 6.5.1.2).

$$T_{i,k+1} = DD * T_{i,k} + \frac{DN * (q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}}$$

$$DN = \text{DAMPD} \text{ and } DD = 1.0 - DN$$

8. Calculation of DRLXCC.
9. Calculation of arithmetic-node temperatures by "successive point" iteration if $NNA > 0$ (refer to Sections 6.2.5.3 and 6.5.1.2).

$$T_{i,k+1} = AD * T_{i,k} + \frac{AN * (q_{i,k} + \sum_{j=1}^i G_{ij,k} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,k} T_{j,k})}{\sum_{j=1}^P G_{ij,k}}$$

10. Calculation of ARLXCC.
11. Checking of DRLXCC and ARLXCC against the relaxation criteria DRLXCA and ARLXCA, respectively, for convergence. If both ARLXCA and DRLXCA are satisfied, iterations cease, otherwise NLOOP iterations are performed.
12. Acceleration of convergence each third iteration, if linear extrapolation criterion is met (refer to Section 6.2.7).
13. Calculation of system energy balance which is stored in ENGBAL.
14. Call VARIABLES 2 and ØUTCAL, print ENGBAL and LØØPCT.
15. Check if $\text{TIMEND} = \text{TIMEN}$.

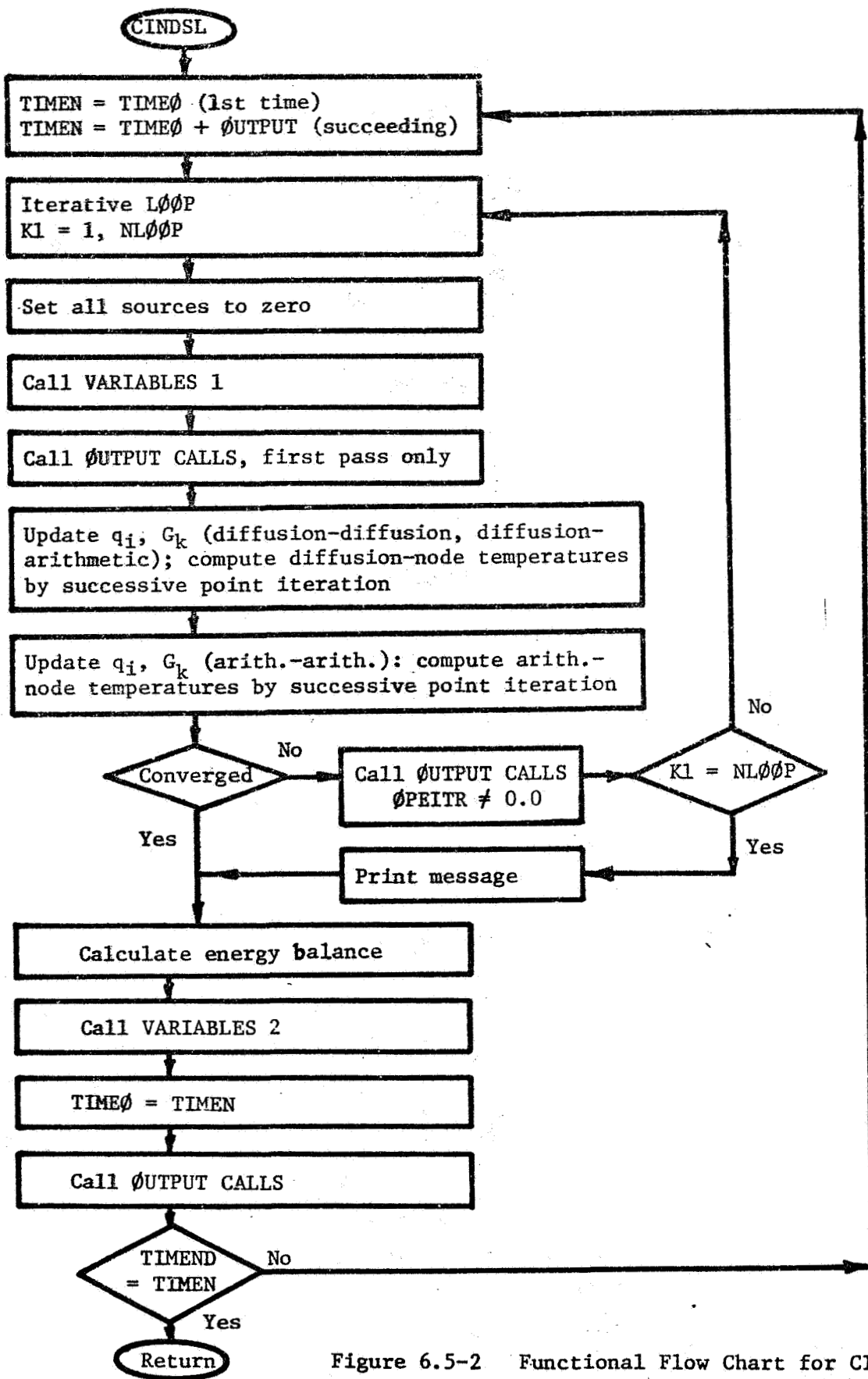


Figure 6.5-2 Functional Flow Chart for CINDSL

6.5.3 Subroutine: CINDSM

6.5.3.1 General Comments

Subroutine CINDSM is a steady state routine specifically generated for radiation dominated problems. CINDSM requires the long pseudo-compute sequence (LPCS) and is considerably different from CINDSL. CINDSM is based on the use of pseudo linear equations which are the result of linearizing the radiation conductors. These equations are solved by using the "successive point" method with LAXFAC iterations. Updating of the properties as well as the linearized conductors occur outside of the iterative loops. Temperature convergence is based on a criterion that is continually tightened until either the NLØØP iterations or the system energy balance criterion stored in BALENG has been satisfied.

The acceleration of convergence by linear extrapolation as used in CINDSM is essentially the same as used in the other SINDA numerical solution routines, but in lieu of limiting the extrapolation by an allowable slope value (refer to Section 6.2.7) the maximum temperature change of the network on the last iteration is used as the allowable value.

Information available at this time indicates that each problem appears to have an optimum combination of NLØØP, DAMPD, and LAXFAC values. An NLØØP of 100, a DAMPD of 0.5 and a LAXFAC of 10 has been successfully applied to spacecraft problems with radiation domination, but the solution time is rather long.

6.5.3.2 Finite Difference Approximation and Computational Algorithm

The set of steady state heat balance equations,

$$q_i + \sum_{j=1}^P a_{ij} (T_j - T_i) + \sum_{j=1}^P \sigma b_{ij} (T_j^4 - T_i^4) = 0$$

$$i = 1, 2, \dots, N$$
$$T_j = \text{constant}, N < j \leq P$$

is solved by a re-iterative "successive point" method after linearization. Linearization is achieved by letting $\sigma b_{ij} (T_j^4 - T_i^4) = G_r (T_j - T_i)$ with $G_r = \sigma b_{ij} (T_j^2 + T_i^2)(T_j + T_i)$. This yields

$$q_i + \sum_{j=1}^P a_{ij} (T_j - T_i) + \sum_{j=1}^P G_r (T_j - T_i) = 0 \quad (6.5-6)$$

Diffusion and Arithmetic Nodes

No distinction is made between diffusion and arithmetic nodes. As a result, the following algorithm applies to both types of nodes,

$$T_{i,k} = DD * T_{i,k} + DN * \frac{(q_{i,L} + \sum_{j=1}^i G_{ij,L} T_{j,k+1} + \sum_{j=i+1}^P G_{ij,L} T_{j,k})}{\sum_{j=1}^P G_{ij,L}} \quad (6.5-7)$$

where, $i = 1, 2, \dots, (NND + NNA)$; $p =$ total number of nodes

$k =$ kth iteration

$L =$ before each LAXFAC iterative loop

$T_{j,k} =$ constant, $(NND + NNA) < j \leq p$

$DN =$ DAMPD (diffusion-node damping factor; DAMPA is not used)

$DD = 1.0 - DAMPD$

$G_{ij,L} = a_{ij,L} + \sigma b_{ij,L} (T_{j,L}^2 + T_{i,L}^2)(T_{j,L} + T_{i,L})$

($G_{ij,L}$ is updated once before each LAXFAC iterative loop)

$NNA =$ number of arithmetic nodes

$NND =$ number of diffusion nodes

$q_i, a_{ij}, b_{ij} =$ may be optionally specified (refer to Tables 6.2-1 - 6.2-4)

6.5.3.3 Comments on the Computational Procedure

A detailed step-by-step computational procedure as used in the steady state routine CINDSM is presented in Table 6.5-3. For a more detailed procedural description, the user must examine the computer listing in Appendix C. A functional flow chart that is compatible with the step-by-step description of Table 6.5-3 is shown in Figure 6.5-3.

CINDSM is considerably different from either CINDSS or CINDSL because of the use of a variable convergence criterion which is internally updated. Overall, from a total system basis, control constants $NL\emptyset\emptyset P$ and $BALENG$ are the ultimate criteria.

It should be particularly noted here that unlike CINDSS or CINDSL, which use both DAMPA and DAMPD, CINDSM uses only DAMPD. The reason for this is that CINDSM does not treat the nodal types as diffusion or arithmetic.

6.5.3.4 Control Constant

Control constants BALENG, LAXFAC AND NLØØP must be specified; otherwise the "run" will terminate with an appropriate error message. Control constant DAMPD may be optionally specified among others. Control constant characteristics are tabulated in Table 6.2.5 and description of these control constants is presented in Section 6.2.3.2. Specification of BALENG, LAXFAC and NLOOP appears to be a trial and error procedure.

6.5.3.5 Error and Other Messages

If control constants BALENG, LAXFAC, and NLØØP are not specified, the following error message will be printed for each,

BALENG	"NØ BALENG"
LAXFAC	"NØ LAXFAC"
NLØØP	"NØ NLØØP"

If the long pseudo-compute sequence LPCS is not specified, the error message will be,

"CINDSM REQUIRES LONG PSEUDØ-CØMPUTE SEQUENCE"

If the dynamic storage allocation is not sufficient, (NDIM < (3* NNA + 3* NND + NGT)), the message will be,

" _____ LOCATIONS AVAILABLE"

Note that the number printed will be negative indicating the additional storage locations required.

If either NLØØP iterations has been made or if ENGBAL < BALENG, the following message is printed,

"LØØPCT = _____ and ENGBAL _____"

Checks on the control constants, the pseudo-compute sequence and the dynamic storage allocation are made in the following order with the "run" terminating if a single check is not satisfied,

NLØØP, LPCS, BALENG, LAXFAC and dynamic storage allocation.

Table 6.5.3. Basic Computational Steps for CINDSM

1. Specification of control constants. Control constants BALENG, LAXFAC and NLØØP must be specified. The long pseudo-compute sequence (LPCS) is required. (Refer to Table 6.2-5 for values and Section 6.2.3.2 for description.)
2. Sufficiency check on dynamic storage. Requirement = 3* (NND + NNA) + NGT (NND = diffusion nodes, NNA = arithmetic nodes and NGT = total number of conductors).
3. Setting of TIMEN for the first and succeeding iterations.

TIMEN = TIMEØ, first iteration
 TIMEN = TIMEØ + ØOUTPUT, succeeding iterations

4. Constants used in CINDSM

NLAX = NLØØP/LAXFAC (both NLØØP and LAXFAC are specified by the user)
 RELAX = .05 (initial value used in CINDSM as the allowable temperature change)
 DELXXX = .05/NLAX (a number used in reducing RELAX for a tighter criterion)
 XXXDUM = .001 (a value of RELAX used in CINDSM for a tighter criterion)
 = .001/5 (a subsequent value of RELAX for a tighter criterion)
 DAMP = DAMPD (damping factor for all nodes; DAMPA is not used)

5. Updating of variables and linearization of radiation .

Variable q_i and G_k are evaluated by calling subroutine NØNLIN.

Linearization means that the radiation exchange expressed as $\sigma b_{ij} (T_j^4 - T_i^4)$. Normally, G_{ij} would be updated each iteration as done in CINDSS or CINDSL, but in CINDSM G_{ij} is not updated within the DØ-LØØP ($k_1 = 1, LAXFAC$) but is updated outside of the loop.

6. Iterative DØ-LØØP ($k_1 = 1, LAXFAC$) is established.

Temperatures of all nodes are calculated by "successive point" iteration with no damping.

$$T_{i,k+1} = \frac{q_i + \sum_{j=1}^i G_{ij} T_{j,k+1} + \sum_{j=i+1}^P G_{ij} T_{j,k}}{\sum_{j=1}^P G_{ij}} \quad (6.5-5)$$

where, $G_{ij} = a_{ij} + \sigma b_{ij} (T_j^2 + T_i^2) (T_j + T_i)$ (q_i and G_{ij} are not updated during the LAXFAC iterations)

Check on temperature convergence. Temperatures have converged if,

$$|T_{i,k+1} - T_{i,k}|_{\max} \leq \text{RELAX} (= .05)$$

If temperatures have converged, the computation goes out of the iteration loop to step (7).

Table 6.5.3. (continued)

Every third iteration, acceleration of convergence is attempted if linear extrapolation criterion is met (refer to Section 6.2.7).

Iteration ceases if LAXFAC iterations have been performed or if the temperatures have converged.

7. Check on NLAX iterations.

If in step (6) the number of iterations, LOOPCT \geq NLAX, the computational procedures go to step (9). However, in step (6) if the number of iterations LOOPCT < NLAX, then a set of temperature calculations is made using "successive point" method with a damping factor and no iterations.

$$T_{i,k+1} = DD * T_{i,k} + DN * \frac{(q_i + \sum_{j=1}^i G_{ij} T_{j,k+1} + \sum_{j=i+1}^P G_{ij} T_{j,k})}{\sum_{j=1}^P G_{ij}}$$

where, DN = DAMPD (diffusion node damping factor; note DAMPA is not used)

G_{ij} = constant

Allowable temperature change criterion RELAX is reduced to,

$$RELAX = .05 - (.05/NLAX)$$

and computational procedure goes to step (5).

8. Repetition of steps (5) through (7) except for temperature convergence criterion.

Temperatures have converged if,

$$|T_{i,k+1} - T_{i,k}|_{\max} \leq RELAX (= .05 - .05/NLAX)$$

9. Assuming step (7) has been satisfied, LOOPCT is checked against NLOOP.

If LOOPCT \geq NLOOP, the computation proceeds to step (12).

If LOOPCT < NLOOP computation proceeds to step (10).

10. Reduce RELAX to .001.

11. Check on temperature convergence.

If $|T_{i,k+1} - T_{i,k}| \leq RELAX (= .001)$ go to step (12).

$|T_{i,k+1} - T_{i,k}| > RELAX (= .001)$, LAXFAC is reduced to

$$LAXFAC = NLOOP - LOOPCT,$$

and steps (5) through (11) are repeated.

12. Compute system energy balance and store in control constant ENGBAL.

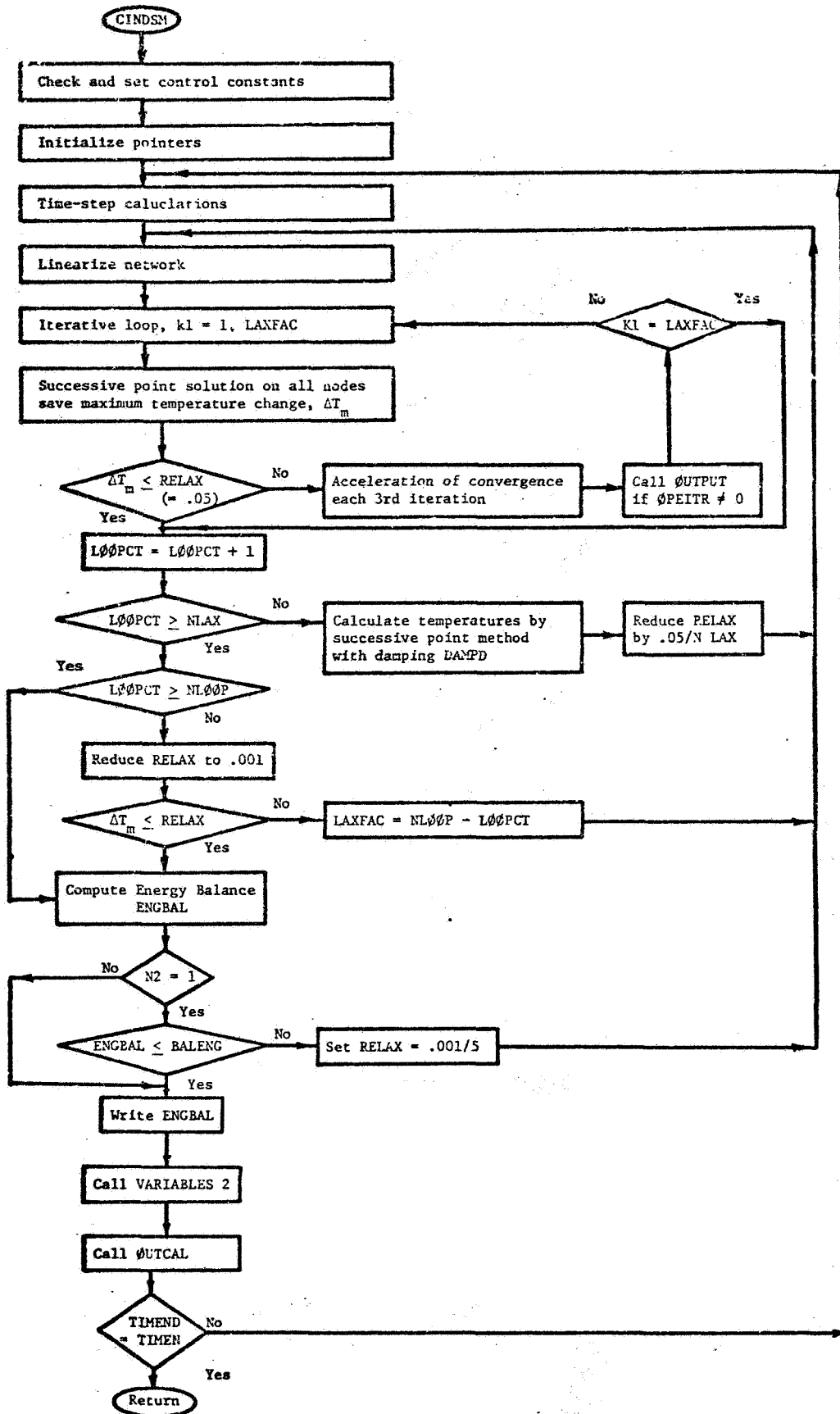
13. If LOOPCT \geq LAXFAC (original user input value), go to step (15)

14. If LOOPCT \leq LAXFAC (original user input value), ENGBAL is checked against BALENG.

If ENGBAL \leq BALENG, go to step (14)

If ENGBAL > BALENG, RELAX is set to, RELAX = .001/5, and steps (5) through (14) are repeated with the new RELAX values.

15. Print ENGBAL; call VARIABLES 2; call OUTCAL; check if TIMEND = TIMEØ.



6.5-3. Functional Flow Chart for CINDSM

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A.

COMPUTER LISTINGS OF SINDA EXPLICIT SOLUTION ROUTINES

	Page
CNFRWD	A-2
CNDRDL	A-13
CNFAST	A-24
CNEXPN	A-33
CNDUFR	A-44
CNQUIK	A-55

Q1W FOR,* CNFRWD,CNFRWD
UNIVAC 1108 FORTRAN V ATHENA VERSION/ 131K-10D CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:37

21 FEB 71

SUBROUTINE CNFRWD ENTRY POINT 003522

STORAGE USED (BLOCK, NAME, LENGTH)

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0000 *COJST+TEMP 000072
0002 *SIMPLE VAR 000050
0004 *ARRAYS 000000
0005 *BLANK 000000
0006 TITLE 000001
0007 TEMP 000001
0010 CAP 000001
0011 SOURCE 000001
0012 COND 000001
0013 PC1 000001
0014 PC2 000001
0015 KONST 000001
0016 ARRAY 000001
0017 FIXCON 000001
0020 XSPACE 000003
0021 DIMENS 000010

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EXTERNAL REFERENCES (BLOCK, NAME)

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0022 VARBL1
0023 DID1WM
0024 PLYAWM
0025 D2D1WM
0026 VARBL2
0027 OUTCAL
0030 EXIT
0031 NERR2$
0032 NWDU$
0033 NI02$
0034 NER10$

```

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

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0016 R 000000 A 0010 R 000000 C 0002 R 000015 CKW
0002 R 000022 C2 0002 R 000041 DD 0002 R 000037 DELTA
0002 R 000034 GV 0002 R 000030 G1 0002 R 000031 G2
0002 I 000002 IE 0002 I 000043 JJ1 0002 I 000044 JJ2
0015 I 000000 K 0017 I 000000 KON 0002 I 000045 L
0002 I 000011 LE 0002 I 000036 LEA 0002 I 000025 LG
0021 I 000007 LSQ2 0002 I 000026 LTA 0002 I 000005 LI
0020 I 000000 NDM 0021 I 000003 NLA 0002 I 000001 NNA
0021 I 000000 NND 0021 I 000002 NNT 0013 I 000000 NSG1
0002 I 000016 NIYPE 0020 N00R02 NX 0002 R 000000 PASS
0002 R 000023 Q1 0002 R 000024 Q2 0002 R 000046 SUBC
0002 R 000014 TCGM 0002 R 000027 TM 0002 R 000047 SUMCV
0002 R 000032 T1 0002 R 000033 T2 0020 R 000002 X
0017 R 000000 CON 0014 I 000000 NSG2
0002 R 000040 DN 0006 R 000000 H 0002 I 000005 NAT
0002 I 000012 J1 0002 I 000017 LA 0002 I 000020 LK
0002 I 000042 LAX 0021 I 000006 LSG1 0021 I 000004 NCT
0002 I 000004 Y 0002 I 000013 J2 0002 I 000001 NNC
0002 R 000000 G 0002 R 000000 G 0020 I 000001 NTH
0002 R 000000 G 0002 R 000000 G 0002 R 000035 QDOT
0002 R 000000 T 0007 R 000000 T
0002 R 000010 TSUM 0002 R 000006 TSTEP
0001 000125 10L 0015 R 000000 XK

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CNFRWD,CNFRWD

0001	003500	1000L	0001	000327	1005L	0001	000350	1010L	0001	003336	10116	0001	000366	1012L
0001	000421	1015L	0001	000427	1020L	0001	000467	1025L	0001	000513	1030L	0001	000534	1032L
0001	000572	1035L	0001	000600	1040L	0001	000643	1045L	0001	000701	105L	0001	002123	105L
0001	002134	120L	0001	002142	125L	0001	002150	130L	0001	002500	135L	0001	003246	140L
0001	003251	145L	0001	000134	15L	0001	003316	165L	0001	003327	175L	0001	003360	180L
0001	003371	185L	0001	003420	190L	0001	003423	195L	0001	000665	1998L	0001	000670	1998L
0001	000206	2006	0001	000673	2000L	0001	001235	2005L	0001	001242	2007L	0001	001262	2010L
0001	001265	2015L	0001	001303	2017L	0001	001337	2020L	0001	001345	2025L	0001	001405	2030L
0001	001412	2032L	0001	001435	2035L	0001	001440	2040L	0001	001461	2042L	0001	001520	2045L
0001	001526	2050L	0001	001571	2055L	0001	001617	2060L	0001	001623	2065L	0001	000225	2126
0001	000256	2326	0001	000146	25L	0001	001627	2998L	0001	001655	2999L	0001	000161	30L
0001	001660	3000L	0001	002566	3005L	0001	002573	3007L	0001	002613	3010L	0001	002616	3015L
0001	002634	3017L	0001	002670	3020L	0001	002676	3025L	0001	002736	3030L	0001	002743	3032L
0001	002766	3035L	0001	002771	3040L	0001	003012	3042L	0001	003051	3045L	0001	003057	3050L
0001	003122	3055L	0001	003150	3060L	0001	003154	3065L	0001	003160	3998L	0001	003206	3999L
0001	003211	4000L	0001	000740	4005L	0001	000745	4010L	0001	000746	4012L	0001	000765	4015L
0001	000766	4017L	0001	001003	4020L	0001	001017	4022L	0001	001034	4025L	0001	001042	4030L
0001	001100	4035L	0001	001114	4037L	0001	001131	4040L	0001	000237	45L	0001	001157	4988L
0001	001143	4999L	0001	000113	5L	0001	001146	50L	0001	001146	5000L	0001	002272	5005L
0001	002277	5010L	0001	002300	5012L	0001	002317	5015L	0001	002320	5017L	0001	002335	5020L
0001	002351	5022L	0001	002366	5025L	0001	002374	5030L	0001	002432	5035L	0001	002446	5037L
0001	002004	5046	0001	002463	5040L	0001	002075	5356	0001	001715	55L	0001	002177	5736
0001	002471	5998L	0001	002475	5999L	0001	001720	60L	0001	002500	6000L	0001	002220	6026
0001	001767	65L	0001	001776	85L	0000	000000	885F	0000	000011	886F	0000	000016	897F
0000	000023	888F	0000	000030	889F	0001	002025	90L	0001	002044	95L	0001	003442	945L
0001	003450	996L	0001	003456	997L	0001	003464	998L	0001	003473	999L			

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00101 1* SUBROUTINE CNFRWD
00101 2* EXPLICIT FORWARD DIFFERENCING EXECUTION SUBROUTINE FOR SINDA F-V
00101 3* THE SHORT PSEUDO-COMPUTE SEQUENCE IS REQUIRED
00103 4* INCLUDE COMMLIST
00104 4* COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/G(1) /COND/G(1)
00105 4* COMMON /PC1/NS01(1) /PC2/NS02(1) /KONST/K(1) /ARRAY/A(1)
00106 4* COMMON /FIXCON/KON(1) /XSPACE/NDIM,NTH,X(1)
00107 4* COMMON /DIMENS/ NND,NMA,NNT,NGT,NCT,NAT,LS01,LS02
00110 4* DIMENSION CON(1),XK(1),NX(1)
00111 4* EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
00112 4* END
00113 5* INCLUDE DEFF,LIST
C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR,EXPLICIT (CSGFAC)
C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)
C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
C CONTROL CONSTANT 17 IS RESERVED FOR THE C/56 MINIMUM (CSGMIN)

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00113 5* C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 5* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 5* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 5* C CC21 IS FOR THE MINIMUM ALLOWED TIME STEP (DTIME1)
00113 5* C CC22 CONTAINS THE C/SG MAXIMUM (CSGMAX)
00113 5* C CC23 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
00113 5* C CC24 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
00113 5* C CC25 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 5* C CC26 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 5* C CC27 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 5* C CC28 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 5* C CC29 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 5* C CC30 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 5* C CC31 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGBAL)
00113 5* C CC32 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 5* C CC33 CONTAINS THE NUCOPY SWITCH FOR MATRIX USERS (NUCOPY)
00113 5* C CC34 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 5* C CC35 CONTAINS RELATIVE NODE NUMBER OF DTMPCC
00113 5* C CC36 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC37 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 5* C CC38-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 5* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (R-S-T-U-VTEST)
00113 5* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (LAXFAC)
00113 5* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM
00113 5* C CC50 IS NOT USED AT PRESENT
00114 5* C END
00115 6* IF(CON(4).LT.1.0) CON(4) = 1.0
00117 7* IF(KON(5).LE.0) KON(5) = 1
00121 8* IF(CON(6).LE.0.) CON(6) = 1.E+8
00125 9* IF(CON(8).LE.0.) CON(8) = 1.E+8
00125 10* IF(CON(9).LE.0.) CON(9) = 1.0
00127 11* IF(CON(11).LE.0.) CON(11) = 1.E+8
00131 12* IF(CON(18).LE.0.) GO TO 999
00133 13* IF(CON(19).LE.0.) CON(19) = 1.E+8
00135 14* IF(KON(31).NE.0) GO TO 995
00137 15* PASS = -1.0
00140 16* NNC = NIND+NNA
00141 17* IE = NTH
00142 18* NLA = NDIM
00143 19* NTH = NTH+NNC
00144 20* NDIM = NDIM+NNC
00144 21* CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00145 22* I = NLA+NNC
00146 23* IF(I.LT.0) GO TO 998
00150 24* LI = NND+1
00151 25* TSTEP = CON(18)
00152 26* TPRINT = CON(13)
00152 27* INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00153 28* TSUM = 0.0
00153 29* 5 DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00154 30* IF(CON(13)+CON(18).LE.CON(3)) GO TO 10
00154 31* DONT EXCEED IT
00156 32* CON(18) = CON(13)-CON(13)
00156 33* 10 IS THE TIME STEP LARGER THAN ALLOWED
00157 34* IF(TSTEP.LE.CON(8)) GO TO 15
00161 35* TSTEP = CON(8)
00161 36* DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
00162 37* 15 IF(TSUM+TSTEP-CON(18)) 25,30,20
00162 38* DONT EXCEED IT
00165 39* 20 TSTEP = CON(18)-TSUM

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40* 00166 60 TO 30
41* 00167 C DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
42* 00168 25 IF(TSUM+2.0*TSTEP.LE.CON(18)) GO TO 30
43* 00169 C APPROACH THE OUTPUT INTERVAL GRADUALLY
44* 00170 TSTEP = (CON(18)-TSUM)/2.0
45* 00171 C STORE DELTA TIME STEP IN THE CONSTANTS
46* 00172 30 CON(2) = TSTEP
47* 00173 C IS THE TIME STEP USED LESS THAN THE TIME STEP ALLOWED
48* 00174 IF(TSTEP.LT.CON(21)) GO TO 997
49* 00175 C CALCULATE THE NEW TIME
50* 00176 CON(1) = TPRINT+TSUM+TSTEP
51* 00177 C COMPUTE THE MEAN TIME BETWEEN ITERATIONS
52* 00178 CON(14) = (CON(1)+CON(13))/2.0
53* 00179 C ZERO OUT ALL SOURCE LOCATIONS AND EXTRA LOCATIONS
54* 00180 DO 35 I = 1,NND
55* 00181 LE = IE+I
56* 00182 X(LE) = 0.0
57* 00183 Q(I) = 0.0
58* 00184 35 CONTINUE
59* 00185 C SHIFT THE ARITHMETIC TEMPERATURES INTO THE EXTRA LOCATIONS
60* 00186 IF(NNA.LE.0) GO TO 45
61* 00187 DO 40 I = L1,NNC
62* 00188 Q(I) = 0.0
63* 00189 LE = IE+I
64* 00190 X(LE) = T(I)
65* 00191 40 CONTINUE
66* 00192 45 KON(12) = 0
67* 00193 CALL VARBL1
68* 00194 IF(KON(12).NE.0) GO TO 10
69* 00195 J1 = 0
70* 00196 J2 = 1
71* 00197 TCGM = 0.0
72* 00198 CPM = 1.E+8
73* 00199 C CALCULATE Q SUM AND 6 SUM
74* 00200 DO 65 I = 1,NND
75* 00201 LE = IE+I
76* 00202 INCLUDE VARC,LIST
77* 00203 IF(FLD(1,1,NS01(J1+1)).EQ.0) GO TO 2000
78* 00204 NTYPE = FLD(0,5,NS02(J2))
79* 00205 LA = FLD(5,17,NS02(J2))
80* 00206 LK = FLD(22,14,NS02(J2))
81* 00207 GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
82* 00208 1005 CALL DDI1WM(T(I),A(LA),XK(LK),C(I))
83* 00209 GO TO 1999
84* 00210 1010 CALL DDI1WM(T(I),A(LA),XK(LK),C1)
85* 00211 1012 J2 = J2+1
86* 00212 LA = FLD(5,17,NS02(J2))
87* 00213 LK = FLD(22,14,NS02(J2))
88* 00214 CALL DDI1WM(T(I),A(LA),XK(LK),C2)
89* 00215 GO TO 1998
90* 00216 1015 C1 = XK(LK)*XK(LA)
91* 00217 GO TO 1012
92* 00218 1020 CALL DDI1WM(T(I),A(LA),XK(LK),C1)
93* 00219 J2 = J2+1
94* 00220 LA = FLD(5,17,NS02(J2))
95* 00221 LK = FLD(22,14,NS02(J2))
96* 00222 C2 = XK(LK)*XK(LA)
97* 00223 GO TO 1998
98* 00224 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
99* 00225 GO TO 1999
00226 76*
00227 71*
00228 72*
00229 73*
00230 74*
00231 75*
00232 76*
00233 76*
00234 76*
00235 76*
00236 76*
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00255 76*
00256 76*
00257 76*
00260 76*
00261 76*
00262 76*
00263 76*
00264 76*
00265 76*

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*NEW
*NEW
*NEW
**-3

*NEW
*NEW
**-2

*NEW
*NEW
**-2

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00266	76*	1030	CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)	*NEW
00267	76*	1032	J2 = J2+1	*NEW
00270	76*		LA = FLD(5,17,NSG2(J2))	***-2
00271	76*		LK = FLD(22,14,NSG2(J2))	
00272	76*		CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)	
00273	76*		GO TO 1998	
00274	76*	1035	C1 = XK(LK)*XK(LA)	
00275	76*		GO TO 1032	
00276	76*	1040	CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)	*NEW
00277	76*		J2 = J2+1	*NEW
00300	76*		LA = FLD(5,17,NSG2(J2))	***-2
00301	76*		LK = FLD(22,14,NSG2(J2))	
00302	76*		C2 = XK(LK)*XK(LA)	
00303	76*		GO TO 1998	
00304	76*	1045	CALL D2DIWM(T(I),CON(14),A(LA),XK(LK),C1))	
00305	76*		GO TO 1999	
00306	76*	1998	C11 = C1+C2	
00307	76*	1999	J2 = J2+1	
00310	76*	2000	CONTINUE	
00311	76*		END	
00312	77*		INCLUDE VARO,LIST	
00313	77*		IF(FLD(4,1,NSO1(J1+1)),EQ,0) GO TO 5000	
00315	77*		NTYPE = FLD(0,5,NSO2(J2))	*NEW
00316	77*		LA = FLD(5,17,NSG2(J2))	*NEW
00317	77*		LK = FLD(22,14,NSG2(J2))	***-3
00320	77*		GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE	
00321	77*	4005	Q11 = XK(LK)+Q11	
00322	77*		GO TO 4999	
00323	77*	4010	Q1 = 0.0	
00324	77*	4012	CALL D1DIWM(T(I),A(LA),XK(LK),Q2)	
00325	77*		GO TO 4998	
00326	77*	4015	Q1 = 0.0	
00327	77*	4017	CALL D1DIWM(CON(14),A(LA),XK(LK),Q2)	
00330	77*		GO TO 4998	
00331	77*	4020	CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)	
00332	77*	4022	J2 = J2+1	
00333	77*		LA = FLD(5,17,NSG2(J2))	
00334	77*		LK = FLD(22,14,NSG2(J2))	
00335	77*		GO TO 4017	
00336	77*	4025	Q1 = XK(LK)*XK(LA)	
00337	77*		GO TO 4022	
00340	77*	4030	CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)	
00341	77*		J2 = J2+1	
00342	77*		LA = FLD(5,17,NSG2(J2))	
00343	77*		LK = FLD(22,14,NSG2(J2))	
00344	77*		Q2 = XK(LK)*XK(LA)	
00345	77*		GO TO 4998	
00346	77*	4035	CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)	
00347	77*	4037	J2 = J2+1	
00350	77*		LA = FLD(5,17,NSG2(J2))	
00351	77*		LK = FLD(22,14,NSG2(J2))	
00352	77*		GO TO 4012	
00353	77*	4040	Q1 = XK(LK)*XK(LA)	
00354	77*		GO TO 4037	
00355	77*	4998	Q(I) = Q1+Q2+Q(I)	
00356	77*	4999	J2 = J2+1	
00357	77*	5000	CONTINUE	
00360	77*		END	
00361	78*	50	J1 = J1+1	
00362	79*	L6	= FLD(5,16,NSO1(J1))	

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00362	80*	C	CHECK FOR LAST CONDUCTOR	*NEW
00363	81*		IF(LG,EO,0) GO TO 85	*NEW
00365	82*		LTA = FLD(22,14,NSQ1(J1))	*NEW
00366	83*		INCLUDE VARG,LIST	**--3
00366	84*	C	CHECK FOR RADIATION CONDUCTOR	
00367	84*		IF(FLD(2,1,NSQ1(J1)),EO,0) GO TO 3000	
00371	84*		NTYPE = FLD(10,5,NSQ2(J2))	
00372	84*		LA = FLD(5,17,NSQ2(J2))	
00373	84*		LK = FLD(22,14,NSQ2(J2))	
00374	84*		GO TO (2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,	
00374	84*		2060,2065), NTYPE	
00375	84*		TM = ((I)+T(LTA))/2.0	
00376	84*		2007 CALL D2D1WM(TM,A(LA),XK(LK),6(LG))	
00377	84*		GO TO 2999	
00400	84*		2010 TM = T(I)	
00401	84*		GO TO 2007	
00402	84*		2015 CALL D2D1WM(T(I),A(LA),XK(LK),61)	
00403	84*		2017 J2 = J2+1	
00404	84*		LA = FLD(5,17,NSQ2(J2))	
00405	84*		LK = FLD(22,14,NSQ2(J2))	
00406	84*		CALL D2D1WM(T(LTA),A(LA),XK(LK),62)	
00407	84*		GO TO 2998	
00410	84*		2020 G1 = XK(LK)*XK(LA)	*NEW
00411	84*		GO TO 2017	*NEW
00412	84*		2025 CALL D2D1WM(T(I),A(LA),XK(LK),61)	**--2
00413	84*		J2 = J2+1	
00414	84*		LA = FLD(5,17,NSQ2(J2))	
00415	84*		LK = FLD(22,14,NSQ2(J2))	
00416	84*		G2 = XK(LK)*XK(LA)	*NEW
00417	84*		GO TO 2998	**--2
00420	84*		2030 TM = ((I)+T(LTA))/2.0	
00421	84*		2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))	*NEW
00422	84*		GO TO 2999	*NEW
00423	84*		2035 TM = T(I)	**--2
00424	84*		GO TO 2032	
00425	84*		2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),61)	
00426	84*		2042 J2 = J2+1	
00427	84*		LA = FLD(5,17,NSQ2(J2))	
00430	84*		LK = FLD(22,14,NSQ2(J2))	
00431	84*		CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)	*NEW
00432	84*		GO TO 2998	*NEW
00433	84*		2045 G1 = XK(LK)*XK(LA)	**--2
00434	84*		GO TO 2042	
00435	84*		2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),61)	*NEW
00436	84*		J2 = J2+1	*NEW
00437	84*		LA = FLD(5,17,NSQ2(J2))	**--2
00440	84*		LK = FLD(22,14,NSQ2(J2))	
00441	84*		G2 = XK(LK)*XK(LA)	
00442	84*		GO TO 2998	
00443	84*		2055 TM = ((I)+T(LTA))/2.0	
00444	84*		CALL D2D1WM(TM,CON(14),A(LA),XK(LK),6(LG))	
00445	84*		GO TO 2999	
00446	84*		2060 TM = T(LTA)	
00447	84*		GO TO 2007	
00450	84*		2065 TM = T(LTA)	
00451	84*		GO TO 2032	
00452	84*		G(LG) = 1./((1./G1)+1./G2)	
00453	84*		IF(FLD(1,1,NSQ1(J1)),EO,1) G(LG) = G1*G2	
00455	84*		2999 J2 = J2+1	
00456	.84*		3000 CONTINUE	

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00457 84*
00460 85*
00462 86*
00463 87*
00464 88*
00465 89*
00466 90*
00466 91*
00467 92*
00470 93*
00470 94*
00471 95*
00471 96*
00472 97*
00472 98*
00474 99*
00475 100*
00476 101*
00476 102*
00477 103*
00501 104*
00501 105*
00503 106*
00506 107*
00506 108*
00507 109*
00510 110*
00512 111*
00513 112*
00513 113*
00514 114*
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00515 116*
00515 117*
00516 118*
00520 119*
00521 120*
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00522 122*
00523 123*
00524 124*
00526 125*
00527 126*
00530 127*
00530 128*
00532 129*
00532 130*
00534 131*
00537 132*
00540 133*
00541 134*
00543 135*
00545 136*
00546 137*
00547 138*
00550 139*
00551 140*
00551 141*
00552 142*
00552 143*

END
IF (FLOI3,1,NSO1(J1)).EQ.0) GO TO 55
T1 = T(I)+460.0
T2 = I(LTA)+460.0
GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
GO TO 60
55 GV = G(LG)
C OBTAIN THE Q RATE THRU THE CONDUCTOR
60 QDOT = GV*(T(LTA)-T(I))
Q(I) = Q(I)+QDOT
C SAVE SUMMATION OF CONDUCTORS
X(LE) = X(LE)+GV
CHECK FOR ADJOINING DIFFUSION NODE
IF (LTA.GT.IJND.OR.FLD(21,1,NSO1(J1)).EQ.1) GO TO 65
C SAVE SUMMATION OF CONDUCTORS FOR ADJOINING NODE
LEA = IE+LTA
X(LEA) = X(LEA)+GV
Q(LTA) = Q(LTA)-QDOT
C CHECK FOR LAST CONDUCTOR
65 IF (IJSO1(J1).GT.0) GO TO 50
85 CONTINUE
C OBTAIN NEW DIFFUSION TEMPERATURES, DTMPC AND CSGMIN
DO 100 I = 1,MND
LE = IE+I
C CALCULATE C/SK MINIMUM
T1 = C(I)/X(LE)
IF (T1.GE.CKM) GO TO 90
CKM = T1
KON(35) = I
C COMPUTE NEW TEMPERATURES USING CALCULATED SOURCE TERMS
90 T1 = TSTEP*Q(I)/C(I)
T2 = ARS(T1)
C CALCULATE THE ABSOLUTE VALUE TEMPERATURE CHANGE
SAVE THE LARGEST TEMPERATURE CHANGE
IF (TCGM.GE.T2) GO TO 95
TCGM = T2
KON(36) = I
C STORE THE TEMPERATURES
95 X(LE) = T(I)
T(I) = T(I)+T1
100 CONTINUE
CON(17) = CKM
DELTA = CKM/CON(4)
IF (CKM.LE.0.0) GO TO 996
C CHECK FOR FIRST PASS
IF (PASS.GT.0.0) GO TO 115
C UNDO THE TEMPERATURE CALCULATIONS
105 DO 110 I = 1,MNC
LE = IE+I
T(I) = X(LE)
110 CONTINUE
IF (PASS.GT.0.0) GO TO 15
PASS = 1.0
CON(1) = TPRINT
CON(2) = 0.0
TSTEP = UELTA*0.95
GO TO 195
C IS THE TIME STEP USED LESS THAN THE TIME STEP CALCULATED
115 IF (TSTEP.LE.DELTA) GO TO 130
C COMPUTE THE TIME STEP

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00554 144* TSTEP = DELTA*0.95
00555 145* GO TO 105
00556 146* 120 TSTEP = 0.95*TSTEP*CON(6)/TCGM
00557 147* GO TO 105
00560 148* 125 TSTEP = 0.95*TSTEP*CON(11)/TCGM
00561 149* GO TO 105
00561 150* C SEE IF THE TEMPERATURE CHANGE WAS TOO LARGE
00562 151* C IF(TCGM.GT.CON(6)) GO TO 120
00562 152* C STORE THE MAXIMUM DIFFUSION TEMPERATURE CHANGE
00564 153* CON(15) = TCGM
00564 154* C CHECK TO SEE IF THERE ARE ANY ARITHMETIC NODES
00565 155* C IF(INNA.LE.0) GO TO 185
00565 156* C COMPUTE ARITHMETIC TEMPERATURES BY SUCCESSIVE POINT OVER-RELAX
00567 157* DN = CON(9)
00570 158* DD = 1.0-DN
00571 159* LAX = KON(5)
00572 160* DO 170 I = 1,LAX
00575 161* JJ1 = J1
00576 162* JJ2 = J2
00577 163* TCGM = 0.0
00600 164* KON(20) = I
00601 165* DO 165 L = L1,MNC
00604 166* SUMCV = 0.0
00605 167* IF(I.GT.1) GO TO 6000
00606 168* INCLUDE VRG2,LIST
00610 169* IF(FLD(4,1,NSQ1(JJ1+1)),EQ.0) GO TO 6000
00613 169* NTYPE = FLD(10,5,NSQ2(JJ2))
00614 169* LA = FLD(5,17,NSQ2(JJ2))
00616 169* LK = FLD(22,14,NSQ2(JJ2))
00617 169* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
00620 169* S005 Q(L) = XK(LK)+Q(L)
00621 169* GO TO 5999
00622 169* S010 Q1 = 0.0
00622 169* S012 CALL DID1WM(T(L),A(LA),XK(LK),Q2)
00623 169* GO TO 5998
00624 169* S015 Q1 = 0.0
00625 169* S017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
00626 169* GO TO 5998
00627 169* S020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00630 169* S022 JJ2 = JJ2+1
00631 169* LA = FLD(5,17,NSQ2(JJ2))
00632 169* LK = FLD(22,14,NSQ2(JJ2))
00633 169* GO TO 5017
00634 169* S025 Q1 = XK(LK)*XK(LA)
00635 169* GO TO 5022
00636 169* S030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00637 169* JJ2 = JJ2+1
00640 169* LA = FLD(5,17,NSQ2(JJ2))
00641 169* LK = FLD(22,14,NSQ2(JJ2))
00642 169* Q2 = XK(LK)*XK(LA)
00643 169* GO TO 5998
00644 169* S035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00645 169* S037 JJ2 = JJ2+1
00646 169* LA = FLD(5,17,NSQ2(JJ2))
00647 169* LK = FLD(22,14,NSQ2(JJ2))
00650 169* GO TO 5012
00651 169* S040 Q1 = XK(LK)*XK(LA)
00652 169* GO TO 5037
00653 169* S996 Q(L) = Q1+Q2+Q(L)

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*NEW
*NEW
**-3

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*NEW
**-2

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*NEW
**-2

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*NEW
**-2

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00654 169* 5999 JJ2 = JJ2+1
 00655 169* 6000 CONTINUE
 00656 169* END
 00657 170* 135 JJ1 = JJ1+1
 00660 171* LG = FLD(5,16,NS01(JJ1))
 00661 172* LTA = FLD(22,14,NS01(JJ1))
 00662 173* IF(I.GT.1) GO TO 4000
 00664 174* INCLUDE VRG2,LIST
 00664 175* CHECK FOR RADIATION CONDUCTOR
 00665 175* IF(FLD(2,1,NS01(JJ1)).EG.0) GO TO 4000
 00667 175* NTYPE = FLD(10,5,NS02(JJ2))
 00670 175* LA = FLD(5,17,NS02(JJ2))
 00671 175* LK = FLD(22,14,NS02(JJ2))
 00672 175* GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,
 00672 175* 3060,3065), NTYPE
 00673 175* 3005 TM = (T(L)+T(LTA))/2.0
 00674 175* 3007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))
 00675 175* GO TO 3999
 00676 175* 3010 TM = T(L)
 00677 175* GO TO 3007
 00700 175* 3015 CALL DID1WM(T(L),A(LA),XK(LK),61)
 00701 175* 3017 JJ2 = JJ2+1
 00702 175* LA = FLD(5,17,NS02(JJ2))
 00703 175* LK = FLD(22,14,NS02(JJ2))
 00704 175* CALL DID1WM(T(LTA),A(LA),XK(LK),62)
 00705 175* GO TO 3998
 00706 175* 3020 G1 = XK(LK)*XK(LA)
 00707 175* GO TO 3017
 00710 175* 3025 CALL DID1WM(T(L),A(LA),XK(LK),61)
 00711 175* JJ2 = JJ2+1
 00712 175* LA = FLD(5,17,NS02(JJ2))
 00713 175* LK = FLD(22,14,NS02(JJ2))
 00714 175* G2 = XK(LK)*XK(LA)
 00715 175* GO TO 3998
 00716 175* 3030 TM = (T(L)+T(LTA))/2.0
 00717 175* 3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))
 00720 175* GO TO 3999
 00721 175* 3035 TM = T(L)
 00722 175* GO TO 3032
 00723 175* 3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
 00724 175* 3042 JJ2 = JJ2+1
 00725 175* LA = FLD(5,17,NS02(JJ2))
 00726 175* LK = FLD(22,14,NS02(JJ2))
 00727 175* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)
 00730 175* GO TO 3998
 00731 175* 3045 G1 = XK(LK)*XK(LA)
 00732 175* GO TO 3042
 00733 175* 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
 00734 175* JJ2 = JJ2+1
 00735 175* LA = FLD(5,17,NS02(JJ2))
 00736 175* LK = FLD(22,14,NS02(JJ2))
 00737 175* G2 = XK(LK)*XK(LA)
 00740 175* GO TO 3998
 00741 175* 3055 TM = (T(L)+T(LTA))/2.0
 00742 175* CALL D2D1WM(TM,CON(14),A(LA),XK(LK),6(LG))
 00743 175* GO TO 3999
 00744 175* 3060 TM = T(LTA)
 00745 175* GO TO 3007
 00746 175* 3065 TM = T(LTA)
 00747 175* GO TO 3032

*NEW
 *NEW
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00750 175*
00751 175*
00753 175*
00754 175*
00755 175*
00756 176*
00760 177*
00761 178*
00762 179*
00763 180*
00764 181*
00765 182*
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00767 185*
00771 186*
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00772 188*
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00773 190*
00773 191*
00774 192*
00776 193*
00777 194*
01000 195*
01000 196*
01002 197*
01004 198*
01004 199*
01006 200*
01006 201*
01007 202*
01010 203*
01013 204*
01014 205*
01015 206*
01017 207*
01020 208*
01021 209*
01021 210*
01023 211*
01025 212*
01026 213*
01027 214*
01027 215*
01030 216*
01030 217*
01032 218*
01033 219*
01034 220*
01034 221*
01035 222*
01035 223*
01037 224*
01041 225*
01042 226*
01042 227*
01043 228*
01044 229*
01044 230*

3998 G(LG) = 1./((1./61+1./62)
IF(FLD(3,1,NSQ(JJ1)),EQ.1) G(LG) = G1*62
3999 JJ2 = JJ2+1
4000 CONTINUE
END
IF(FLD(3,1,NSQ(JJ1)),EQ.0) GO TO 140
T1 = T(L)+460.0
T2 = T(LTA)+460.0
GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
GO TO 145
140 GV = G(LG)
145 SUMC = SUMC+GV
SUMCV = SUMCV+GV*(LTA)
CHECK FOR LAST CONDUCTOR
IF(NSQ(JJ1).GT.0) GO TO 135
T2 = DD*(L)+DN*(SUMCV+Q(L))/SUMC
OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
T1 = ABS(T(L)-T2)
STORE THE NEW TEMPERATURE
T(L) = T2
SAVE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
IF(TCGM.GE.T1) GO TO 165
TCGM = T1
KON(37) = L
165 CONTINUE
SEE IF RELAXATION CRITERIA WAS MET
IF(TCGM.LE.CON(19)) GO TO 175
170 CONTINUE
STORE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
CON(30) = TCGM
175 COMPUTE THE ARITHMETIC TEMPERATURE CHANGE
TCGM = 0.0
DO 180 I = L1,NC
LE = IE+I
T1 = ARS(T(I)-X(LE))
IF(T1.LT.TCGM) GO TO 180
TCGM = T1
KON(38) = I
180 CONTINUE
SEE IF ATMPCA WAS SATISFIED
IF(TCGM.GT.CON(11)) GO TO 125
CON(16) = TCGM
KON(12) = 0
CALL VARBL2
CHECK THE BACKUP SWITCH
IF(KON(12).NE.0) GO TO 105
ADVANCE TIME
CON(13) = CON(11)
TSUM = TSUM+TSTEP
TSTEP = DELTA*0.95
CHECK FOR TIME TO PRINT
IF(TSUM.GE.CON(18)) GO TO 190
CHECK FOR PRINT EVERY ITERATION
IF(KON(7).EQ.0) GO TO 10
CALL OUTCAL
GO TO 10
TRY TO EVEN THE OUTPUT INTERVALS
190 TPRINT = TPRINT+TSUM
195 CALL OUTCAL
C IS TIME GREATER THAN END COMPUTE TIME

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01045 231* IFCON(1)*1.000001.LT.CON(3)) GO TO 5
01047 232* NTH = IE
01050 233* NDIM = NLA
01051 234* RETURN
01052 235* 995 WRITE(6,885)
01054 236* GO TO 1000
01055 237* 996 WRITE(6,886)
01057 238* GO TO 1000
01060 239* 997 WRITE(6,887)
01062 240* GO TO 1000
01063 241* 998 WRITE(6,888) I
01066 242* GO TO 1000
01067 243* 999 WRITE(6,889)
01071 244* 1000 CALL OUTCAL
01072 245* CALL EXIT
01073 246* 885 FORMAT(46H CNFRWD REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
01074 247* 886 FORMAT(24H CSGMIN ZERO OR NEGATIVE)
01075 248* 887 FORMAT(20H TIME STEP TOO SMALL)
01076 249* 888 FORMAT(18,20H LOCATIONS AVAILABLE)
01077 250* 889 FORMAT(19H NO OUTPUT INTERVAL)
01100 251* END

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END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
 CNFRWD SYMBOLIC
 CNFRWD CODE RELOCATABLE

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Q1W FOR,* CNFRDL,CNFRDL
 UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-100 CREATED ON 20 AUG 70
 THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:31

21 FEB 71

SUBROUTINE CNFRDL ENTRY POINT 003463

STORAGE USED (BLOCK, NAME, LENGTH)

0001 *CODE 003476
 0000 *CONST+TEMP 000072
 0002 *SIMPLE VAR 000046
 0004 *ARRAYS 000000
 0005 *BLANK 000000
 0006 TITLE 000001
 0007 TEMP 000001
 0010 CAP 000001
 0011 SOURCE 000001
 0012 COMD 000001
 0013 PC1 000001
 0014 PC2 000001
 0015 KONST 000001
 0016 ARRAY 000001
 0017 FIXCON 000001
 0020 XSPACE 000003
 0021 DIMENS 000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022 VARL1
 0023 D1DIWM
 0024 PLYAWM
 0025 D2DIWM
 0026 VARUL2
 0027 OUTCAL
 0030 EXIT
 0031 NERR25
 0032 NWDU5
 0033 NIO25
 0034 NER105

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016 R 000000 A	0010 R 000000 C	0002 R 000015 CKM	0017 R 000000 CON	0002 R 000021 C1
0002 R 000022 C2	0002 R 000037 DD	0002 R 000035 DELTA	0002 R 000036 DN	0012 R 000000 G
0002 R 000034 GV	0002 R 000030 G1	0002 R 000031 G2	0006 R 000000 H	0002 I 000004 I
0002 I 000002 IE	0002 I 000041 JJ1	0002 I 000042 JJ2	0002 I 000012 J1	0002 I 000013 J2
0015 I 000000 K	0017 I 000000 KON	0002 I 000043 L	0002 I 000017 LA	0002 I 000040 LAX
0002 I 000011 LE	0002 I 000025 LG	0002 I 000020 LK	0021 I 000006 L501	0021 I 000007 L502
0002 I 000026 LYA	0002 I 000005 LI	0021 I 000005 NAT	0020 I 000004 NCT	0020 I 000000 N71M
0021 I 000003 NGT	0002 I 000003 NLA	0021 I 000001 NNA	0002 I 000001 NRC	0002 I 000016 NTYPE
0021 I 000002 NNT	0013 I 000000 NS01	0014 I 000000 NS02	0020 I 000001 NTH	0002 R 000024 G2
0020 000002 NX	0002 R 000000 PASS	0011 R 000000 Q	0002 R 000023 O1	0002 R 000027 TM
0002 R 000044 SUMC	0002 R 000045 SUMCV	0007 R 000000 T	0002 R 000014 YCGM	0002 R 000033 T2
0002 R 000007 TPRI;T	0002 R 000046 TSTEP	0002 R 000010 T5UM	0002 R 000032 T1	0001 003277 I0016
0020 R 000002 X	0015 R 000000 XK	0001 000126 I0L	0001 003441 I000L	

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0001 000330 1005L 0001 000351 1010L 0001 000367 1012L 0001 000422 1015L 0001 000430 1020L
0001 000470 1025L 0001 000514 1030L 0001 000535 1032L 0001 000573 1035L 0001 000601 1040L
0001 000644 1045L 0001 002032 105L 0001 002064 115L 0001 002075 120L 0001 002103 125L
0001 002111 130L 0001 002441 135L 0001 000135 15L 0001 003207 155L 0001 003212 160L
0001 003257 165L 0001 003270 175L 0001 003321 180L 0001 003332 185L 0001 003361 190L
0001 003364 195L 0001 000666 1998L 0001 000671 1999L 0001 000207 2006 0001 000674 2000L
0001 001233 2005L 0001 001240 2007L 0001 001260 2010L 0001 001263 2015L 0001 001301 2017L
0001 001335 2020L 0001 001343 2025L 0001 001403 2030L 0001 001410 2032L 0001 001433 2035L
0001 001436 2040L 0001 001457 2042L 0001 001516 2045L 0001 001524 2050L 0001 001567 2055L
0001 001615 2060L 0001 001621 2065L 0001 000226 2126 0001 000257 2326 0001 000147 25L
0001 001625 2998L 0001 001653 2999L 0001 000162 30L 0001 001656 3000L 0001 002527 3005L
0001 002534 3007L 0001 002554 3010L 0001 002557 3015L 0001 002575 3017L 0001 002631 3020L
0001 002637 3025L 0001 002677 3030L 0001 002704 3032L 0001 002727 3035L 0001 002732 3040L
0001 002753 3042L 0001 003012 3045L 0001 003020 3050L 0001 003063 3055L 0001 003111 3060L
0001 003115 3065L 0001 003121 3098L 0001 003147 3099L 0001 003152 4000L 0001 000741 4005L
0001 000746 4010L 0001 000747 4012L 0001 000766 4015L 0001 000767 4017L 0001 001004 4020L
0001 001020 4022L 0001 001035 4025L 0001 001043 4030L 0001 001101 4035L 0001 001115 4037L
0001 001132 4040L 0001 000240 45L 0001 001745 4746 0001 001140 4998L 0001 001144 4999L
0001 000114 5L 0001 001147 5000L 0001 002233 5005L 0001 002240 5010L 0001 002241 5012L
0001 002260 5015L 0001 002261 5017L 0001 002276 5020L 0001 002312 5022L 0001 002327 5025L
0001 002335 5030L 0001 002373 5035L 0001 002407 5037L 0001 002424 5040L 0001 002036 5256
0001 002140 5636 0001 002161 5726 0001 002432 5998L 0001 002436 5999L 0001 002441 6000L
0001 001147 70L 0001 001713 75L 0001 001716 80L 0000 000000 885F 0000 000011 886F
0000 000016 887F 0000 000023 888F 0000 000030 889F 0000 002005 95L
0001 003403 995L 0001 003411 996L 0001 003417 997L 0001 003425 998L 0001 003434 999L

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1* SUBROUTINE CNFRDL

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00101 C EXPLICIT FORWARD DIFFERENCING EXECUTION SUBROUTINE FOR SINDA F-V
00102 C THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED
00103 C INCLUDE COMM,LIST
00104 COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
00105 COMMON /FIXCON/KON(1) /PC2/NSQ2(1) /KONST/K(1) /ARRAY/A(1)
00106 COMMON /DIMENS/ NND,NNA,NNT,NGT,NCT,NAT,LSQ1,LSQ2
00107 DIMENSION CON(1),XK(1),NX(1)
00108 EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
00109 END
00110 INCLUDE DEFF,LIST
00111 C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00112 C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEIN)
00113 C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00114 C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEO)
00115 C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR*EXPLICIT (CSGFAC)
00116 C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00117 C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00118 C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
00119 C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
00120 C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
00121 C CC10 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (DAMPD)
00122 C CC11 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (ATMPCA)
00123 C CC12 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEO)
00124 C CC13 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)
00125 C CC14 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
00126 C CONTROL CONSTANT 17 IS RESERVED FOR THE C/56 MINIMUM (CSGMIN)
00127 C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)

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00113 5* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 5* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED,INTEGER (LOOPCT)
00113 5* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEL)
00113 5* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
00113 5* C CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
00113 5* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
00113 5* C CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
00113 5* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 5* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 5* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 5* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 5* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 5* C CC31 IS INDICATOR, 0=THERMAL SPCS,1=THERMAL LPCS,2=GENERAL (LSPCS)
00113 5* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGBAL)
00113 5* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 5* C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 5* C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 5* C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 5* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 5* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 5* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 5* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDS (LAXFAC)
00113 5* C CC50 IS NOT USED AT PRESENT
00114 5* C END
00115 6* IF(CON(4).LT.1.0) CON(4) = 1.0
00117 7* IF(KON(5).LE.0) KON(5) = 1
00121 8* IF(CON(6).LE.0.) CON(6) = 1.E+8
00123 9* IF(CON(8).LE.0.) CON(8) = 1.E+8
00125 10* IF(CON(9).LE.0.) CON(9) = 1.0
00127 11* IF(CON(11).LE.0.) CON(11) = 1.E+8
00131 12* IF(CON(18).LE.0.) GO TO 999
00133 13* IF(CON(19).LE.0.) CON(19) = 1.E+8
00135 14* IF(KON(31).NE.1) GO TO 995
00137 15* PASS = -1.0
00140 16* NNC = MND+RNA
00141 17* IE = NTH
00142 18* NLA = NDIM
00143 19* NTH = NTH+INC
00144 20* NDIM = NDIM+NNC
00144 21* CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00145 22* I = NLA+NNC
00146 23* IF(I.LT.0) GO TO 998
00150 24* LI = NND+1
00151 25* TSTEP = CON(18)
00152 26* TPRINT = CON(13)
00152 27* INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00153 28* 5 TSUM = 0.0
00153 29* C DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00154 30* IF(CON(13)+CON(18).LE.CON(31)) GO TO 10
00156 31* C DONT EXCEED IT
00156 32* CON(18) = CON(13)-CON(13)
00156 33* C IS THE TIME STEP LARGER THAN ALLOWED
00157 34* TSTEP = CON(18)
00161 35* C DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
00162 37* 15 IF(TSUM+TSTEP-CON(18)) 25,30,20
00165 38* C DONT EXCEED IT
00165 39* 20 TSTEP = CON(16)-TSUM
00166 40* GO TO 30

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00166 41* C DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
00167 42* C IF(TSUM+2.0*TSTEP,LE,CON(18)) GO TO 30
00167 43* C APPROACH THE OUTPUT INTERVAL GRADUALLY
00171 44* C TSTEP = (CON(18)-TSUM)/2.0
00171 45* C STORE DELTA TIME STEP IN THE CONSTANTS
00172 46* C CON(2) = TSTEP
00172 47* C IS THE TIME STEP USED LESS THAN THE TIME STEP ALLOWED
00173 48* C IF(TSTEP,LT,CON(21)) GO TO 997
00173 49* C CALCULATE THE NEW TIME
00175 50* C CON(1) = TPRINT+TSUM+TSTEP
00175 51* C COMPUTE THE MEAN TIME BETWEEN ITERATIONS
00176 52* C CON(14) = (CON(1)+CON(13))/2.0
00176 53* C ZERO OUT ALL SOURCE LOCATIONS AND EXTRA LOCATIONS
00177 54* C DO 35 I = 1,NND
00202 55* C LE = IE+1
00203 56* C X(LE) = 0.0
00204 57* C G(I) = 0.0
00205 58* C 35 CONTINUE
00205 59* C SHIFT THE ARITHMETIC TEMPERATURES INTO THE EXTRA LOCATIONS
00207 60* C IF(NNA,LE,0) GO TO 45
00211 61* C DO 40 I = L1,NNC
00214 62* C G(I) = 0.0
00215 63* C LE = IE+1
00216 64* C X(LE) = T(I)
00217 65* C 40 CONTINUE
00221 66* C 45 KON(12) = 0
00222 67* C CALL VARBLI
00223 68* C IF(KON(12),NE,0) GO TO 10
00225 69* C J1 = 0
00226 70* C J2 = 1
00227 71* C TCGM = 0.0
00230 72* C CKM = 1.E+8
00230 73* C CALCULATE 0 SUM AND 6 SUM
00231 74* C DO 85 I = 1,NND
00234 75* C LE = IE+1
00235 76* C INCLUDE VARC,LIST
00236 76* C IF(FLD(1,1,NSQ1(J1+1),EG,0) GO TO 2000
00240 76* C NTYPE = FLD(0,5,NSQ2(J2))
00241 76* C LA = FLD(5,17,NSQ2(J2))
00242 76* C LK = FLD(22,14,NSQ2(J2))
00243 76* C GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
00244 76* C 1005 CALL DID1WM(T(I),A(LA),XK(LK),C(I))
00245 76* C GO TO 1999
00246 76* C 1010 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00247 76* C 1012 J2 = J2+1
00250 76* C LA = FLD(5,17,NSQ2(J2))
00251 76* C LK = FLD(22,14,NSQ2(J2))
00252 76* C CALL DID1WM(T(I),A(LA),XK(LK),C2)
00253 76* C GO TO 1998
00254 76* C 1015 C1 = XK(LK)*XK(LA)
00255 76* C GO TO 1012
00256 76* C 1020 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00257 76* C J2 = J2+1
00260 76* C LA = FLD(5,17,NSQ2(J2))
00261 76* C LK = FLD(22,14,NSQ2(J2))
00262 76* C C2 = XK(LK)*XK(LA)
00263 76* C GO TO 1998
00264 76* C 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
00265 76* C GO TO 1999
00266 76* C 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)

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*NEW
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CNFRDL,CNFRDL

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00267 76* 1032 J2 = J2+1
00270 76* LA = FLD(5,17,NSQ2(J2))
00271 76* LK = FLD(22,14,NSQ2(J2))
00272 76* CALL PLYAW(A(LA),T(I),A(LA+1),XK(LK),C2)
00273 76* GO TO 1998
00274 76* C1 = XK(LK)*XK(LA)
00275 76* GO TO 1032
00276 76* 1040 CALL PLYAW(A(LA),T(I),A(LA+1),XK(LK),C1)
00277 76* J2 = J2+1
00300 76* LA = FLD(5,17,NSQ2(J2))
00301 76* LK = FLD(22,14,NSQ2(J2))
00302 76* C2 = XK(LK)*XK(LA)
00303 76* GO TO 1998
00304 76* 1045 CALL D2D1WM(T(I),CON(14),A(LA),XK(LK),C(I))
00305 76* GO TO 1999
00306 76* C(I) = C1+C2
00307 76* 1999 J2 = J2+1
00310 76* 2000 CONTINUE
00311 76* END
00312 77* INCLUDE VARO,LIST
00313 77* IF(FLD(4,1,NSQ1(J1+1)),EG,0) GO TO 5000
00315 77* NTYPE = FLD(0,5,NSQ2(J2))
00316 77* LA = FLD(5,17,NSQ2(J2))
00317 77* LK = FLD(22,14,NSQ2(J2))
00320 77* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4050), NTYPE
00321 77* Q(I) = XK(LK)+Q(I)
00322 77* GO TO 4999
00323 77* 4010 Q1 = 0.0
00324 77* 4012 CALL D1D1WM(T(I),A(LA),XK(LK),Q2)
00325 77* GO TO 4998
00326 77* 4015 Q1 = 0.0
00327 77* 4017 CALL D1D1WM(CON(14),A(LA),XK(LK),Q2)
00330 77* GO TO 4998
00331 77* 4020 CALL D1D1WM(CON(14),A(LA),XK(LK),Q1)
00332 77* 4022 J2 = J2+1
00333 77* LA = FLD(5,17,NSQ2(J2))
00334 77* LK = FLD(22,14,NSQ2(J2))
00335 77* GO TO 4017
00336 77* 4025 Q1 = XK(LK)*XK(LA)
00337 77* GO TO 4022
00340 77* 4030 CALL D1D1WM(CON(14),A(LA),XK(LK),Q1)
00341 77* J2 = J2+1
00342 77* LA = FLD(5,17,NSQ2(J2))
00343 77* LK = FLD(22,14,NSQ2(J2))
00344 77* Q2 = XK(LK)*XK(LA)
00345 77* GO TO 4998
00346 77* 4035 CALL D1D1WM(CON(14),A(LA),XK(LK),Q1)
00347 77* 4037 J2 = J2+1
00350 77* LA = FLD(5,17,NSQ2(J2))
00351 77* LK = FLD(22,14,NSQ2(J2))
00352 77* GO TO 4012
00353 77* 4040 Q1 = XK(LK)*XK(LA)
00354 77* GO TO 4037
00355 77* 4998 Q(I) = Q1+Q2+Q(I)
00356 77* 4999 J2 = J2+1
00357 77* 5000 CONTINUE
00360 77* END
00361 78* 70 J1 = J1+1
00362 79* LG = FLD(5,10,NSQ1(J1))
00363 80* LTA = FLD(22,14,NSQ1(J1))

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CNFRDL,CNFPDL

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00461 85* T2 = T(LTA)+460.0
00462 86* GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
00463 87* GO TO 80
00464 88* GV = G(LG)
00465 89* C OBTAIN THE Q RATE THRU THE CONDUCTOR
00466 90* G(I) = G(I)+GV*(T(LTA)-T(I))
00467 91* C SAVE SUMMATION OF CONDUCTORS
00468 92* X(LE) = X(LE)+GV
00469 93* C CHECK FOR LAST CONDUCTOR
00470 94* IF(NSQI(J1).GT.0) GO TO 70
00471 95* C CONTINUE
00472 96* C OBTAIN NEW DIFFUSION TEMPERATURES, OTMPC AND CSGMIN
00473 97* DO 100 I = 1,NND
00474 98* LE = IE+I
00475 99* C CALCULATE C/SK MINIMUM
00476 100* T1 = C(I)/X(LE)
00477 101* IF(T1.GE.CKM) GO TO 90
00478 102* CKM = T1
00479 103* KON(35) = I
00480 104* C COMPUTE NEW TEMPERATURES USING CALCULATED SOURCE TERMS
00481 105* T1 = TSTEP*(I)/C(I)
00482 106* C CALCULATE THE ABSOLUTE VALUE TEMPERATURE CHANGE
00483 107* T2 = ABS(T1)
00484 108* C SAVE THE LARGEST TEMPERATURE CHANGE
00485 109* IF(TCGM.GE.T2) GO TO 95
00486 110* TCGM = T2
00487 111* KON(36) = I
00488 112* C STORE THE TEMPERATURES
00489 113* X(LE) = T(I)
00490 114* T(I) = T(I)+T1
00491 115* C CONTINUE
00492 116* CON(17) = CKM
00493 117* DELTA = CKM/CON(4)
00494 118* IF(CKM.LE.0.0) GO TO 996
00495 119* C CHECK FOR FIRST PASS
00496 120* IF(PASS.GT.0.0) GO TO 115
00497 121* C UNDO THE TEMPERATURE CALCULATIONS
00498 122* DO 110 I = 1,NIC
00499 123* LE = IE+I
00500 124* T(I) = X(LE)
00501 125* C CONTINUE
00502 126* IF(PASS.GT.0.0) GO TO 15
00503 127* PASS = 1.0
00504 128* CON(1) = TPRINT
00505 129* CON(2) = 0.0
00506 130* TSTEP = DELTA*0.95
00507 131* GO TO 195
00508 132* C IS THE TIME STEP USED LESS THAN THE TIME STEP CALCULATED
00509 133* IF(TSTEP.LE.DELTA) GO TO 130
00510 134* C COMPUTE THE TIME STEP
00511 135* TSTEP = DELTA*0.95
00512 136* GO TO 105
00513 137* C TSTEP = 0.95*TSTEP+CON(6)/TCGM
00514 138* GO TO 105
00515 139* C TSTEP = 0.95*TSTEP+CON(11)/TCGM
00516 140* GO TO 105
00517 141* C SEE IF THE TEMPERATURE CHANGE WAS TOO LARGE
00518 142* IF(TCGM.GT.CON(6)) GO TO 120
00519 143* C STORE THE MAXIMUM DIFFUSION TEMPERATURE CHANGE
00520 144* CON(15) = TCGM

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C 145* CHECK TO SEE IF THERE ARE ANY ARITHMETIC NODES
C 146* IF (NNA.LE.0) GO TO 185
C 147* COMPUTE ARITHMETIC TEMPERATURES BY SUCCESSIVE POINT OVER-RELAX
C 148* DN = CON(9)
C 149* DD = 1.0-DN
C 150* LAX = KON(5)
C 151* DO 170 I = 1,LAX
C 152* JJ1 = J1
C 153* JJ2 = J2
C 154* TCGM = 0.0
C 155* KON(20) = I
C 156* DO 165 L = L1,NMC
C 157* SUMC = 0.0
C 158* SUMCV = 0.0
C 159* IF (I.GT.1) GO TO 6000
C 160* INCLUDE VRQ2,LIST
C 161* IF (FLD(4,1,NSQ1(JJ1+1)),EQ.0) GO TO 6000
C 162* NTYPE = FLD(0,5,NSQ2(JJ2))
C 163* LA = FLD(5,17,NSQ2(JJ2))
C 164* LK = FLD(22,14,NSQ2(JJ2))
C 165* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5050), NTYPE
C 166* 5005 G(L) = XK(LK)+O(L)
C 167* GO TO 5999
C 168* 5010 G1 = 0.0
C 169* 5012 CALL DID1WM(T(L),A(LA),XK(LK),O2)
C 170* GO TO 599B
C 171* 5015 G1 = 0.0
C 172* 5017 CALL DID1WM(CON(14),A(LA),XK(LK),O2)
C 173* GO TO 599B
C 174* 5020 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
C 175* 5022 JJ2 = JJ2+1
C 176* LA = FLD(5,17,NSQ2(JJ2))
C 177* LK = FLD(22,14,NSQ2(JJ2))
C 178* GO TO 5017
C 179* 5025 G1 = XK(LK)+XK(LA)
C 180* GO TO 5022
C 181* 5030 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
C 182* JJ2 = JJ2+1
C 183* LA = FLD(5,17,NSQ2(JJ2))
C 184* LK = FLD(22,14,NSQ2(JJ2))
C 185* G2 = XK(LK)+XK(LA)
C 186* GO TO 599B
C 187* 5035 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
C 188* 5037 JJ2 = JJ2+1
C 189* LA = FLD(5,17,NSQ2(JJ2))
C 190* LK = FLD(22,14,NSQ2(JJ2))
C 191* GO TO 5012
C 192* 5040 G1 = XK(LK)+XK(LA)
C 193* GO TO 5037
C 194* 599B G(L) = G1+O2+O(L)
C 195* 5999 JJ2 = JJ2+1
C 196* 6000 CONTINUE
C 197* ENO
C 198* 135 JJ1 = JJ1+1
C 199* LG = FLD(5,16,NSQ1(JJ1))
C 200* L1A = FLD(22,14,NSQ1(JJ1))
C 201* IF (I.GT.1) GO TO 4000
C 202* INCLUDE VRG2,LIST
C 203* CHECK FOR RADIATION CONDUCTOR
C 204* IF (FLD(2,1,NSQ1(JJ1)),EQ.0) GO TO 4000
C 205* 166*
C 206* 00554
C 207* 00555
C 208* 00556
C 209* 00557
C 210* 00560
C 211* 00561
C 212* 00562
C 213* 00565
C 214* 00566
C 215* 00567
C 216* 00570
C 217* 00571
C 218* 00574
C 219* 00575
C 220* 00576
C 221* 00600
C 222* 00601
C 223* 00603
C 224* 00604
C 225* 00605
C 226* 00606
C 227* 00607
C 228* 00610
C 229* 00611
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C 232* 00614
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C 250* 00636
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C 252* 00640
C 253* 00641
C 254* 00642
C 255* 00643
C 256* 00644
C 257* 00645
C 258* 00646
C 259* 00647
C 260* 00650
C 261* 00651
C 262* 00652
C 263* 00654
C 264* 00655

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CNFRDL,CNFRDL

00657	166*	NTYPE = FLD(0,5,NSQ2(JJ2))	*NEW
00660	166*	LA = FLD(15,17,NSQ2(JJ2))	*NEW
00661	166*	LK = FLD(22,14,NSQ2(JJ2))	*NEW
00662	166*	GOTO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,	***3
00663	166*	3060,3065), NTYPE	
00663	166*	TM = (T(L)+T(LTA))/2.0	
00664	166*	3007 CALL DD1WM(TM,A(LA),XK(LK),6(LG))	
00665	166*	GO TO 3999	
00666	166*	TM = T(L)	
00667	166*	GO TO 3007	
00670	166*	3015 CALL DD1WM(T(L),A(LA),XK(LK),61)	
00671	166*	3017 JJ2 = JJ2+1	
00672	166*	LA = FLD(15,17,NSQ2(JJ2))	
00673	166*	LK = FLD(22,14,NSQ2(JJ2))	
00674	166*	CALL DD1WM(T(LTA),A(LA),XK(LK),62)	
00675	166*	GO TO 3998	
00676	166*	61 = XK(LK)*XK(LA)	
00677	166*	GO TO 3017	
00700	166*	3025 CALL DD1WM(T(L),A(LA),XK(LK),61)	
00701	166*	JJ2 = JJ2+1	
00702	166*	LA = FLD(15,17,NSQ2(JJ2))	
00703	166*	LK = FLD(22,14,NSQ2(JJ2))	
00704	166*	G2 = XK(LK)*XK(LA)	
00705	166*	GO TO 3998	
00706	166*	TM = (T(L)+T(LTA))/2.0	
00707	166*	3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))	
00710	166*	GO TO 3999	
00711	166*	TM = T(L)	
00712	166*	GO TO 3032	
00713	166*	3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
00714	166*	3042 JJ2 = JJ2+1	
00715	166*	LA = FLD(15,17,NSQ2(JJ2))	
00716	166*	LK = FLD(22,14,NSQ2(JJ2))	
00717	166*	CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)	
00720	166*	GO TO 3998	
00721	166*	61 = XK(LK)*XK(LA)	
00722	166*	GO TO 3042	
00723	166*	3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
00724	166*	JJ2 = JJ2+1	
00725	166*	LA = FLD(15,17,NSQ2(JJ2))	
00726	166*	LK = FLD(22,14,NSQ2(JJ2))	
00727	166*	G2 = XK(LK)*XK(LA)	
00730	166*	GO TO 3998	
00731	166*	TM = (T(L)+T(LTA))/2.0	
00732	166*	CALL DD1WM(TM,CON(14),A(LA),XK(LK),6(LG))	
00733	166*	GO TO 3999	
00734	166*	TM = T(LTA)	
00735	166*	GO TO 3007	
00736	166*	3065 TM = T(LTA)	
00737	166*	GO TO 3032	
00740	166*	3998 G(LG) = 1./(1./61+1./62)	
00741	166*	IF(FLD(3,1,NSQ1(JJ1)).EQ.1) G(LG) = 61*62	
00743	166*	3999 JJ2 = JJ2+1	
00744	166*	4000 CONTINUE	
00745	166*	END	
00746	167*	IF(FLD(3,1,NSQ1(JJ1)).EQ.0) GO TO 155	
00750	168*	T1 = T(L)+460.0	
00751	169*	T2 = T(LTA)+460.0	
00752	170*	GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)	
00753	.171*	GO TO 160	

CNFRDL,CNFRDL

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00754 172* 155 GV = 6(LG)
00755 173* 160 SUMC = SUMC+GV
00756 174* SUMCV = SUMCV+GV*(ILTA)
00757 175* CHECK FOR LAST CONDUCTOR
00757 176* IF(1/501*(JJ1)-GT.0) GO TO 135
00761 177* T2 = DD*(L)+DN*(SUMCV+Q(L))/SUMC
00761 178* OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
00762 179* T1 = ABS(T(L)-T2)
00762 180* STORE THE NEW TEMPERATURE
00763 181* T(L) = T2
00763 182* SAVE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
00764 183* IF(TCGM.GE.T1) GO TO 165
00766 184* TCGM = T1
00767 185* KON(37) = L
00770 186* 165 CONTINUE
00770 187* SEE IF RELAXATION CRITERIA WAS MET
00772 188* IF(TCGM.LE.CON(19)) GO TO 175
00774 189* CONTINUE
00774 190* 170 STORE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
00776 191* CON(30) = TCGM
00776 192* 175 COMPUTE THE ARITHMETIC TEMPERATURE CHANGE
00777 193* TCGM = 0.0
00777 193* DO 180 I = 1,INC
01000 194* LE = IE+1
01003 195* T1 = ABS(T(I)-X(LE))
01004 196* IF(T1.LT.TCGM) GO TO 180
01005 197* TCGM = T1
01007 198* KON(38) = I
01010 199* 180 CONTINUE
01011 200* SEE IF ATMPCA WAS SATISFIED
01011 201* IF(TCGM.GT.CON(11)) GO TO 125
01013 202* CON(11) = TCGM
01015 203* KON(12) = 0
01016 204* 185 CALL VARBL2
01017 205* CHECK THE BACKUP SWITCH
01017 206* IF(KON(12).NE.0) GO TO 105
01020 207* ADVANCE TIME
01020 208* CON(13) = CON(1)
01022 209* TSUM = TSUM+TSTEP
01023 210* TSTEP = DELTA*0.95
01024 211* CHECK FOR TIME TO PRINT
01024 212* IF(TSUM.GE.CON(18)) GO TO 190
01025 213* 190 CHECK FOR PRINT EVERY ITERATION
01025 214* IF(KON(7).EQ.0) GO TO 10
01027 215* CALL OUTCAL
01031 216* GO TO 10
01032 217* 190 TRY TO EVEN THE OUTPUT INTERVALS
01032 218* TPRINT = TPRINT+TSUM
01033 219* 195 CALL OUTCAL
01034 220* IS TIME GREATER THAN END COMPUTE TIME
01034 221* IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
01035 222* NTH = IE
01037 223* NDIM = NLA
01040 224* RETURN
01041 225* 995 WRITE(6,885)
01042 226* GO TO 1000
01044 227* 996 WRITE(6,886)
01045 228* GO TO 1000
01047 229* 997 WRITE(6,887)
01050 230* GO TO 1000
01052 231*

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CNFRDL.CNFRDL

01053	232*	998	WRITE(6,888) I
01056	233*		GO TO 1000
01057	234*	999	WRITE(6,889)
01061	235*	1000	CALL OUTCAL
01062	236*		CALL EXIT
01063	237*	885	FORMAT(45H CNFRDL REQUIRES LONG PSEUDO-COMPUTE SEQUENCE)
01064	238*	886	FORMAT(24H CSGMIN ZERO OR NEGATIVE)
01065	239*	887	FORMAT(20H TIME STEP TOO SMALL)
01066	240*	888	FORMAT(18,20H LOCATIONS AVAILABLE)
01067	241*	889	FORMAT(19H NO OUTPUT INTERVAL)
01070	242*		END

END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
CNFRDL SYMBOLIC
CNFRDL CODE RELOCATABLE

21 FEB 71

CNFAST,CNFAST

DLW FOR,* CNFAST,CNFAST
URIVAC 1108 FORTRAN V ATHENA VERSION 131K-10D CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:25

SUBROUTINE CNFAST ENTRY POINT 003360

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	003373
0000	*CONST+TEMP	000070
0002	*SIMPLE VAR	000047
0004	*ARRAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COHO	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARBL1
0023	OUTCAL
0024	D1D1WM
0025	PLYAWM
0026	D2D1WM
0027	VARBL2
0030	EXIT
0031	NERR2\$
0032	NWDU\$
0033	NIO2\$
0034	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C	0002	R	000034	CKM
0002	R	000020	C2	0002	R	000037	DAMPN	0002	R	000040	DAHPO
0002	R	000026	G1	0002	R	000027	G2	0006	R	000000	H
0002	I	000042	JJ1	0002	I	000043	JJ2	0002	I	000012	J1
0017	I	000000	KON	0002	I	000044	L	0002	I	000015	LA
0002	I	000033	LEA	0002	I	000023	LG	0002	I	000016	LK
0002	I	000024	LTA	0021	I	000005	PIAT	0021	I	000004	RCT
0002	I	000003	HLA	0002	I	000004	HN	0021	I	000001	RMA
0020	I	000002	NIJ	0013	I	000000	NSQ1	0014	I	000000	NSQ2
0002	R	000041	HLX	0002	K	000000	PASS	0011	R	000000	Q
0002	R	000025	TA	0002	K	000045	SUMC	0002	R	000046	SUMCV
0002	R	000031	T2	0002	R	000005	TPRINT	0002	R	000006	TSTEP
				0020	R	000002	X	0015	R	000000	XK
0002	R	000017	C1	0017	R	000000	CON	0020	I	000000	NDIM
0002	R	000032	GV	0012	R	000000	G	0020	I	000001	NTH
0002	I	000002	IE	0002	I	000010	I	0020	I	000021	O1
0015	I	000000	K	0002	I	000013	J2	0002	R	000021	O1
0002	I	000011	LE	0002	I	000036	LAX	0007	R	000000	I
0021	I	000007	L5Q2	0021	I	000006	L5Q1	0002	R	000007	TSUM
0021	I	000003	NS1	0020	I	000000	NDIM	0002	R	000035	TCGM
0021	I	000000	ND	0002	I	000001	NIC	0002	R	000030	T1
0002	I	000014	NTYPE	0020	I	000000	NSQ2	0001	R	000104	I0L
0002	R	000022	Q2	0020	I	000000	NDIM				
0002	R	000035	TCGM								
0002	R	000030	T1								
0001	R	0002036	I0L								

CNFAST,CNFAST

0001	003336	1000L	0001	000322	1005L	0001	000343	1018L	0001	000361	1012L	0001	000414	1015L
0001	000422	1020L	0001	000462	1025L	0001	000506	1030L	0001	000527	1032L	0001	000565	1035L
0001	000573	1040L	0001	000636	1045L	0001	002055	105L	0001	002405	110L	0001	003153	115L
0001	003161	120L	0001	003223	140L	0001	003236	155L	0001	003240	160L	0001	000170	1706
0001	000660	1998L	0001	000663	1999L	0001	000135	20L	0001	000666	2000L	0001	001226	2005L
0001	001233	2007L	0001	000210	2016	0001	001253	2010L	0001	001256	2015L	0001	001274	2017L
0001	001330	2020L	0001	001336	2025L	0001	001376	2030L	0001	001403	2032L	0001	001426	2035L
0001	001431	2040L	0001	001452	2042L	0001	001511	2045L	0001	001517	2050L	0001	001562	2055L
0001	001610	2060L	0001	001614	2065L	0001	000251	2256	0001	000150	25L	0001	001620	2998L
0001	001646	2099L	0001	001651	3000L	0001	002473	3005L	0001	002500	3007L	0001	002520	3010L
0001	002523	3015L	0001	002541	3017L	0001	002575	3020L	0001	002603	3025L	0001	002643	3030L
0001	002650	3032L	0001	002673	3035L	0001	002676	3040L	0001	002717	3042L	0001	002756	3045L
0001	002764	3050L	0001	003027	3055L	0001	003055	3060L	0001	003061	3065L	0001	003065	3998L
0001	003113	3999L	0001	000212	40L	0001	003116	4000L	0001	000731	4005L	0001	000736	4010L
0001	000737	4012L	0001	000756	4015L	0001	000757	4017L	0001	000774	4020L	0001	001010	4022L
0001	001025	4025L	0001	001033	4030L	0001	001071	4035L	0001	001105	4037L	0001	001122	4040L
0001	000237	45L	0001	001130	4998L	0001	001134	4999L	0001	000072	5L	0001	001137	50L
0001	001773	5006	0001	001137	5000L	0001	002177	5005L	0001	002204	5010L	0001	002205	5012L
0001	002224	5015L	0001	002225	5017L	0001	002242	5020L	0001	002256	5022L	0001	002273	5025L
0001	002301	5030L	0001	002337	5035L	0001	002353	5037L	0001	002370	5040L	0001	002104	5376
0001	002125	5466	0001	001702	55L	0001	002376	5998L	0001	002402	5999L	0001	001705	60L
0001	002405	6000L	0001	001753	65L	0001	001762	85L	0000	000000	885F	0000	000011	886F
0000	000016	887F	0000	000023	888F	0000	000026	889F	0001	002014	90L	0001	002032	95L
0001	003300	995L	0001	003306	996L	0001	003314	997L	0001	003323	998L	0001	003331	999L

00101	1*	SUBROUTINE CNFAST
00101	2*	AN EXPLICIT EXECUTION SUBROUTINE FOR SINDA FORTRAN V
00101	3*	THE SHORT PSEUDO COMPUTE SEQUENCE IS REQUIRED
00101	4*	NODES WITH CSG BELOW DTIMEI RECEIVE STEADY STATE SOLUTION
00101	5*	NO BACKING UP IS DONE OR ALLOWED.
00103	6*	INCLUDE COMM,LIST
00104	6*	COMMON /TITLE/H(I) /TEMP/T(I) /CAP/C(I) /SOURCE/Q(I) /COND/G(I)
00105	6*	COMMON /PCI/NSO(I) /PC2/NSQ2(I) /KONST/K(I) /ARRAY/A(I)
00106	6*	COMMON /FIXCON/KON(I) /XSPACE/NDIM,NTH,X(I)
00107	6*	COMMON /DI:HEHS/ NMD,NNA,:NIT,NGT,NCT,NAT,LSO1,LSO2
00110	6*	DIMENSION CON(I),XK(I),NX(I)
00111	6*	EQUIVALENCE (KON(I),CON(I)),(K(I),XK(I)),(X(I),NX(I))
00112	6*	END
00113	7*	INCLUDE DEFF,LIST
00113	7*	C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00113	7*	C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
00113	7*	C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00113	7*	C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
00113	7*	C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR,EXPLICIT (CSGFAC)
00113	7*	C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00113	7*	C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00113	7*	C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
00113	7*	C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
00113	7*	C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
00113	7*	C CC10 CONTAINS THE MAX DIFFUSION TEMP. DAMPING FACTOR (DAMPO)
00113	7*	C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
00113	7*	C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
00113	7*	C CC13 CONTAINS THE NEAR TIME OR PROBLEM START TIME (TIMEO)
00113	7*	C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
00113	7*	C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)
00113	7*	C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
00113	7*	C CONTROL CONSTANT 17 IS RESERVED FOR THE C/5G MINIMUM (CSGMIN)

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00113 C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED,INTEGER (LOOPCT)
00113 C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (CSGMAX)
00113 C CC23 CONTAINS THE C/SG MAXIMUM (CSGRAL)
00113 C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRCL)
00113 C CC25 CONTAINS THE C/SG RANGE CALCULATED (DRLXCA)
00113 C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCC)
00113 C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (LINECT)
00113 C CC28 CONTAINS THE LINE COUNTER, INTEGER (PAGECT)
00113 C CC29 CONTAINS THE PAGE COUNTER, INTEGER (ARLXCC)
00113 C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 C CC31 IS INDICATOR, 0=THERMAL SPCS,1=THERMAL LPCS,2=GENERAL (LSPCS)
00113 C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGRAL)
00113 C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPC
00113 C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 C CC50 IS NOT USED AT PRESENT
00114 C END
00115 8 IF(KON(5).LE.0) KON(5) = 1
00117 9 IF(CON(18).LE.0.) CON(18) = 1.E+8
00121 10 IF(CON(9).LE.0.) CON(9) = 1.0
00123 11 IF(CON(18).LE.0.) GO TO 999
00125 12 IF(CON(19).LE.0.) CON(19) = 1.E+8
00127 13 IF(CON(21).LE.0.) GO TO 998
00131 14 IF(KON(31).NE.0) GO TO 995
00133 15 PASS = -1.0
00134 16 NNC = NNA+NND
00135 17 IE = NTH
00136 18 NLA = NDIM
00137 19 NTH = NTH+NND
00140 20 NDIM = NDIM+NND
00141 21 IF(NDIM.LT.0) GO TO 997
00143 22 NN = NND+1
00144 23 TPRINT = CON(13)
00145 24 TSTEP = CON(21)
00146 25 TSUM = 0.0
00147 26 IF(CON(13)+CON(18).GT.CON(3)) CON(18) = CON(3)-CON(13)
00151 27 IF(TSTEP.GT.CON(8)) TSTEP = CON(8)
00153 28 IF(TSTEP.LT.CON(21)) TSTEP = CON(21)+1.000001
00155 29 IF(TSUM+TSTEP-CON(18)) 20,25,15
00160 30 TSTEP = CON(18)-TSUM
00161 31 GO TO 25
00162 32 IF(TSUM+2.0*TSTEP.GT.CON(18)) TSTEP = 0.5*(CON(18)-TSUM)
00164 33 CON(2) = TSTEP
00165 34 CON(1) = TPRINT+TSUM+TSTEP
00166 35 CON(14) = 0.5*(CON(1)+CON(13))
00172 36 DO 30 I = 1,NND
00173 37 O(I) = 0.0
00174 38 LL = IE+1
00176 39 IF(NNA.LE.0) GO TO 40
00200 40 DO 35 I = NN,NP,C
00200 41*

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42* 00203 G(I) = 0.0
43* 00204 35 CONTINUE
44* 00206 40 KON(I2) = 0
45* 00207 CALL VARBL1
46* 00210 IF(KON(I2).NE.0) GO TO 10
47* 00212 IF(PASS.GT.0.) GO TO 45
48* 00214 PASS = 1.0
49* 00215 CON(1) = TPRINT
50* 00216 CON(2) = 0.0
51* 00217 CALL OUTCAL
52* 00220 CON(1) = TPRINT+TSTEP
53* 00221 CON(2) = TSTEP
54* 00222 J1 = 0
55* 00223 J2 = 1
56* 00224 DO 85 I = 1,NND
57* 00227 LE = IE+1
58* 00230 INCLUDE VARC*LIST
58* 00231 IF(FLD(I,1,NSQ1(J1+1),EQ.0) GO TO 2000
58* 00233 NTYPE = FLD(I,5,NSQ2(J2))
58* 00234 LA = FLD(5,17,NSQ2(J2))
58* 00235 LK = FLD(22,14,NSQ2(J2))
58* 00236 GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
58* 00237 1005 CALL DID1WM(T(I),A(LA),XK(LK),C(I))
58* 00240 GO TO 1999
58* 00241 1010 CALL DID1WM(T(I),A(LA),XK(LK),C1)
58* 00242 1012 J2 = J2+1
58* 00243 LA = FLD(5,17,NSQ2(J2))
58* 00244 LK = FLD(22,14,NSQ2(J2))
58* 00245 CALL DID1WM(T(I),A(LA),XK(LK),C2)
58* 00246 GO TO 1998
58* 00247 1015 C1 = XK(LK)*XK(LA)
58* 00250 GO TO 1012
58* 00251 1020 CALL DID1WM(T(I),A(LA),XK(LK),C1)
58* 00252 J2 = J2+1
58* 00253 LA = FLD(5,17,NSQ2(J2))
58* 00254 LK = FLD(22,14,NSQ2(J2))
58* 00255 C2 = XK(LK)*XK(LA)
58* 00256 GO TO 1998
58* 00257 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
58* 00260 GO TO 1999
58* 00261 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
58* 00262 1032 J2 = J2+1
58* 00263 LA = FLD(5,17,NSQ2(J2))
58* 00264 LK = FLD(22,14,NSQ2(J2))
58* 00265 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
58* 00266 GO TO 1998
58* 00267 1035 C1 = XK(LK)*XK(LA)
58* 00270 GO TO 1032
58* 00271 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
58* 00272 J2 = J2+1
58* 00273 LA = FLD(5,17,NSQ2(J2))
58* 00274 LK = FLD(22,14,NSQ2(J2))
58* 00275 C2 = XK(LK)*XK(LA)
58* 00276 GO TO 1998
58* 00277 1045 CALL U2D1WM(T(I),CON(I4),A(LA),XK(LK),C(I))
58* 00300 GO TO 1999
58* 00301 1998 C(I) = C1+C2
58* 00302 1999 J2 = J2+1
58* 00303 2000 CONTINUE
58* 00304 END

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CNFAST,CNFAST

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00305 59* INCLUDE VARG,LIST
00306 59* IF(FLD(4,1,NSQ(J1+1),EQ,0) GO TO 5000
00310 59* NTYPE = FLD(0,5,NSQ(J2))
00311 59* LA = FLD(5,17,NSQ(J2))
00312 59* LK = FLD(22,14,NSQ(J2))
00313 59* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4050), NTYPE
00314 59* 4005 Q(I) = XK(LK)+0(I)
00315 59* GO TO 4999
00316 59* 4010 OI = 0+0
00317 59* 4012 CALL DID1WM(T(I),A(LA),XK(LK),02)
00320 59* GO TO 4998
00321 59* 4015 OI = 0+0
00322 59* 4017 CALL DID1WM(CON(14),A(LA),XK(LK),02)
00323 59* GO TO 4998
00324 59* 4020 CALL DID1WM(CON(14),A(LA),XK(LK),01)
00325 59* 4022 J2 = J2+1
00326 59* LA = FLD(5,17,NSQ(J2))
00327 59* LK = FLD(22,14,NSQ(J2))
00330 59* GO TO 4017
00331 59* 4025 OI = XK(LK)*XK(LA)
00332 59* GO TO 4022
00333 59* 4030 CALL DID1WM(CON(14),A(LA),XK(LK),01)
00334 59* J2 = J2+1
00335 59* LA = FLD(5,17,NSQ(J2))
00336 59* LK = FLD(22,14,NSQ(J2))
00337 59* OI = XK(LK)*XK(LA)
00340 59* GO TO 4998
00341 59* 4035 CALL DID1WM(CON(14),A(LA),XK(LK),01)
00342 59* 4037 J2 = J2+1
00343 59* LA = FLD(5,17,NSQ(J2))
00344 59* LK = FLD(22,14,NSQ(J2))
00345 59* OI = XK(LK)*XK(LA)
00346 59* GO TO 4012
00347 59* 4040 OI = XK(LK)*XK(LA)
00350 59* GO TO 4037
00351 59* 4998 Q(I) = OI+02+0(I)
00352 59* 4999 J2 = J2+1
00353 59* 5000 CONTINUE
00354 60* END
00355 61* 50 J1 = J1+1
00356 62* LG = FLD(5,16,NSQ(J1))
00360 63* IF(LG,EQ,0) GO TO 85
00361 64* LTA = FLD(22,14,NSQ(J1))
00362 64* INCLUDE VARG,LIST
00364 64* IF(FLD(2,1,NSQ(J1),EQ,0) GO TO 3000
00365 64* NTYPE = FLD(0,5,NSQ(J2))
00366 64* LA = FLD(5,17,NSQ(J2))
00367 64* LK = FLD(22,14,NSQ(J2))
00370 64* $ GOT0(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00371 64* 2060,2065), NTYPE
00372 64* 2005 TM = (T(I)+LTA)/2.0
00373 64* 2007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))
00374 64* GO TO 2999
00375 64* 2010 TM = T(I)
00376 64* GO TO 2007
00377 64* 2015 CALL DID1WM(T(I),A(LA),XK(LK),61)
00400 64* 2017 J2 = J2+1
00401 64* LA = FLD(5,17,NSQ(J2))
00402 64* LK = FLD(22,14,NSQ(J2))
00403 64* CALL DID1WM(LTA),A(LA),XK(LK),62)
00404 64* GO TO 2998
00405 64*

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CNFAST,CNFAST

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00403 64* 2020 G1 = XK(LK)*XK(LA)
00404 64* GO TO 2017
00405 64* 2025 CALL D2DIWM(T(I),A(LA),XK(LK),G1)
00406 64* J2 = J2+1
00407 64* LA = FLD(5,17,NSQ2(J2))
00410 64* LK = FLD(22,14,NSQ2(J2))
00411 64* G2 = XK(LK)*XK(LA)
00412 64* GO TO 2998
00413 64* TM = (T(I)+T(LTA))/2.0
00414 64* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00415 64* GO TO 2999
00416 64* 2035 TM = T(I)
00417 64* GO TO 2032
00420 64* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00421 64* 2042 J2 = J2+1
00422 64* LA = FLD(5,17,NSQ2(J2))
00423 64* LK = FLD(22,14,NSQ2(J2))
00424 64* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00425 64* GO TO 2998
00426 64* 2045 G1 = XK(LK)*XK(LA)
00427 64* GO TO 2042
00430 64* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00431 64* J2 = J2+1
00432 64* LA = FLD(5,17,NSQ2(J2))
00433 64* LK = FLD(22,14,NSQ2(J2))
00434 64* G2 = XK(LK)*XK(LA)
00435 64* GO TO 2998
00436 64* 2055 TM = (T(I)+T(LTA))/2.0
00437 64* CALL D2DIWM(TM,CON(I4),A(LA),XK(LK),G(LG))
00440 64* GO TO 2999
00441 64* 2060 TM = T(LTA)
00442 64* GO TO 2007
00443 64* 2065 TM = T(LTA)
00444 64* GO TO 2032
00445 64* 2998 G(LG) = 1./(1./G1+1./G2)
00446 64* IF(FLD(3,1,NSQ1(J1)),EQ.1) G(LG) = G1*G2
00450 64* 2999 J2 = J2+1
00451 64* 3000 CONTINUE
00452 64* END
00453 65* IF(FLD(3,1,NSQ1(J1)),EQ.0) GO TO 55
00455 66* T1 = T(I)+460.0
00456 67* T2 = T(LTA)+460.0
00457 68* GV = G(LG)*(T1*T1+T2*T2)*(T1+T2)
00460 69* GO TO 60
00461 70* 55 GV = G(LG)
00462 71* 60 Q(I) = Q(I)+GV*T(LTA)
00463 72* X(LE) = X(LE)+GV
00464 73* IF(LTA.GT.RND.OR.FLD(21,1,NSQ1(J1)),EQ.1) GO TO 65
00466 74* LEA = IE+LTA
00467 75* X(LEA) = X(LEA)+GV
00470 76* Q(LTA) = Q(LTA)+GV*T(I)
00471 77* 65 IF(NSQ1(J1).GT.0) GO TO 50
00473 78* 85 CONTINUE
00475 79* CKM = 1.E+8
00476 80* TCGM = 0.0
00477 81* DO 105 I = 1,MND
00502 82* LE = IE+I
00503 83* T1 = C(I)/X(LE)
00504 84* IF(T1.GE.CKM) GO TO 90
00506 85* CKM = T1

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CNFAST,CNFAST

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00507 86* KON(35) = I
00510 87* 90 IF(TSTEP.GT.T1) GO TO 95
00512 88* T1 = T(1)+TSTEP*(Q11)-X(LE)*T(11)/C(11)
00513 89* GO TO 100
00514 90* 95 T1 = Q(1)/X(LE)
00515 91* 100 T2 = ABS(T1-T(1))
00516 92* T(1) = T1
00517 93* IF(T2.LI.TCGM) GO TO 105
00521 94* TCGM = T2
00522 95* KOH(15) = I
00523 96* 105 CONTINUE
00525 97* CON(15) = TCGM
00526 98* CON(17) = CKM
00527 99* IF(CKM.LE.0.) GO TO 996
00531 100* IF(MNA.LE.0) GO TO 160
00533 101* LAX = KON(5)
00534 102* DAMPN = CON(9)
00535 103* DAMPO = 1.0-DAMPN
00536 104* DO 150 I = 1,LAX
00541 105* KOH(20) = I
00542 106* RLX = 0.0
00543 107* JJ1 = J1
00544 108* JJ2 = J2
00545 109* DO 145 L = NN,MNC
00550 110* SUMC = 0.0
00551 111* SUMCV = 0.0
00552 112* IF(I.GT.1) GO TO 6000
00554 113* INCLUDE VR02,LIST
00555 113* IF(FLD(4,1,NS01(JJ1+1),EQ.0) GO TO 6000
00557 113* NTYPE = FLD(0,5,NS02(JJ2))
00560 113* LA = FLD(5,17,NS02(JJ2))
00561 113* LK = FLD(22,14,NS02(JJ2))
00562 113* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
00563 113* 5005 Q(L) = XK(ILK)+C(L)
00564 113* GO TO 5999
00565 113* 5010 Q1 = 0.0
00566 113* 5012 CALL DIDIWM(T(L),A(LA),XK(ILK),Q2)
00567 113* GO TO 5998
00570 113* 5015 Q1 = 0.0
00571 113* 5017 CALL DIDIWM(CON(14),A(LA),XK(ILK),Q2)
00572 113* GO TO 5998
00573 113* 5020 CALL DIDIWM(CON(14),A(LA),XK(ILK),Q1)
00574 113* 5022 JJ2 = JJ2+1
00575 113* LA = FLD(5,17,NS02(JJ2))
00576 113* LK = FLD(22,14,NS02(JJ2))
00577 113* GO TO 5017
00600 113* 5025 Q1 = XK(ILK)*XK(LA)
00601 113* GO TO 5022
00602 113* 5030 CALL DIDIWM(CON(14),A(LA),XK(ILK),Q1)
00603 113* JJ2 = JJ2+1
00604 113* LA = FLD(5,17,NS02(JJ2))
00605 113* LK = FLD(22,14,NS02(JJ2))
00606 113* Q2 = XK(ILK)*XK(LA)
00607 113* GO TO 5998
00610 113* 5035 CALL DIDIWM(CON(14),A(LA),XK(ILK),Q1)
00611 113* 5037 JJ2 = JJ2+1
00612 113* LA = FLD(5,17,NS02(JJ2))
00613 113* LK = FLD(22,14,NS02(JJ2))
00614 113* GO TO 5012
00615 113* 5040 Q1 = XK(ILK)*XK(LA)

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00616 113* GO TO 5037
 00617 113* 5998 0(L) = 01+02+0(L)
 00620 113* 5999 JJ2 = JJ2+1
 00621 113* 6000 CONTINUE
 00622 113* END
 00623 114* 110 JJ1 = JJ1+1
 00624 115* LG = FLD(5,16,NSQ1(JJ1))
 00625 116* LTA = FLD(22,14,NSQ1(JJ1))
 00626 117* IF(I.GT.1) GO TO 4000
 00630 118* INCLUDE VRG2,LIST
 00631 119* CHECK FOR RADIATION CONDUCTOR
 00633 119* IF(FLD(2,1,NSQ1(JJ1)).EQ.0) GO TO 4000
 00635 119* NTYPE = FLD(0,5,NSQ2(JJ2))
 00636 119* LA = FLD(5,17,NSQ2(JJ2))
 00637 119* LK = FLD(22,14,NSQ2(JJ2))
 00640 119* GOTO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,
 00641 119* 3060,3065), NTYPE
 00642 119* 3005 TM = (T(L)+T(LTA))/2.0
 00643 119* 3007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))
 00644 119* GO TO 3999
 00645 119* 3010 TM = T(L)
 00646 119* GO TO 3007
 00647 119* 3015 CALL DID1WM(T(L),A(LA),XK(LK),61)
 00650 119* 3017 JJ2 = JJ2+1
 00651 119* LA = FLD(5,17,NSQ2(JJ2))
 00652 119* LK = FLD(22,14,NSQ2(JJ2))
 00653 119* CALL DID1WM(T(LTA),A(LA),XK(LK),62)
 00654 119* GO TO 3998
 00655 119* 3020 G1 = XK(LK)*XK(LA)
 00656 119* GO TO 3017
 00657 119* 3025 CALL DID1WM(T(L),A(LA),XK(LK),61)
 00660 119* JJ2 = JJ2+1
 00661 119* LA = FLD(5,17,NSQ2(JJ2))
 00662 119* LK = FLD(22,14,NSQ2(JJ2))
 00663 119* G2 = XK(LK)*XK(LA)
 00664 119* GO TO 3998
 00665 119* 3030 TM = (T(L)+T(LTA))/2.0
 00666 119* 3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))
 00667 119* 3035 TM = T(L)
 00668 119* GO TO 3032
 00669 119* 3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
 00670 119* 3042 JJ2 = JJ2+1
 00671 119* LA = FLD(5,17,NSQ2(JJ2))
 00672 119* LK = FLD(22,14,NSQ2(JJ2))
 00673 119* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)
 00674 119* GO TO 3998
 00675 119* 3045 G1 = XK(LK)*XK(LA)
 00676 119* GO TO 3042
 00677 119* 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
 00700 119* JJ2 = JJ2+1
 00701 119* LA = FLD(5,17,NSQ2(JJ2))
 00702 119* LK = FLD(22,14,NSQ2(JJ2))
 00703 119* G2 = XK(LK)*XK(LA)
 00704 119* GO TO 3998
 00705 119* 3055 TM = (T(L)+T(LTA))/2.0
 00706 119* CALL DID1WM(TM,CON(14),A(LA),XK(LK),6(LG))
 00707 119* GO TO 3999
 00710 119* 3060 TM = T(LTA)
 00711 119* GO TO 3007

*NEW
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CNFAST,CNFAST

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00712 119*
00713 119*
00714 119*
00715 119*
00716 119*
00717 119*
00718 119*
00719 119*
00720 119*
00721 119*
00722 120*
00723 121*
00724 122*
00725 122*
00726 123*
00727 124*
00728 124*
00729 125*
00730 125*
00731 126*
00732 127*
00733 128*
00734 129*
00735 130*
00736 131*
00737 132*
00738 133*
00739 134*
00740 134*
00741 135*
00742 135*
00743 136*
00744 137*
00745 138*
00746 138*
00747 139*
00748 140*
00749 141*
00750 142*
00751 143*
00752 143*
00753 144*
00754 145*
00755 146*
00756 147*
00757 148*
00758 149*
00759 150*
00760 151*
00761 152*
00762 153*
00763 154*
00764 154*
00765 155*
00766 156*
00767 157*
00768 157*
00769 158*
00770 159*
00771 160*
00772 161*
00773 162*
00774 163*
00775 164*
00776 165*
00777 166*
00778 167*
00779 168*

3065 TM = T(LTA)
    GO TO 3032
3998 G(LG) = 1./(1./G1+1./G2)
    IF(FLD(3,1,NSQ1(JJ1)).EQ.1) 6(L6) = 61*62
3999 JJ2 = JJ2+1
4000 CONTINUE
    END
    IF(FLD(3,1,NSQ1(JJ1)).EQ.0) GO TO 115
    T1 = T(L)+460.0
    T2 = T(LTA)+460.0
    GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
    GO TO 120
115 GV = G(LG)
    T2 = T(LTA)
120 SUMC = SUMC+GV
    SUMCV = SUMCV+GV*T2
    CHECK FOR LAST CONDUCTOR TO THIS NODE
    IF(NSQ1(JJ1).GT.0) GO TO 110
    T1 = DAMPN*(SUMCV+G(L))/SUMC+DAMP0*T(L)
    T2 = ABS(T(L)-T1)
    IF(RLX.GE.T2) GO TO 140
    RLX = T2
    KON(37) = L
140 T(L) = T1
145 CONTINUE
    IF(RLX.LE.CON(19)) GO TO 155
150 CONTINUE
155 CON(30) = RLX
160 CALL VARBL2
    CON(13) = CON(1)
    TSUM = TSUM+TSTEP
    TSTEP = CKM
    IF(TSUM.LT.CON(18)) GO TO 10
    TPRINT = TPRINT+TSUM
    CALL OUTCAL
    IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
    NTH = IE
    NDIM = NLA
    RETURN
995 WRITE(6,885)
    GO TO 1000
996 WRITE(6,886)
    GO TO 1000
997 WRITE(6,887) NDIM
    GO TO 1000
998 WRITE(6,888)
    GO TO 1000
999 WRITE(6,889)
1000 CALL OUTCAL
    CALL EXIT
885 FORMAT(4H CNFAST REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
886 FORMAT(22H C/SK ZERO OR NEGATIVE)
887 FORMAT(18,20H LOCATIONS AVAILABLE)
888 FORMAT(10H NO DTIMEL)
889 FORMAT(19H NO OUTPUT INTERVAL)
    END

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END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
CNFAST SYMBOLIC
CNFAST CODE RELOCATABLE

21 FEB 71

CNEXPN,CNEXPN

QIW FOR,* CNEXPN,CNEXPN
UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-10D CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:19

SUBROUTINE CNEXPN ENTRY POINT 003536

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	003551
0000	*COST+TEMP	000072
0002	*SIMPLE VAR	000047
0004	*ARWAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COND	000001
0013	PCI	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARBL1
0023	DID1WM
0024	PLYAWM
0025	D2D1WM
0026	VARBL2
0027	OUTCAL
0030	EX11
0031	MERR25
0032	EXP
0033	NWDUS
0034	NI025
0035	NER105

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C
0002	R	000022	C2	0002	R	000040	DD
0002	R	000034	GV	0002	R	000030	G1
0002	I	000002	IE	0002	I	000042	JJ1
0015	I	000000	K	0017	I	000000	KON
0002	I	000011	LE	0002	I	000035	LEA
0021	I	000007	LSQ2	0002	I	000026	LTA
0020	I	000000	MDIM	0021	I	000003	MGT
0021	I	000000	N40	0021	I	000002	NMT
0002	I	000016	NTYPE	0020	I	000002	FIX
0002	R	000024	Q2	0002	R	000045	SUMC
0002	R	000027	TM	0002	R	000007	TPRINT
0002	R	000021	C1	0002	R	000015	CKM
0012	R	000000	G	0002	R	000036	DELTA
0002	I	000004	I	0002	R	000031	G2
0002	I	000013	J2	0002	I	000043	JJ2
0002	I	000041	LAX	0002	I	000044	L
0021	I	000006	LSQ1	0002	I	000025	LG
0021	I	000004	NCT	0002	I	000005	L1
0002	I	000001	NNA	0002	I	000003	NLA
0020	I	000001	NTH	0013	I	000000	NSQ1
0002	R	000023	O1	0014	I	000000	NSQ2
0002	R	000014	TCGM	0011	R	000000	Q
0002	R	000032	T1	0007	R	000000	T
				0002	R	000010	TSUM
				0017	R	000000	CON
				0002	R	000037	DN
				0006	R	000000	H
				0002	I	000012	J1
				0002	I	000017	LA
				0002	I	000020	LK
				0021	I	000005	NAT
				0021	I	000001	NNA
				0014	I	000000	NSQ2
				0011	R	000000	Q
				0007	R	000000	T
				0002	R	000010	TSUM

CNEXPN.CNEXPN

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0002 R 000033 Y2      0020 R 000002 X      0015 R 006000 XK      000124 IBL      003916 I000L
0001 000326 1005L    0001 000347 1010L    0001 003352 10116    000126 1012L    000420 1015L
0001 000426 1020L    0001 000466 1025L    0001 006512 1030L    000126 1032L    000471 1035L
0001 000577 1040L    0001 000642 1045L    0001 002105 105L    000126 105L    000571 105L
0001 002156 125L    0001 002164 130L    0001 002514 135L    000126 140L    000621 120L
0001 000133 15L     0001 003332 165L    0001 003343 175L    000126 145L    000671 145L
0001 003434 190L    0001 003437 195L    0001 000664 1998L    000126 150L    000721 150L
0001 000672 2000L   0001 001234 2005L    0001 001241 2007L    000126 155L    000771 155L
0001 001302 2017L   0001 001336 2020L    0001 001344 2025L    000126 160L    000821 160L
0001 001434 2035L   0001 001437 2040L    0001 001460 2042L    000126 165L    000871 165L
0001 001570 2055L   0001 001616 2060L    0001 001622 2065L    000126 170L    000921 170L
0001 000145 25L     0001 001626 2998L    0001 001654 2999L    000126 175L    000971 175L
0001 002602 3005L   0001 002607 3007L    0001 002627 3010L    000126 180L    001021 180L
0001 002704 3020L   0001 002712 3025L    0001 002752 3030L    000126 185L    001071 185L
0001 003005 3040L   0001 003026 3042L    0001 003065 3045L    000126 190L    001121 190L
0001 003164 3060L   0001 003170 3065L    0001 003174 3068L    000126 195L    001171 195L
0001 000737 4005L   0001 000744 4010L    0001 000745 4012L    000126 200L    001221 200L
0001 001002 4020L   0001 001016 4022L    0001 001033 4025L    000126 205L    001271 205L
0001 001113 4037L   0001 001130 4040L    0001 000236 45L    000126 210L    001321 210L
0001 000112 5L      0001 001145 50L    0001 001145 5000L    000126 215L    001371 215L
0001 002314 5012L   0001 002333 5015L    0001 002334 5017L    000126 220L    001421 220L
0001 002402 5025L   0001 002402 5036    0001 002410 5030L    000126 225L    001471 225L
0001 002477 5040L   0001 002111 5356    0001 001714 55L    000126 230L    001521 230L
0001 002511 5999L   0001 001717 60L    0001 002514 6000L    000126 235L    001571 235L
0001 001774 85L     0001 000000 885F    0000 000011 886F    000126 240L    001621 240L
0000 000030 889F    0001 002023 90L    0001 002061 95L    000126 245L    001671 245L
0001 003472 997L    0001 003500 998L    0001 003507 999L    000126 250L    001721 250L

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00101 1* C SUBROUTINE CNEXPN
00101 2* C EXPLICIT EXPONENTIAL DIFFERENCING ROUTINE FOR SINDA FORTRAN V
00101 3* C THE SHORT PSEUDO-COMPUTE SEQUENCE IS REQUIRED
00103 4* C INCLUDE COMM,LIST
00104 4* C COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/O(1)
00105 4* C COMMON /PC1/NSG(1) /PC2/NSG2(1) /KONST/K(1) /ARRAY/A(1)
00106 4* C COMMON /FIXCON/KON(1) /XSPACE/NDIM,HTH,X(1)
00107 4* C COMMON /DIMENS/ NND,NNA,NNT,NGT,NCT,NAT,LSO1,LSO2
00110 4* C DIMENSION CON(1),XK(1),NX(1)
00111 4* C EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
00112 4* C
00113 5* C INCLUDE DEFF,LIST
00113 5* C ***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00113 5* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
00113 5* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00113 5* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEHD)
00113 5* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR, EXPLICIT (CSGFAC)
00113 5* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00113 5* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00113 5* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
00113 5* C CCA CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
00113 5* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPFA)
00113 5* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
00113 5* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
00113 5* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUPC)
00113 5* C CC13 CONTAINS THE PRESET TIME OR PROBLEM START TIME (TIMEO)
00113 5* C CC14 CONTAINS THE NEAR TIME BETWEEN AN ITERATION (TIMEM)
00113 5* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)
00113 5* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)

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CNEXPX,CNEXPX

00113 C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSGMIN)
00113 C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (CSGMAX)
00113 C CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
00113 C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
00113 C CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
00113 C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINCT)
00113 C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 C CC31 IS INDICATOR, 0=THERMAL SPCS,1=THERMAL LPCS,2=GENERAL (LSPCS)
00113 C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (EPGBAL)
00113 C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 C CC50 IS NOT USED AT PRESENT
00114 EVID
00115 IF(CON(4).LE.0.0) CON(4) = 1.0
00117 IF(KON(5).LE.0) KON(5) = 1
00121 IF(CON(6).LE.0.) CON(6) = 1.E+8
00123 IF(CON(8).LE.0.) CON(8) = 1.E+8
00125 IF(CON(9).LE.0.) CON(9) = 1.0
00127 IF(CON(11).LE.0.) CON(11) = 1.E+8
00131 IF(CON(1A).LE.0.) GO TO 999
00133 IF(CON(19).LE.0.) CON(19) = 1.E+8
00135 IF(KON(31).NE.0) GO TO 995
00137 PASS = -1.0
00140 NNC = I*ND+NNA
00141 IE = NTH
00142 NLA = NDIM
00143 NTH = NTH+NNC
00144 NDIM = NDIM+NNC
00144 C CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00145 I = NLA+NNC
00146 IF(I.LT.0) GO TO 998
00150 L1 = NID+1
00151 TSTEP = CON(18)
00152 TPRINT = CON(13)
00152 C INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00153 TSUM = 0.0
00153 C DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00154 IF(CON(13)+CON(18).LE.CON(3)) GO TO 10
00154 C DONT EXCEED IT
00156 CON(18) = CON(13)-CON(13)
00156 C IS THE TIME STEP LARGER THAN ALLOWED
00157 IF(TSTEP.LE.CON(18)) GO TO 15
00161 TSTEP = CON(18)
00161 C DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
00162 IF(TSUM+TSTEP>CON(18)) 25,30,20
00162 C DONT EXCEED IT

CNEXPN,CNEXPN

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00165 39* 20 TSTEP = CON(18)-TSUM
00166 40* GO TO 30
00167 41* C DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
00168 42* 25 IF(TSUM*2.0+TSTEP*LE.CON(18)) GO TO 30
00169 43* C APPROACH THE OUTPUT INTERVAL GRADUALLY
00170 44* TSTEP = (CON(18)-TSUM)/2.0
00171 45* C STORE DELTA TIME STEP IN THE CONSTANTS
00172 46* 30 CON(2) = TSTEP
00173 47* C IS THE TIME STEP USED LESS THAN THE TIME STEP ALLOWED
00174 48* IF(TSTEP.LT.CON(21)) GO TO 997
00175 49* C CALCULATE THE NEW TIME
00176 50* CON(1) = TPRINT+TSUM+TSTEP
00177 51* C COMPUTE THE MEAN TIME BETWEEN ITERATIONS
00178 52* CON(14) = (CON(1)+CON(13))/2.0
00179 53* C ZERO OUT ALL SOURCE LOCATIONS AND EXTRA LOCATIONS
00180 54* DO 35 I = 1,NND
00181 55* LE = IE+1
00182 56* X(LE) = 0.0
00183 57* Q(I) = 0.0
00184 58* 35 CONTINUE
00185 59* C SHIFT THE ARITHMETIC TEMPERATURES INTO THE EXTRA LOCATIONS
00186 60* IF(INNALE.0) GO TO 45
00187 61* DO 40 I = 1,NND
00188 62* Q(I) = 0.0
00189 63* LE = IE+1
00190 64* X(LE) = T(I)
00191 65* 40 CONTINUE
00192 66* 45 KORN(12) = 0
00193 67* CALL VARBL1
00194 68* IF(KORN(12).NE.0) GO TO 10
00195 69* J1 = 0
00196 70* J2 = 1
00197 71* TCGM = 0.0
00198 72* CKM = 1.E+8
00199 73* C CALCULATE Q SUM AND G SUM
00200 74* DO 85 I = 1,NND
00201 75* LE = IE+1
00202 76* INCLUDE VARC,LIST
00203 76* IF(FLD(1,1,NS01(J1+1)).EQ.0) GO TO 2000
00204 76* NTYPE = FLD(10,5,NS02(J2))
00205 76* LA = FLD(5,17,NS02(J2))
00206 76* LK = FLD(22,14,NS02(J2))
00207 76* GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
00208 76* 1005 CALL DID1WM(T(I),A(LA),XK(LK),C(I))
00209 76* GO TO 1999
00210 76* 1010 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00211 76* 1012 J2 = J2+1
00212 76* LA = FLD(5,17,NS02(J2))
00213 76* LK = FLD(22,14,NS02(J2))
00214 76* CALL DID1WM(T(I),A(LA),XK(LK),C2)
00215 76* GO TO 1998
00216 76* 1015 C1 = XK(LK)*XK(LA)
00217 76* GO TO 1012
00218 76* 1020 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00219 76* J2 = J2+1
00220 76* LA = FLD(5,17,NS02(J2))
00221 76* LK = FLD(22,14,NS02(J2))
00222 76* C2 = XK(LK)*XK(LA)
00223 76* GO TO 1998
00224 76* 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))

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*NEW
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00265 76* CNEXPN,CNEXPN
00266 76* GO TO 1999
00267 76* 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),CI)
00270 76* 1032 J2 = J2+1
00271 76* LA = FLD(5,17,NSQ2(J2))
00272 76* LK = FLD(22,14,NSQ2(J2))
00273 76* CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00274 76* GO TO 1998
00275 76* C1 = XK(LK)*XK(LA)
00276 76* GO TO 1032
00277 76* 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),CI)
00300 76* J2 = J2+1
00301 76* LA = FLD(5,17,NSQ2(J2))
00302 76* LK = FLD(22,14,NSQ2(J2))
00303 76* C2 = XK(LK)*XK(LA)
00304 76* GO TO 1998
00305 76* 1045 CALL D2D1WM(T(I),CON(14),A(LA),XK(LK),C(I))
00306 76* 1998 C(I) = C1+C2
00307 76* 1999 J2 = J2+1
00310 76* 2000 CONTINUE
00311 76* END
00312 77* INCLUDE VARG,LIST
00313 77* IF(FLD(4,1,NSQ1(J1+1)),EQ,0) GO TO 5000
00315 77* NTYPE = FLD(0,5,NSQ2(J2))
00316 77* LA = FLD(5,17,NSQ2(J2))
00317 77* LK = FLD(22,14,NSQ2(J2))
00320 77* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE
00321 77* 4005 G(I) = XK(LK)*G(I)
00322 77* GO TO 4999
00323 77* 4010 G1 = 0.0
00324 77* 4012 CALL D1D1WM(T(I),A(LA),XK(LK),G2)
00325 77* GO TO 4998
00326 77* 4015 G1 = 0.0
00327 77* 4017 CALL D1D1WM(CON(14),A(LA),XK(LK),G2)
00330 77* GO TO 4998
00331 77* 4020 CALL D1D1WM(CON(14),A(LA),XK(LK),G1)
00332 77* 4022 J2 = J2+1
00333 77* LA = FLD(5,17,NSQ2(J2))
00334 77* LK = FLD(22,14,NSQ2(J2))
00335 77* GO TO 4017
00336 77* 4025 G1 = XK(LK)*XK(LA)
00337 77* GO TO 4022
00340 77* 4030 CALL D1D1WM(CON(14),A(LA),XK(LK),G1)
00341 77* J2 = J2+1
00342 77* LA = FLD(5,17,NSQ2(J2))
00343 77* LK = FLD(22,14,NSQ2(J2))
00344 77* G2 = XK(LK)*XK(LA)
00345 77* GO TO 4998
00346 77* 4035 CALL D1D1WM(CON(14),A(LA),XK(LK),G1)
00347 77* 4037 J2 = J2+1
00350 77* LA = FLD(5,17,NSQ2(J2))
00351 77* LK = FLD(22,14,NSQ2(J2))
00352 77* GO TO 4012
00353 77* 4040 G1 = XK(LK)*XK(LA)
00354 77* GO TO 4037
00355 77* 4998 G(I) = G1+G2+G(I)
00356 77* 4999 J2 = J2+1
00357 77* 5000 CONTINUE
00360 77* END
00361 78* 50 J1 = J1+1

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CNEXPN,CNEXPN

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00362 79* L6 = FLD(5,16,NS01(J1))
00363 80* IF(L6.EQ.0) GO TO 85
00365 81* LTA = FLD(22,14,NS01(J1))
00366 82* INCLUDE VARG-LIST
00366 83* CHECK FOR RADIATION CONDUCTOR
00367 83* IF(FLD(2,1,NS01(J1)).EQ.0) GO TO 5000
00371 83* NTYPE = FLD(10,5,NS02(J2))
00372 83* LA = FLD(5,17,NS02(J2))
00373 83* LK = FLD(22,14,NS02(J2))
00374 83* GOT0(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00374 83* 2060,2065), NTYPE
00375 83* TM = (T(I)+T(LTA))/2.0
00376 83* 2007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
00377 83* GO TO 2999
00400 83* 2010 TM = T(I)
00401 83* GO TO 2007
00402 83* 2015 CALL DID1WM(T(I),A(LA),XK(LK),G(LG))
00403 83* 2017 J2 = J2+1
00404 83* LA = FLD(5,17,NS02(J2))
00405 83* LK = FLD(22,14,NS02(J2))
00406 83* CALL DID1WM(T(LTA),A(LA),XK(LK),G(LG))
00407 83* GO TO 2998
00410 83* 2020 G1 = XK(LK)*XK(LA)
00411 83* GO TO 2017
00412 83* 2025 CALL DID1WM(T(I),A(LA),XK(LK),G(LG))
00413 83* J2 = J2+1
00414 83* LA = FLD(5,17,NS02(J2))
00415 83* LK = FLD(22,14,NS02(J2))
00416 83* G2 = XK(LK)*XK(LA)
00417 83* GO TO 2998
00420 83* 2030 TM = (T(I)+T(LTA))/2.0
00421 83* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00422 83* GO TO 2999
00423 83* 2035 TM = T(I)
00424 83* GO TO 2032
00425 83* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G(LG))
00426 83* 2042 J2 = J2+1
00427 83* LA = FLD(5,17,NS02(J2))
00430 83* LK = FLD(22,14,NS02(J2))
00431 83* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G(LG))
00432 83* GO TO 2998
00433 83* 2045 G1 = XK(LK)*XK(LA)
00434 83* GO TO 2042
00435 83* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G(LG))
00436 83* J2 = J2+1
00437 83* LA = FLD(5,17,NS02(J2))
00440 83* LK = FLD(22,14,NS02(J2))
00441 83* G2 = XK(LK)*XK(LA)
00442 83* GO TO 2998
00443 83* 2055 TM = (T(I)+T(LTA))/2.0
00444 83* CALL D2D1WM(TM,CON(14),A(LA),XK(LK),G(LG))
00445 83* GO TO 2999
00446 83* 2060 TM = T(LTA)
00447 83* GO TO 2007
00450 83* 2065 TM = T(LTA)
00451 83* GO TO 2032
00452 83* 2998 G(LG) = 1./(1./G1+1./G2)
00453 83* IF(FLD(3,1,NS01(J1)).EQ.1) G(LG) = G1*G2
00455 83* 2999 J2 = J2+1
00456 83* 3000 CONTINUE

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CNEXPN,CNEXPN

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00457 83* END
00460 84* IF(FLD(3,1,NSQ1(J1)),EQ,0) GO TO 55
00462 85* T1 = T(I)+460.0
00463 86* T2 = T(LTA)+460.0
00464 87* GV = G(LG)*(T1+T2)*(T1+T2)
00465 88* GO TO 60
00466 89* 55 GV = G(LG)
00467 90* C OBTAIN THE G RATE THRU THE CONDUCTOR
00468 91* G(I) = G(I)+GV*(LTA)
00469 92* C SAVE SUMMATION OF CONDUCTORS
00470 93* X(LE) = X(LE)+GV
00471 94* C CHECK FOR ADJOINING DIFFUSION NODE
00472 95* IF(LTA.GT.NMD.OR.FLD(21,1,NSQ1(J1)),EQ,1) GO TO 65
00473 96* C SAVE SUMMATION OF CONDUCTORS FOR ADJOINING NODE
00474 97* LEA = IE+LTA
00475 98* X(LEA) = X(LEA)+GV
00476 99* G(LTA) = G(LTA)+GV*(I)
00477 100* C CHECK FOR LAST CONDUCTOR
00478 101* 65 IF(NSQ1(J1).GT,0) GO TO 50
00479 102* 85 CONTINUE
00480 103* C OBTAIN NEW DIFFUSION TEMPERATURES, DTMPC AND CSGMIN
00481 104* DO 100 I = 1,NND
00482 105* LE = IE+I
00483 106* C CALCULATE C/SK MINIMUM
00484 107* T1 = C(I)/X(LE)
00485 108* IF(T1.GE.CKM) GO TO 90
00486 109* CKM = T1
00487 110* KON(35) = I
00488 111* C COMPUTE NEW TEMPERATURES USING CALCULATED SOURCE TERMS
00489 112* T2 = 1.0/EXP(TSTEP*X(LE)/C(I))
00490 113* T1 = ((1.0-T2)*Q(I)/X(LE)+T2*T(I))
00491 114* C CALCULATE THE ABSOLUTE VALUE TEMPERATURE CHANGE
00492 115* T2 = ABS(T1-T(I))
00493 116* C SAVE THE LARGEST TEMPERATURE CHANGE
00494 117* IF(TCGM.GE.T2) GO TO 95
00495 118* TCGM = T2
00496 119* KON(36) = I
00497 120* C STORE THE TEMPERATURES
00498 121* 95 X(LE) = T(I)
00499 122* T(I) = T1
00500 123* 100 CONTINUE
00501 124* CON(17) = CKM
00502 125* DELTA = CKM*CON(4)
00503 126* IF(CKM.LE.0.0) GO TO 996
00504 127* C CHECK FOR FIRST PASS
00505 128* IF(PASS.GT,0.0) GO TO 115
00506 129* C UNDO THE TEMPERATURE CALCULATIONS
00507 130* 105 DO 110 I = 1,NMC
00508 131* LE = IE+I
00509 132* T(I) = X(LE)
00510 133* 110 CONTINUE
00511 134* IF(PASS.GT,0.0) GO TO 15
00512 135* PASS = 1.0
00513 136* CON(1) = TPRINT
00514 137* CON(2) = 0.0
00515 138* TSTEP = DELTA*0.95
00516 139* GO TO 195
00517 140* C IS THE TIME STEP USED LESS THAN THE TIME STEP CALCULATED
00518 141* 115 IF(TSTEP.LE.DELTA) GO TO 130
00519 142* C COMPUTE THE TIME STEP

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CNEXPN,CNEXPN

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00601 164*
00604 165*
00605 166*
00606 167*
00610 168*
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00653 168*

TSTEP = DELTA*0.95
GO TO 105
120 TSTEP = 0.95*TSTEP*CON(61)/TCGM
GO TO 105
125 TSTEP = 0.95*TSTEP*CON(11)/TCGM
GO TO 105
C SEE IF THE TEMPERATURE CHANGE WAS TOO LARGE
130 IF(TCGM.GT.CON(6)) GO TO 120
C STORE THE MAXIMUM DIFFUSION TEMPERATURE CHANGE
CON(15) = TCGM
C CHECK TO SEE IF THERE ARE ANY ARITHMETIC NOTES
IF(RNA.LE.0) GO TO 185
C COMPUTE ARITHMETIC TEMPERATURES BY SUCCESSIVE POINT OVER-RELAX
DN = CON(9)
DD = 1.0-DN
LAX = KON(5)
DO 170 I = 1,LAX
JJ1 = J1
JJ2 = J2
TCGM = 0.0
KON(20) = I
DO 165 L = 1,NNC
SUMC = 0.0
SUMCV = 0.0
IF(I.GT.1) GO TO 6000
INCLUDE VR02,LIST
IF(FLD(4),NSQ(JJ1+1),EQ.0) GO TO 6000
NTYPE = FLD(0,5,NSQ(JJ2))
LA = FLD(15,17,NSQ(JJ2))
LK = FLD(22,14,NSQ(JJ2))
GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
5005 0(L) = XK(LK)+0(L)
GO TO 5999
5010 01 = 0.0
5012 CALL DIDIWM(T(L),A(LA),XK(LK),02)
GO TO 5998
5015 01 = 0.0
5017 CALL DIDIWM(CON(14),A(LA),XK(LK),02)
GO TO 5998
5020 CALL DIDIWM(CON(14),A(LA),XK(LK),01)
5022 JJ2 = JJ2+1
LA = FLD(15,17,NSQ(JJ2))
LK = FLD(22,14,NSQ(JJ2))
GO TO 5017
5025 01 = XK(LK)*XK(LA)
GO TO 5022
5030 CALL DIDIWM(CON(14),A(LA),XK(LK),01)
JJ2 = JJ2+1
LA = FLD(15,17,NSQ(JJ2))
LK = FLD(22,14,NSQ(JJ2))
02 = XK(LK)*XK(LA)
GO TO 5998
5035 CALL DIDIWM(CON(14),A(LA),XK(LK),01)
5037 JJ2 = JJ2+1
LA = FLD(15,17,NSQ(JJ2))
LK = FLD(22,14,NSQ(JJ2))
GO TO 5012
5040 01 = XK(LK)*XK(LA)
GO TO 5037
5998 0(L) = 01+02+0(L)

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CNEXPN,CNEXPN

00654 168* 5999 JJ2 = JJ2+1
00655 168* 6000 CONTINUE
00656 168* END
00657 169* 135 JJ1 = JJ1+1
00660 170* LG = FLD(5,16,NSQ1(JJ1))
00661 171* LTA = FLD(22,14,NSQ1(JJ1))
00662 172* IF(1.GT.1) GO TO 4000
00664 173* INCLUDE VRG2,LIST
00664 174* CHECK FOR RADIATION CONDUCTOR
00665 174* IF(FLD(2,1,NSQ1(JJ1)).EQ.0) GO TO 4000
00667 174* NTYPE = FLD(10,5,NSQ2(JJ2))
00670 174* LA = FLD(5,17,NSQ2(JJ2))
00671 174* LK = FLD(22,14,NSQ2(JJ2))
00672 174* GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,
3060,3065), NTYPE
00673 174* 3005 TM = (T(L)+T(LTA))/2.0
00674 174* 3007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))
00675 174* GO TO 3999
00676 174* 3010 TM = T(L)
00677 174* GO TO 3007
00700 174* 3015 CALL DID1WM(T(L),A(LA),XK(LK),61)
00701 174* 3017 JJ2 = JJ2+1
00702 174* LA = FLD(5,17,NSQ2(JJ2))
00703 174* LK = FLD(22,14,NSQ2(JJ2))
00704 174* CALL DID1WM(T(LTA),A(LA),XK(LK),62)
00705 174* GO TO 3998
00706 174* 3020 G1 = XK(LK)*XK(LA)
00707 174* GO TO 3017
00710 174* 3025 CALL DID1WM(T(L),A(LA),XK(LK),61)
00711 174* JJ2 = JJ2+1
00712 174* LA = FLD(5,17,NSQ2(JJ2))
00713 174* LK = FLD(22,14,NSQ2(JJ2))
00714 174* G2 = XK(LK)*XK(LA)
00715 174* GO TO 3998
00716 174* 3030 TM = (T(L)+T(LTA))/2.0
00717 174* 3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))
00720 174* GO TO 3999
00721 174* 3035 TM = T(L)
00722 174* GO TO 3032
00723 174* 3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
00724 174* 3042 JJ2 = JJ2+1
00725 174* LA = FLD(5,17,NSQ2(JJ2))
00726 174* LK = FLD(22,14,NSQ2(JJ2))
00727 174* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)
00730 174* GO TO 3998
00731 174* 3045 G1 = XK(LK)*XK(LA)
00732 174* GO TO 3042
00733 174* 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)
00734 174* JJ2 = JJ2+1
00735 174* LA = FLD(5,17,NSQ2(JJ2))
00736 174* LK = FLD(22,14,NSQ2(JJ2))
00737 174* G2 = XK(LK)*XK(LA)
00740 174* GO TO 3998
00741 174* 3055 TM = (T(L)+T(LTA))/2.0
00742 174* CALL D2D1WM(TM,CON(14),A(LA),XK(LK),6(LG))
00743 174* GO TO 3999
00744 174* 3060 TM = T(LTA)
00745 174* GO TO 3007
00746 174* 3065 TM = T(LTA)
00747 174* GO TO 3032

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CNEXPN,CNEXPN

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00750 174* 3998 G(LG) = 1./(1./G1+1./G2)
00751 174* IF(FLD(3,1,NS01(JJ1)).EQ.1) G(LG) = G1*G2
00753 174* 3999 JJ2 = JJ2+1
00754 174* 4000 CONTINUE
00755 174* END
00756 175* IF(FLD(3,1,NS01(JJ1)).EQ.0) GO TO 140
00760 176* T1 = T(L)+460.0
00761 177* T2 = T(LTA)+460.0
00762 178* GV = G(LG)*(T1+T1+T2*T2)*(T1+T2)
00763 179* GO TO 145
00764 180* 140 GV = G(LG)
00765 181* 145 SUMC = SUMC+GV
00766 182* SUMCV = SUMCV+GV*T(LTA)
00767 183* CHECK FOR LAST CONDUCTOR
00771 185* IF(NS01(JJ1).GT.0) GO TO 135
00771 186* T2 = DD*(L)+DH*(SUMCV+G(L))/SUMC
00772 187* T1 = ABS(T(L)-T2)
00772 188* STORE THE NEW TEMPERATURE
00773 189* T(L) = T2
00773 190* C SAVE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
00774 191* IF(TCGM.GE.T1) GO TO 165
00776 192* TCGM = T1
00777 193* KON(37) = L
01000 194* 165 CONTINUE
01000 195* C SEE IF RELAXATION CRITERIA WAS MET
01002 196* IF(TCGM.LE.CON(19)) GO TO 175
01004 197* 170 CONTINUE
01004 198* C STORE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
01006 199* TCGM = TCGM
01007 200* C COMPUTE THE ARITHMETIC TEMPERATURE CHANGE
01010 201* TCGM = 0.0
01010 202* DO 180 I = L1,NNC
01013 203* LE = IE+1
01014 204* T1 = ABS(T(I)-X(LE))
01015 205* IF(T1.LT.TCGM) GO TO 180
01017 206* TCGM = T1
01020 207* KON(38) = I
01021 208* 180 CONTINUE
01021 209* C SEE IF ATMPCA WAS SATISFIED
01023 210* IF(TCGM.GT.CON(11)) GO TO 125
01025 211* CON(16) = TCGM
01026 212* 185 KON(12) = 0
01027 213* CALL VARBL2
01027 214* C CHECK THE BACKUP SWITCH
01030 215* IF(KON(12).NE.0) GO TO 105
01030 216* C ADVANCE TIME
01032 217* CON(13) = CON(11)
01033 218* TSUM = TSUM+TSTEP
01034 219* TSTEP = DELTA*0.95
01034 220* C CHECK FOR TIME TO PRINT
01035 221* IF(TSUM.GE.CON(18)) GO TO 190
01035 222* C CHECK FOR PRINT EVERY ITERATION
01037 223* IF(KON(7).EQ.0) GO TO 10
01041 224* CALL OUTCAL
01042 225* GO TO 10
01042 226* C TRY TO EVEN THE OUTPUT INTERVALS
01043 227* 190 TPRINT = TPRINT+TSUM
01044 228* 195 CALL OUTCAL
01044 229* C 15 TIME GREATER THAN END COMPUTE TIME

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CNEXPX,CNEXPX

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01045 230*
01047 231*
01050 232*
01051 233*
01052 234*
01054 235*
01055 236*
01057 237*
01060 238*
01062 239*
01063 240*
01066 241*
01067 242*
01071 243*
01072 244*
01073 245*
01074 246*
01075 247*
01076 248*
01077 249*
01100 250*

IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
NTH = IE
NDIM = NLA
RETURN
      995 WRITE(6,885)
        GO TO 1000
      996 WRITE(6,886)
        GO TO 1000
      997 WRITE(6,887)
        GO TO 1000
      998 WRITE(6,888) I
        GO TO 1000
      999 WRITE(6,889)
        GO TO 1000
      1000 CALL OUTCAL
           CALL EXIT
      885 FORMAT(46H CNEXPX REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
      886 FORMAT(24H CSG*IN ZERO OR NEGATIVE)
      887 FORMAT(20H TIME STEP TOO SMALL)
      888 FORMAT(18,20H LOCATIONS AVAILABLE)
      889 FORMAT(19H NO OUTPUT INTERVAL)
           END

```

END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
 CNEXPX SYMBOLIC
 CNEXPX CODE RELOCATABLE

91 FOR: CNDUFR
UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-100 CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 23:15:00

18 FEB 71

SUBROUTINE CNDUFR ENTRY POINT 003620

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	003635
0000	*CONST+TEMP	000100
0002	*SIMPLE VAR	000054
0004	*ARHAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COND	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARBL1
0023	DIDJWM
0024	PLYAWM
0025	D2DIWM
0026	VARBL2
0027	OUTCAL
0030	EXIT
0031	NERR2\$
0032	NWDUS
0033	NIQ2\$
0034	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C
0002	R	000024	C2	0002	R	000043	DELTA
0002	R	000041	DT2	0002	R	000036	GV
0006	R	000000	H	0002	I	000002	IE
0002	I	000050	JJ2	0002	I	000015	J2
0002	I	000051	L	0002	I	000046	LAX
0002	I	000042	LEH	0002	I	000022	LK
0002	I	000030	LTA	0021	I	000005	NAT
0021	I	000003	NGT	0021	I	000001	NNA
0021	I	000002	NNT	0014	I	000000	NSQ2
0020	R	000002	NX	0011	R	000000	Q
0002	R	000052	SUMC	0007	R	000000	T
0002	R	000011	TFRINT	0002	R	000010	TSTEPO
0002	R	000000	A	0002	R	000017	CKM
0002	R	000040	DT1	0002	R	000044	DN
0002	R	000033	G2	0002	R	000032	G1
0002	I	000047	JJ1	0002	I	000003	IEM
0017	I	000000	KON	0015	I	000000	K
0002	I	000037	LEA	0002	I	000013	LE
0021	I	000007	LSQ2	0002	I	000006	LSQ1
0020	I	000000	NDIM	0021	I	000004	NCT
0021	I	000000	NDIM	0002	I	000001	NMC
0002	R	000026	O2	0020	I	000001	NTH
0002	R	000031	T1	0002	R	000025	O1
0002	R	000034	T1	0002	R	000016	TCGM
0002	R	000023	C1	0002	R	000000	CON

CNDUFR

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00113 5* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSGMIN)
00113 5* C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 5* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 5* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 5* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 5* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (CSGMAX)
00113 5* C CC23 CONTAINS THE C/SG MAXIMUM (CSGRAL)
00113 5* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRCL)
00113 5* C CC25 CONTAINS THE C/SG RANGE CALCULATION (DRLXCA)
00113 5* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCC)
00113 5* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (LIMECT)
00113 5* C CC28 CONTAINS THE LINE COUNTER, INTEGER (PAGECT)
00113 5* C CC29 CONTAINS THE ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 5* C CC30 IS INDICATOR, 0=THERMAL SPCS,1=THERMAL LPCS,2=GENERAL (LSPCS)
00113 5* C CC31 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGRAL)
00113 5* C CC32 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 5* C CC33 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 5* C CC34 CONTAINS RELATIVE NODE NUMBER OF DTMPCC
00113 5* C CC35 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 5* C CC36 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC38 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 5* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 5* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 5* C CC50 IS NOT USED AT PRESENT
00114 5* C END
00115 6* IF(CON(4).LT.1.0) CON(4) = 1.0
00117 7* IF(KON(5).LE.0) KON(5) = 1
00121 8* IF(CON(6).LE.0) CON(6) = 1.E+8
00123 9* IF(CON(8).LE.0) CON(8) = 1.E+8
00125 10* IF(CON(9).LE.0) CON(9) = 1.0
00127 11* IF(CON(11).LE.0) CON(11) = 1.E+8
00131 12* IF(CON(18).LE.0) GO TO 999
00133 13* IF(CON(19).LE.0) CON(19) = 1.E+8
00135 14* IF(CON(31).NE.0) GO TO 995
00137 15* PASS = -1.0
00140 16* NNC = NND+NNA
00141 17* IE = NTH
00142 18* IEH = NTH+NNC
00143 19* NLA = NOIM
00144 20* NTH = NTH+NNC+NND
00145 21* NOIM = NNDIM+NNC+NND
00145 22* CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00146 23* I = NLA+NNC+NND
00147 24* IF(I.LT.0) GO TO 998
00151 25* L1 = NND+1
00152 26* TSTEP = CON(18)
00153 27* TSTEP0 = 0.0
00154 28* TPRINT = CON(13)
00154 29* INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00155 30* TSUM = 0.0
00155 31* 5 DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00156 32* IF(CON(13)+CON(18).LE.CON(3)) GO TO 10
00156 33* DONT EXCEED IT
00160 34* CON(18) = CON(13)-CON(13)
00160 35* IS THE TIME STEP LARGER THAN ALLOWED
00161 36* 10 IF(TSTEP.LE.CON(8)) GO TO 15
00163 37* TSTEP = CON(8)
00163 38* DOES THE TIME STEP PLUS THE TIME STEP EXCEED OUTPUT INTERVAL

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00164 39* 15 IF(TSUM+TSTEP-CON(18)) 25,30,20
00164 40* C DONT EXCEED IT
00167 41* C 20 TSTEP = CON(18) - TSUM
00170 42* C GO TO 30
00170 43* C DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
00171 44* C 25 IF(TSUM+2.0*TSTEP.LE.CON(18)) GO TO 30
00171 45* C APPROACH THE OUTPUT INTERVAL GRADUALLY
00173 46* C TSTEP = (CON(18)-TSUM)/2.0
00173 47* C STORE DELTA TIME STEP IN THE CONSTANTS
00174 48* C 30 CON(2) = TSTEP
00174 49* C IS THE TIME STEP USED LESS THAN THE TIME STEP ALLOWED
00175 50* C IF(TSTEP.LT.CON(21)) GO TO 997
00175 51* C CALCULATE THE NEW TIME
00177 52* C CON(1) = TPRINT+TSUM+TSTEP
00177 53* C COMPUTE THE MEAN TIME BETWEEN ITERATIONS
00200 54* C CON(14) = (CON(1)+CON(13))/2.0
00200 55* C ZERO OUT ALL SOURCE LOCATIONS AND EXTRA LOCATIONS
00201 56* C DO 35 I = 1,MND
00204 57* C LE = IE+1
00205 58* C X(LE) = 0.0
00205 59* C Q(I) = 0.0
00207 60* C 35 CONTINUE
00207 61* C SHIFT THE ARITHMETIC TEMPERATURES INTO THE EXTRA LOCATIONS
00211 62* C IF(INNA.LE.0) GO TO 45
00213 63* C DO 40 I = 1,MNC
00216 64* C Q(I) = 0.0
00217 65* C LE = IE+1
00220 66* C X(LE) = T(I)
00221 67* C 40 CONTINUE
00223 68* C 45 KON(12) = 0
00224 69* C CALL VARBL1
00225 70* C IF(KON(12).NE.0) GO TO 10
00227 71* C J1 = 0
00230 72* C J2 = 1
00231 73* C TCGM = 0.0
00232 74* C CKM = 1.E+10
00232 75* C CALCULATE Q SUM AND G SUM
00233 76* C DO 85 I = 1,MND
00236 77* C LE = IE+1
00237 78* C INCLUDE VARC+LIST
00240 78* C IF(FLD(1,1,NS01(J1+1)).EQ.0) GO TO 2000
00242 78* C NTYPE = FLD(0,5,NS02(J2))
00243 78* C LA = FLD(5,17,NS02(J2))
00244 78* C LK = FLD(22,14,NS02(J2))
00245 78* C GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
00246 78* C 1005 CALL DID1WM(T(I),A(LA),XK(LK),C(I))
00247 78* C GO TO 1999
00250 78* C 1010 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00251 78* C 1012 J2 = J2+1
00252 78* C LA = FLD(5,17,NS02(J2))
00253 78* C LK = FLD(22,14,NS02(J2))
00254 78* C CALL DID1WM(T(I),A(LA),XK(LK),C2)
00255 78* C GO TO 1998
00256 78* C 1015 C1 = XJ(LK)*XK(LA)
00257 78* C GO TO 1012
00260 78* C 1020 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00261 78* C J2 = J2+1
00262 78* C LA = FLD(5,17,NS02(J2))
00263 78* C LK = FLD(22,14,NS02(J2))
00264 78* C C2 = XK(LK)*XK(LA)

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00265 78* 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
00266 78* 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00267 78* 1032 J2 = J2+1
00270 78* LA = FLD(5,17,NSQ2(J2))
00271 78* LK = FLD(22,14,NSQ2(J2))
00272 78* CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00273 78* 1035 C1 = XK(LK)*XK(LA)
00274 78* 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00275 78* J2 = J2+1
00276 78* LA = FLD(5,17,NSQ2(J2))
00277 78* LK = FLD(22,14,NSQ2(J2))
00300 78* C2 = XK(LK)*XK(LA)
00301 78* 1045 CALL D2DIWM(T(I),CON(14),A(LA),XK(LK),C(I))
00302 78* 1998 C(I) = C1+C2
00303 78* 1999 J2 = J2+1
00304 78* 2000 CONTINUE
00305 78* END
00306 78* INCLUDE VARO,LIST
00307 78* IF(FLD(4,1,NSQ1(J1+1),E0,0) GO TO 5000
00310 78* NTYPE = FLD(0,5,NSQ2(J2))
00311 78* LA = FLD(5,17,NSQ2(J2))
00312 78* LK = FLD(22,14,NSQ2(J2))
00313 78* 4005 GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4050), NTYPE
00314 78* Q(I) = XK(LK)+Q(I)
00315 78* 4010 Q1 = 0.0
00316 78* 4012 CALL D1DIWM(T(I),A(LA),XK(LK),Q2)
00317 78* 4015 Q1 = 0.0
00318 78* 4017 CALL D1DIWM(CON(14),A(LA),XK(LK),Q2)
00319 78* 4020 CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)
00320 78* 4022 J2 = J2+1
00321 78* LA = FLD(5,17,NSQ2(J2))
00322 78* LK = FLD(22,14,NSQ2(J2))
00323 78* 4025 Q1 = XK(LK)*XK(LA)
00324 78* 4030 CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)
00325 78* J2 = J2+1
00326 78* LA = FLD(5,17,NSQ2(J2))
00327 78* LK = FLD(22,14,NSQ2(J2))
00328 78* 4035 CALL D1DIWM(CON(14),A(LA),XK(LK),Q1)
00329 78* 4037 J2 = J2+1
00330 78* LA = FLD(5,17,NSQ2(J2))
00331 78* LK = FLD(22,14,NSQ2(J2))
00332 78* 4040 Q1 = XK(LK)*XK(LA)
00333 78* 4998 Q(I) = Q1+Q2+G(I)
00334 78* 4999 J2 = J2+1
00335 78* 5000 CONTINUE
00336 78*
00337 78*
00338 78*
00339 78*
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CNDUFR

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00362 79* 50 J1 = J1+1
00363 80* LG = FLD(5,16,NSQ1(J1))
00364 81* IF(LG.EQ.0) GO TO 65
00365 82* LTA = FLD(22,14,NSQ1(J1))
00366 83* INCLUDE VARG,LIST
00370 84* CHECK FOR RADIATION CONDUCTOR
00370 85* IF(FLD(2,1,NSQ1(J1)).EQ.0) GO TO 3000
00371 85* NTYPE = FLD(0,5,NSQ2(J2))
00373 85* LA = FLD(5,17,NSQ2(J2))
00374 85* LK = FLD(22,14,NSQ2(J2))
00375 85* GO TO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00376 85* 2060,2065), NTYPE
00376 85* 2005 TM = (T(I)+T(LTA))/2.0
00377 85* 2007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
00400 85* GO TO 2999
00401 85* 2010 TM = T(I)
00402 85* GO TO 2007
00403 85* 2015 CALL DID1WM(T(I),A(LA),XK(LK),G1)
00404 85* 2017 J2 = J2+1
00405 85* LA = FLD(5,17,NSQ2(J2))
00406 85* LK = FLD(22,14,NSQ2(J2))
00407 85* CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
00410 85* GO TO 2998
00411 85* 2020 G1 = XK(LK)*XK(LA)
00412 85* GO TO 2017
00413 85* 2025 CALL DID1WM(T(I),A(LA),XK(LK),G1)
00414 85* J2 = J2+1
00415 85* LA = FLD(5,17,NSQ2(J2))
00416 85* LK = FLD(22,14,NSQ2(J2))
00417 85* G2 = XK(LK)*XK(LA)
00420 85* GO TO 2998
00421 85* 2030 TM = (T(I)+T(LTA))/2.0
00422 85* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00423 85* GO TO 2999
00424 85* 2035 TM = T(I)
00425 85* GO TO 2032
00426 85* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00427 85* 2042 J2 = J2+1
00430 85* LA = FLD(5,17,NSQ2(J2))
00431 85* LK = FLD(22,14,NSQ2(J2))
00432 85* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00433 85* GO TO 2998
00434 85* 2045 G1 = XK(LK)*XK(LA)
00435 85* GO TO 2042
00436 85* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00437 85* J2 = J2+1
00440 85* LA = FLD(5,17,NSQ2(J2))
00441 85* LK = FLD(22,14,NSQ2(J2))
00442 85* G2 = XK(LK)*XK(LA)
00443 85* GO TO 2998
00444 85* 2055 TM = (T(I)+T(LTA))/2.0
00445 85* CALL D2D1WM(TM,COU(14),A(LA),XK(LK),G(LG))
00446 85* GO TO 2999
00447 85* 2060 TM = T(LTA)
00450 85* GO TO 2007
00451 85* 2065 TM = T(LTA)
00452 85* GO TO 2032
00453 85* G(LG) = 1./(1./G1+1./G2)
00454 85* 2998 IF(FLD(3,1,NSQ1(J1)).EQ.1) G(LG) = G1*G2
00455 85*

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CMDUFR

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00457 85* 2999 J2 = J2+1
00460 85* 3000 CONTINUE
00461 85* END
00462 86* IF(FLD(I,1,NSQ(I,J1)),EQ,0) GO TO 55
00464 87* T1 = T(I)+460.0
00465 88* T2 = T(LTA)+460.0
00466 89* GV = G(LG)*(T1+T2)*(T1+T2)
00467 90* GO TO 60
00470 91* 55 GV = G(LG)
00471 92* C OBTAIN THE G PLUS SUMMATION G*TA TERM
00472 93* Q(I) = Q(I)+GV*(LTA)
00473 94* C SAVE SUMMATION OF CONDUCTORS
00474 95* X(LE) = X(LE)+GV
00475 96* C CHECK FOR ADJOINING DIFFUSION NODE OR ONE-WAY CONDUCTOR
00476 97* IF(LTA.GT.MND.OR.FLD(I,1,NSQ(I,J1)),EQ,1) GO TO 65
00477 98* LEA = IE+LTA
00478 99* X(LEA) = X(LEA)+GV
00479 100* Q(LTA) = Q(LTA)+GV*(LTA)
00480 101* C CHECK FOR LAST CONDUCTOR
00481 102* 65 IF(NSQ(I,J1).GT,0) GO TO 50
00482 103* 85 CONTINUE
00483 104* DT1 = TSTEP/(TSTEP+TSTEP0)
00484 105* DT2 = TSTEP/(TSTEP+TSTEP0)
00485 106* C OBTAIN NEW DIFFUSION TEMPERATURES, DTMPC AND CSGMIN
00486 107* DO 100 I = 1,NND
00487 108* LE = IE+I
00488 109* LEH = IEH+I
00489 110* C CALCULATE C/SK MINIMUM
00490 111* T1 = C(I)/X(LE)
00491 112* IF(T1.GE,CKM) GO TO 90
00492 113* CKM = T1
00493 114* KON(35) = I
00494 115* C COMPUTE NEW TEMPERATURES USING CALCULATED SOURCE TERMS
00495 116* T1=(DT1*X(LEH)*(C(I)/TSTEP-X(LE))+Q(I))/(C(I)*(1.-DT2)/TSTEP+DT2*
00496 117* X(LE))
00497 118* C CALCULATE THE ABSOLUTE VALUE TEMPERATURE CHANGE
00498 119* T2 = ABS(T1-T1)
00499 120* C SAVE THE LARGEST TEMPERATURE CHANGE
00500 121* IF(TCGM.GE,T2) GO TO 95
00501 122* TCGM = T2
00502 123* KON(36) = I
00503 124* C STORE THE TEMPERATURES
00504 125* X(LE) = T(I)
00505 126* T(I) = T1
00506 127* 100 CONTINUE
00507 128* CON(17) = CKM
00508 129* DELTA = CKM*CON(4)
00509 130* IF(CKM.LE,0.0) GO TO 996
00510 131* C CHECK FOR FIRST PASS
00511 132* IF(PASS.GT,0.0) GO TO 115
00512 133* C UNDO THE TEMPERATURE CALCULATIONS
00513 134* DO 110 I = 1,NNC
00514 135* LE = IE+I
00515 136* T(I) = X(LE)
00516 137* 110 CONTINUE
00517 138* IF(PASS.GT,0.0) GO TO 15
00518 139* PASS = 1.0
00519 140* CON(1) = TPRINT
00520 141* CON(2) = 0.0
00521 142*

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CNDUFR

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00554 143* TSTEP = CKM*0.9
00555 144* DO 112 I = 1,NND
00560 145* LEH = IEH+I
00561 146* 112 X(LEH) = T(I)
00563 147* GO TO 195
00563 148* C IS THE TIME STEP USED LESS THAN THE TIME STEP CALCULATED
00564 149* C 115 IF(TSTEP.LE.DELTA) GO TO 130
00564 150* C COMPUTE THE TIME STEP
00566 151* TSTEP = DELTA*0.95
00567 152* GO TO 105
00570 153* 120 TSTEP = 0.95*TSTEP*CON(6)/TCGM
00571 154* GO TO 105
00572 155* 125 TSTEP = 0.95*TSTEP*CON(11)/TCGM
00573 156* GO TO 105
00573 157* C SEE IF THE TEMPERATURE CHANGE WAS TOO LARGE
00574 158* C 130 IF(TCGM.GT.CON(6)) GO TO 120
00574 159* C STORE THE MAXIMUM DIFFUSION TEMPERATURE CHANGE
00576 160* CON(15) = TCGM
00576 161* C CHECK TO SEE IF THERE ARE ANY ARITHMETIC NODES
00577 162* C IF(MNA.LE.0) GO TO 185
00577 163* C COMPUTE ARITHMETIC TEMPERATURES BY SUCCESSIVE POINT OVER-RELAX
00601 164* DN = CON(9)
00602 165* DD = 1.0-DN
00603 166* LAX = KON(5)
00604 167* DO 170 I = 1,LAX
00607 168* JJ1 = J1
00610 169* JJ2 = J2
00611 170* TCGM = 0.0
00612 171* KON(20) = I
00613 172* DO 165 L = LI,NNC
00616 173* SUMC = 0.0
00617 174* SUMCV = 0.0
00620 175* IF(I.GT.1) GO TO 6000
00622 176* INCLUDE VRG2,LIST
00623 176* IF(FLD(4,1,NSQ1(JJ1+1)).EQ.0) GO TO 6000
00625 176* NTYPE = FLD(0,5,NSQ2(JJ2))
00626 176* LA = FLD(5,17,NSQ2(JJ2))
00627 176* LK = FLD(22,14,NSQ2(JJ2))
00630 176* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
00631 176* 5005 Q(L) = XK(LK)+Q(L)
00632 176* GO TO 5999
00633 176* 5010 Q1 = 0.0
00634 176* 5012 CALL DID1WM(T(L),A(LA),XK(LK),Q2)
00635 176* GO TO 5998
00636 176* 5015 Q1 = 0.0
00637 176* 5017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
00640 176* GO TO 5998
00641 176* 5020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00642 176* 5022 JJ2 = JJ2+1
00643 176* LA = FLD(5,17,NSQ2(JJ2))
00644 176* LK = FLD(22,14,NSQ2(JJ2))
00645 176* GO TO 5017
00646 176* 5025 Q1 = XK(LK)*XK(LA)
00647 176* GO TO 5022
00650 176* 5030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00651 176* JJ2 = JJ2+1
00652 176* LA = FLD(5,17,NSQ2(JJ2))
00653 176* LK = FLD(22,14,NSQ2(JJ2))
00654 176* Q2 = XK(LK)*XK(LA)
00655 176* GO TO 5998

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00656 176*	5035 CALL DID1WM(CON(14),A(LA),XK(LK),61)	*NEW
00657 176*	5037 JJ2 = JJ2+1	*NEW
00660 176*	LA = FLD(15,17,NSQ2(JJ2))	**--2
00661 176*	LK = FLD(22,14,NSQ2(JJ2))	
00662 176*	GO TO 5012	
00663 176*	GO TO 5037	
00664 176*	GO TO 5012	
00665 176*	GO TO 5037	
00666 176*	G(L) = G1+02*0(L)	
00667 176*	5999 JJ2 = JJ2+1	
00670 176*	6000 CONTINUE	
00671 177*	END	
00672 178*	135 JJI = JJI+1	
00673 179*	LG = FLD(15,16,NSQ1(JJI))	
00674 180*	LTA = FLD(22,14,NSQ1(JJI))	
00676 181*	IF(I.GT.1) GO TO 4000	
00677 182*	INCLUDE VRG2,LIST	
00701 182*	C	
00702 182*	CHECK FOR RADIATION CONDUCTOR	
00703 182*	IF(FLD(2,1,NSQ1(JJI)).EQ.0) GO TO 4000	
00704 182*	NTYPE = FLD(0,5,NSQ2(JJ2))	
00705 182*	LA = FLD(15,17,NSQ2(JJ2))	
00706 182*	LK = FLD(22,14,NSQ2(JJ2))	
00707 182*	GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,	
00710 182*	3060,3065), NTYPE	
00711 182*	3005 TM = (TL)+T(LTA))/2.0	
00712 182*	3007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))	
00713 182*	GO TO 3999	
00714 182*	3010 TM = T(L)	
00715 182*	GO TO 3007	
00716 182*	3015 CALL DID1WM(T(L),A(LA),XK(LK),61)	
00717 182*	3017 JJ2 = JJ2+1	
00720 182*	LA = FLD(15,17,NSQ2(JJ2))	
00721 182*	LK = FLD(22,14,NSQ2(JJ2))	
00722 182*	CALL DID1WM(T(LTA),A(LA),XK(LK),62)	
00723 182*	GO TO 3998	
00724 182*	3020 G1 = XK(LK)*XK(LA)	
00725 182*	GO TO 3017	
00726 182*	3025 CALL DID1WM(T(L),A(LA),XK(LK),61)	
00727 182*	JJ2 = JJ2+1	
00730 182*	LA = FLD(15,17,NSQ2(JJ2))	
00731 182*	LK = FLD(22,14,NSQ2(JJ2))	
00732 182*	G2 = XK(LK)*XK(LA)	
00733 182*	GO TO 3998	
00734 182*	3030 TM = (TL)+T(LTA))/2.0	
00735 182*	3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))	
00736 182*	GO TO 3999	
00737 182*	3035 TM = T(L)	
00740 182*	GO TO 3032	
00741 182*	3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
00742 182*	3042 JJ2 = JJ2+1	
00743 182*	LA = FLD(15,17,NSQ2(JJ2))	
00744 182*	LK = FLD(22,14,NSQ2(JJ2))	
00745 182*	CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)	
00746 182*	GO TO 3998	
00747 182*	3045 G1 = XK(LK)*XK(LA)	
00748 182*	GO TO 3042	
00749 182*	3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
00750 182*	JJ2 = JJ2+1	
00751 182*	LA = FLD(15,17,NSQ2(JJ2))	
	LK = FLD(22,14,NSQ2(JJ2))	
	G2 = XK(LK)*XK(LA)	

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00752 182* GO TO 3998
00753 182* TM = (T(L)+T(LTA))/2.0
00754 182* CALL D2DIM(TM,CON(14),A(LA),XK(LK),G(LG))
00755 182* GO TO 3999
00756 182* TM = T(LTA)
00757 182* GO TO 3007
00760 182* TM = T(LTA)
00761 182* GO TO 3032
00762 182* G(LG) = 1./(1./G1+1./G2)
00763 182* IF(FLD(3,1,NSQ1(JJ1)),EQ.1) G(LG) = 61*G2
00765 182* J2 = JJ2+1
00766 182* 4000 CONTINUE
00767 182* END
00770 183* IF(FLD(3,1,NSQ1(JJ1)),EQ.0) GO TO 140
00772 184* T1 = T(L)+460.0
00773 185* T2 = T(LTA)+460.0
00774 186* GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
00775 187* GO TO 145
00776 188* 140 GV = G(LG)
00777 189* 145 SUMC = SUMC+GV
01000 190* SUMCV = SUMCV+GV*T(LTA)
01000 191* CHECK FOR LAST CONDUCTOR
01001 192* IF(NSQ1(JJ1).GT.0) GO TO 135
01003 193* T2 = OD+T(L)+DN*(SUMCV+G(L))/SUMC
01003 194* OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
01004 195* T1 = ARS(T(L)-T2)
01004 196* STORE THE NEW TEMPERATURE
01005 197* T(L) = T2
01005 198* SAVE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
01006 199* IF(TCGM.GE.T1) GO TO 165
01010 200* TCGM = T1
01011 201* KON(37) = L
01012 202* 165 CONTINUE
01012 203* SEE IF RELAXATION CRITERIA WAS MET
01014 204* IF(TCGM.LE.CON(19)) GO TO 175
01016 205* 170 CONTINUE
01016 206* C STORE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
01020 207* CON(30) = TCGM
01021 208* C COMPUTE THE ARITHMETIC TEMPERATURE CHANGE
01021 209* TCGM = 0.0
01022 210* DO 180 I = L1,NNC
01025 211* LE = IE+1
01026 212* T1 = ABS(T(I)-X(LE))
01027 213* IF(T1.LT.TCGM) GO TO 180
01031 214* TCGM = T1
01032 215* KON(38) = I
01033 216* 180 CONTINUE
01033 217* C SEE IF ATMPCA WAS SATISFIED
01035 218* IF(TCGM.GT.CON(11)) GO TO 125
01037 219* CON(16) = TCGM
01040 220* KON(12) = 0
01041 221* CALL VARBL2
01041 222* CHECK THE RACKUP SWITCH
01042 223* IF(KON(12).NE.0) GO TO 105
01042 224* ADVANCE TIME
01044 225* CON(13) = CON(1)
01045 226* TSUM = TSUM+TSTEP
01046 227* TSTEP = TSTEP
01047 228* DO 200 I = 1,NNID
01052 229* LE = IE+1
```

CNDUFR

```
01053 230* LEH = IEH*I
01054 231* 200 X(LEH) = X(LE)
01056 232* TSTEP = DELTA*0.95
01056 233* CHECK FOR TIME TO PRINT
01057 234* IF(TSUM.GE.CON(18)) GO TO 190
01057 235* CHECK FOR PRINT EVERY ITERATION
01061 236* IF(KON(7).EQ.0) GO TO 10
01063 237* CALL OUTCAL
01064 238* GO TO 10
01064 239* TRY TO EVEN THE OUTPUT INTERVALS
01065 240* 190 TPRINT = TPRINT+TSUM
01066 241* 195 CALL OUTCAL
01066 242* IS TIME GREATER THAN END COMPUTE TIME
01067 243* IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
01071 244* NTH = IE
01072 245* NDIM = NLA
01073 246* RETURN
01074 247* 995 WRITE(6,885)
01076 248* GO TO 1000
01077 249* 996 WRITE(6,886)
01101 250* GO TO 1000
01102 251* 997 WRITE(6,887)
01104 252* GO TO 1000
01105 253* 998 WRITE(6,888) I
01110 254* GO TO 1000
01111 255* 999 WRITE(6,889)
01113 256* 1000 CALL OUTCAL
01114 257* CALL EXIT
01115 258* 885 FORMAT(46H CNDUFR REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
01116 259* 886 FORMAT(24H CS6MIN ZERO OR NEGATIVE)
01117 260* 887 FORMAT(20H TIME STEP TOO SMALL)
01120 261* 888 FORMAT(18,20H LOCATIONS AVAILABLE)
01121 262* 889 FORMAT(19H NO OUTPUT INTERVAL)
01122 263* END
```

END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
CNDUFR CODE RELOCATABLE

CNQUIK

01 FOR: CNQUIK
 UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-10D CREATED ON 20 AUG 70
 THIS COMPILATION WAS DONE ON 09 JUN 70 AT 23:15:06

16 FEB 71

SUBROUTINE CNQUIK ENTRY POINT 003643

STORAGE USED (BLOCK, NAME, LENGTH)

```

0001 *CODE 003660
0000 *CONST+TEMP 000100
0002 *SIMPLE VAR 00054
0004 *ARRAYS 000000
0005 *BLANK 000000
0006 TITLE 000001
0007 TEMP 000001
0010 CAP 000001
0011 SOURCE 000001
0012 COND 000001
0013 PC1 000001
0014 PC2 000001
0015 KONST 000001
0016 ARRAY 000001
0017 FIXCON 000001
0020 XSPACE 000003
0021 DIMENS 000010
  
```

EXTERNAL REFERENCES (BLOCK, NAME)

```

0022 VARBL1
0023 DID1WM
0024 PLYAWM
0025 D2D1WM
0026 VARBL2
0027 OUTCAL
0030 EXIT
0031 NERR2$
0032 EXP
0033 NWDUS$
0034 NIO2$
0035 NER10$
  
```

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

```

0016 R 000000 A 0010 R 000000 C 0002 R 000017 CKM
0002 R 000024 C2 0002 R 000045 DD 0002 R 000043 DELTA
0002 R 000041 DT2 0012 R 000000 G 0002 R 000036 GV
0006 R 000000 H 0002 I 000005 I 0002 I 000002 IE
0002 I 000050 JJ2 0002 I 000014 J1 0002 I 000015 J2
0002 I 000051 L 0002 I 000021 LA 0002 I 000046 LAX
0002 I 000042 LEH 0002 I 000027 LG 0002 I 000022 LK
0002 I 000030 LTA 0002 I 000006 L1 0002 I 000005 NAT
0021 I 000003 NGT 0002 I 000004 NLA 0021 I 000001 NNA
0021 I 000002 INT 0013 I 000000 NSG1 0014 I 000000 NS32
0020 000002 IAX 0002 R 000000 PASS 0011 R 000000 Q
0002 R 000052 SUMC 0002 R 000053 SUMCV 0007 R 000000 T
0017 R 000000 COM 0002 R 000044 DN 0002 R 000023 C1
0002 R 000032 G1 0002 R 000033 G2 0002 R 000040 DT1
0002 I 000003 IEH 0002 I 000007 JJI 0002 I 000047 JJJ
0015 I 000000 K 0017 I 000000 KON 0017 I 000000 K0N
0002 I 000013 LE 0002 I 000037 LEA 0002 I 000007 LEA
0021 I 000006 LSG1 0021 I 000007 LSG2 0021 I 000007 LSG2
0021 I 000004 NCT 0020 I 000000 NDM 0020 I 000000 NDM
0002 I 000001 NNC 0021 I 000000 NND 0021 I 000000 NND
0020 I 000001 NTH 0002 I 000020 NTYPE 0002 I 000020 NTYPE
0002 R 000025 Q1 0002 R 000026 Q2 0002 R 000026 Q2
0002 R 000016 TCGM 0002 R 000031 TM
  
```

CNGUIK

```

0002 R 000011 TPRINT 0002 R 000007 TSTEP 0002 R 000010 TSTEPO 0002 R 000012 TSUM 0002 R 000034 T1
0002 R 000035 T2 0002 R 000002 X 0002 R 000000 XK 0001 000132 10L 0001 003617 1000L
0001 003444 1005L 0001 000365 1010L 0001 000403 1012L 0001 000436 1015L 0001 000444 1020L
0001 003432 1025G 0001 000504 1025L 0001 000530 1030L 0001 000551 1032L 0001 000607 1035L
0001 000615 1040L 0001 000660 1045L 0001 002150 105L 0001 002150 105L 0001 003505 1052G 0001 002217 115L
0001 002230 120L 0001 002236 125L 0001 002244 130L 0001 002244 130L 0001 002574 135L 0001 003342 140L
0001 003345 145L 0001 000141 15L 0001 003412 165L 0001 003423 175L 0001 003454 180L
0001 003465 185L 0001 003537 190L 0001 003542 195L 0001 000705 1999L 0001 000705 1999L
0001 000710 2000L 0001 001250 2005L 0001 001255 2007L 0001 001275 2010L 0001 001300 2015L
0001 001316 2017L 0001 000213 202G 0001 001352 2020L 0001 001360 2025L 0001 001420 2030L
0001 001425 2032L 0001 001450 2035L 0001 001453 2040L 0001 001474 2042L 0001 001533 2045L
0001 001541 2050L 0001 001604 2055L 0001 001632 2060L 0001 001636 2065L 0001 001533 2045L
0001 000273 234G 0001 000153 25L 0001 001642 2998L 0001 001670 2999L 0001 001670 2999L 0001 000232 214G
0001 001673 3000L 0001 002662 3005L 0001 002667 3007L 0001 002707 3010L 0001 002712 3015L
0001 002730 3017L 0001 002764 3020L 0001 002772 3025L 0001 002707 3030L 0001 003037 3032L
0001 003062 3035L 0001 003065 3040L 0001 003106 3042L 0001 003145 3045L 0001 003153 3050L
0001 003216 3055L 0001 003244 3060L 0001 003250 3065L 0001 003254 3998L 0001 003302 3999L
0001 003305 4000L 0001 000753 4005L 0001 000760 4010L 0001 000761 4012L 0001 001015 4030L
0001 001001 4017L 0001 001016 4020L 0001 001032 4022L 0001 001047 4025L 0001 001047 4025L
0001 001113 4035L 0001 001127 4037L 0001 001144 4040L 0001 000244 45L 0001 001152 4998L
0001 001156 4999L 0001 000120 5L 0001 001161 50L 0001 001161 5000L 0001 002366 5005L
0001 002373 5100L 0001 002374 5012L 0001 002413 5015L 0001 002414 5017L 0001 002431 5020L
0001 002445 5022L 0001 002462 5025L 0001 002470 5030L 0001 002526 5035L 0001 002542 5037L
0001 002557 5040L 0001 002020 507G 0001 002154 543G 0001 001724 55L 0001 002205 560G
0001 002565 5998L 0001 002571 5999L 0001 001727 60L 0001 002574 6000L 0001 002273 607G
0001 002314 616G 0001 001775 65L 0001 000000 885F 0001 000011 886F 0001 000016 897F
0000 000023 888F 0001 002044 90L 0001 002044 90L 0001 002124 95L 0001 003561 995L
0001 003567 996L 0001 003575 997L 0001 003603 998L 0001 003612 999L

```

1* SUBROUTINE CNGUIK
2* COMBINED EXPONENTIAL PREDICTION AND DUFORT-FRANKEL SINDA ROUTINE
3* THE SHORT PSEUDO-COMPUTE SEQUENCE IS REQUIRED
4* INCLUDE COMMON LIST
5* COMMON /FC1/NS01(1) /TEMP/T(1) /CAP/C(1) /SOURCE/Q(1) /COND/G(1)
6* COMMON /FIXCON/KON(1) /XSPACE/NDIM,NTH,X(1)
7* COMMON /DIMENS/ IJND,NNA,NNT,NGT,NCT,NAT,LS01,LS02
8* DIMENSION CON(1),XK(1),NX(1)
9* EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
10* END
11* INCLUDE DEF*LIST
12* C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
13* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
14* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
15* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
16* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR, EXPLICIT (INLOOP)
17* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (DTMPCA)
18* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DPEITR)
19* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (DTIMEH)
20* C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DAMPA)
21* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPO)
22* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPO)
23* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
24* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
25* C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
26* C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
27* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)

CNOUIK

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00113 5* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
00113 5* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CS6MIN)
00113 5* C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 5* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 5* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 5* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 5* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
00113 5* C CC23 CONTAINS THE C/SG MAXIMUM (CS6MAX)
00113 5* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CS6RCL)
00113 5* C CC25 CONTAINS THE C/SG RANGE CALCULATED (CS6RCL)
00113 5* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 5* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DPLXCC)
00113 5* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 5* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 5* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 5* C CC31 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 5* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGBAL)
00113 5* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 5* C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 5* C CC35 CONTAINS RELATIVE NODE NUMBER OF CS6MIN
00113 5* C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 5* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 5* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMFCC
00113 5* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 5* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 5* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 5* C CC50 IS NOT USED AT PRESENT
00114 5* C END
00115 6* IF(CON(4).LT.1.0) CON(4) = 1.0
00117 7* IF(KON(5).LE.0) KON(5) = 1
00121 8* IF(CON(6).LE.0) CON(6) = 1.E+8
00123 9* IF(CON(8).LE.0) CON(8) = 1.E+8
00125 10* IF(CON(9).LE.0) CON(9) = 1.0
00127 11* IF(CON(11).LE.0) CON(11) = 1.E+8
00131 12* IF(CON(18).LE.0) GO TO 999
00133 13* IF(CON(19).LE.0) CON(19) = 1.E+8
00135 14* IF(KON(31).NE.0) GO TO 995
00137 15* PASS = -1.0
00140 16* NNC = NND+RNA
00141 17* IE = NTH
00142 18* IEH = NTH+NNC
00143 19* NLA = NDIM
00144 20* NTH = NTH+NNC+IND
00145 21* NDIM = NDIM-NNC-NND
00145 22* CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00146 23* I = NLA-NNC-NND
00147 24* IF(I.LT.0) GO TO 998
00151 25* LI = NND+1
00152 26* TSTEP = CON(18)
00153 27* TSTEPO = 0.0
00154 28* TPRINT = CON(13)
00154 29* INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00155 30* TSUM = 0.0
00155 31* DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00156 32* IF(CON(13)+CON(18).LE.CON(13)) GO TO 10
00156 33* DONT EXCEED IT
00160 34* CON(18) = CON(3)-CON(13)
00160 35* IS THE TIME STEP LARGER THAN ALLOWED
00161 36* IF(TSTEP.LE.CON(16)) GO TO 15
00163 37* TSTEP = CON(6)

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00-2

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00264 78* C2 = XK(LK)*XK(LA)
00265 78* GO TO 1998
00266 78* 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
00267 78* GO TO 1999
00270 78* 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00271 78* 1032 J2 = J2+1
00272 78* LA = FLD(5,17,NSQ2(J2))
00273 78* LK = FLD(22,14,NSQ2(J2))
00274 78* CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00275 78* GO TO 1998
00276 78* 1035 C1 = XK(LK)*XK(LA)
00277 78* GO TO 1032
00300 78* 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00301 78* J2 = J2+1
00302 78* LA = FLD(5,17,NSQ2(J2))
00303 78* LK = FLD(22,14,NSQ2(J2))
00304 78* C2 = XK(LK)*XK(LA)
00305 78* GO TO 1998
00306 78* 1045 CALL D2DIWM(T(I),CON(14),A(LA),XK(LK),C(I))
00307 78* GO TO 1999
00310 78* 1998 C(I) = C1+C2
00311 78* 1999 J2 = J2+1
00312 78* 2000 CONTINUE
00313 78* END
00314 79* INCLUDE VARQ,LIST
00315 79* IF(FLD(4,1,NSQ1(J1+1),E0,0) GO TO 5000
00317 79* NTYPE = FLD(0,5,NSQ2(J2))
00320 79* LA = FLD(5,17,NSQ2(J2))
00321 79* LK = FLD(22,14,NSQ2(J2))
00322 79* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE
00323 79* 4005 G(I) = XK(LK)+G(I)
00324 79* GO TO 4999
00325 79* 4010 G1 = 0.0
00326 79* 4012 CALL DIDIWM(T(I),A(LA),XK(LK),G2)
00327 79* GO TO 4998
00330 79* 4015 G1 = 0.0
00331 79* 4017 CALL DIDIWM(CON(14),A(LA),XK(LK),G2)
00332 79* GO TO 4998
00333 79* 4020 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00334 79* 4022 J2 = J2+1
00335 79* LA = FLD(5,17,NSQ2(J2))
00336 79* LK = FLD(22,14,NSQ2(J2))
00337 79* GO TO 4017
00340 79* 4025 G1 = XK(LK)*XK(LA)
00341 79* GO TO 4022
00342 79* 4030 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00343 79* J2 = J2+1
00344 79* LA = FLD(5,17,NSQ2(J2))
00345 79* LK = FLD(22,14,NSQ2(J2))
00346 79* G2 = XK(LK)*XK(LA)
00347 79* GO TO 4998
00350 79* 4035 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00351 79* 4037 J2 = J2+1
00352 79* LA = FLD(5,17,NSQ2(J2))
00353 79* LK = FLD(22,14,NSQ2(J2))
00354 79* GO TO 4012
00355 79* 4040 G1 = XK(LK)*XK(LA)
00356 79* GO TO 4037
00357 79* 4998 G(I) = G1+G2+G(I)
00360 79* 4999 J2 = J2+1

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00361 79* 5000 CONTINUE
00362 79* END
00363 80* 50 J1 = J1+1
00364 81* LG = FLD(5,16,NSQ1(J1))
00365 82* IF(LG.EQ.0) GO TO 65
00366 83* LTA = FLD(22,14,NSQ1(J1))
00367 84* INCLUDE VARGLIST
00370 85* CHECK FOR RADIATION CONDUCTOR
00371 85* IF(FLD(2,1,NSQ1(J1)).EQ.0) GO TO 3000
00372 85* NTYPE = FLD(0,5,NSQ2(J2))
00374 85* LA = FLD(5,17,NSQ2(J2))
00375 85* LK = FLD(22,14,NSQ2(J2))
00376 85* GOTO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00375 85* 2060,2065), NTYPE
00377 85* TM = (T(I)+T(LTA))/2.0
00400 85* 2005 CALL DIDIWM(TM,A(LA),XK(LK),6(LG))
00401 85* GO TO 2999
00402 85* 2010 TM = T(I)
00403 85* GO TO 2007
00404 85* 2015 CALL DIDIWM(T(I),A(LA),XK(LK),61)
00405 85* 2017 J2 = J2+1
00406 85* LA = FLD(5,17,NSQ2(J2))
00407 85* LK = FLD(22,14,NSQ2(J2))
00410 85* CALL DIDIWM(T(LTA),A(LA),XK(LK),62)
00411 85* GO TO 2998
00412 85* 2020 G1 = XK(LK)*XK(LA)
00413 85* GO TO 2017
00414 85* 2025 CALL DIDIWM(T(I),A(LA),XK(LK),61)
00415 85* J2 = J2+1
00416 85* LA = FLD(5,17,NSQ2(J2))
00417 85* LK = FLD(22,14,NSQ2(J2))
00420 85* G2 = XK(LK)*XK(LA)
00421 85* GO TO 2998
00422 85* 2030 TM = (T(I)+T(LTA))/2.0
00423 85* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))
00424 85* GO TO 2999
00425 85* 2035 TM = T(I)
00426 85* GO TO 2032
00427 85* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),61)
00430 85* 2042 J2 = J2+1
00431 85* LA = FLD(5,17,NSQ2(J2))
00432 85* LK = FLD(22,14,NSQ2(J2))
00433 85* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)
00434 85* GO TO 2998
00435 85* 2045 G1 = XK(LK)*XK(LA)
00436 85* GO TO 2042
00437 85* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),61)
00440 85* J2 = J2+1
00441 85* LA = FLD(5,17,NSQ2(J2))
00442 85* LK = FLD(22,14,NSQ2(J2))
00443 85* G2 = XK(LK)*XK(LA)
00444 85* GO TO 2998
00445 85* 2055 TM = (T(I)+T(LTA))/2.0
00446 85* CALL D2DIWM(TM,CON(14),A(LA),XK(LK),6(LG))
00447 85* GO TO 2999
00450 85* 2060 TM = T(LTA)
00451 85* GO TO 2007
00452 85* 2065 TM = T(LTA)
00453 85* GO TO 2032
00454 85* 2996 6(LG) = 1./(1./61+1./62)

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*NEW
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00455 IF(FLD(3,1,NSO1(J1)).EQ.1) G(1,6) = 61*G2
00456 J2 = J2+1
00460 3000 CONTINUE
00461 END
00462 IF(FLD(3,1,NSO1(J1)).EQ.0) GO TO 55
00464 T1 = T(I)+460.0
00465 T2 = T(LTA)+460.0
00466 GV = G(LG)*(T1+T1+T2)*T1+T2
00467 GO TO 60
00470 55 GV = G(LG)
00470 C OBTAIN THE Q PLUS SUMMATION G*TA TERM
00471 Q(I) = Q(I)+GV*T(LTA)
00471 C SAVE SUMMATION OF CONDUCTORS
00472 X(LE) = X(LE)+GV
00472 C CHECK FOR ADJOINING DIFFUSION NODE OR ONE-WAY CONDUCTOR
00473 IF(LTA.GT.HND.OR.FLD(2,1,NSO1(J1)).EQ.1) GO TO 65
00475 LEA = IE+LTA
00476 X(LEA) = X(LEA)+GV
00477 Q(LTA) = Q(LTA)+GV*T(I)
00477 C CHECK FOR LAST CONDUCTOR
00500 65 IF(NSO1(J1).GT.0) GO TO 50
00502 65 CONTINUE
00504 DT1 = TSTEP/(TSTEP+TSTEP0)
00505 DT2 = TSTEP0/(TSTEP+TSTEP0)
00506 DO 100 I = 1,NND
00511 LE = IE+I
00512 LEH = IEH+I
00512 C CALCULATE C/SK MINIMUM
00513 T1 = C(I)/X(LE)
00514 IF(T1.GE.CKM) GO TO 90
00516 CKM = T1
00517 KON(35) = I
00517 C COMPUTE NEW TEMPERATURES USING CALCULATED SOURCE TERMS
00520 T1=(DT1+X(LEH)*(C(I)/(TSTEP-X(LE))+0(I))/(C(I)*(1.-DT2)/TSTEP+DT2*
00521 $X(LE))
00522 T2 = 1.0/EXP(TSTEP*X(LE)/C(I))
00522 T1 = (T1+(1.-T2)*0(I)/X(LE)+T2*T(I))*0.5
00522 C CALCULATE THE ABSOLUTE VALUE TEMPERATURE CHANGE
00523 T2 = ABS(T(I)-T1)
00523 C SAVE THE LARGEST TEMPERATURE CHANGE
00524 IF(TCGM.GE.T2) GO TO 95
00526 TCGM = T2
00527 KON(36) = I
00527 C STORE THE TEMPERATURES
00527 T(I) = T(I)
00530 T(I) = T1
00532 100 CONTINUE
00534 CON(17) = CKM
00534 DELTA = CKM*CON(17)
00535 IF(CKM.LE.0.0) GO TO 996
00536 C CHECK FOR FIRST PASS
00536 IF(PASS.GT.0.0) GO TO 115
00540 C UNDO THE TEMPERATURE CALCULATIONS
00542 DO 110 I = 1,N/C
00545 LE = IE+I
00545 T(I) = X(LE)
00547 110 CONTINUE
00551 IF(PASS.GT.0.0) GO TO 15

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00553 142* PASS = 1.0
00554 143* CON(1) = TPRINT
00555 144* CON(2) = 0.0
00556 145* TSTEP = CKM*0.9
00557 146* DO 112 I = 1,NJD
00562 147* LEH = IEH+I
00563 148* X(LEH) = T(I)
      GO TO 195
C      IS THE TIME STEP USED LESS THAN THE TIME STEP CALCULATED
115 IF(TSTEP.LE.DELTA) GO TO 130
C      COMPUTE THE TIME STEP
120 TSTEP = DELTA*0.95
      GO TO 105
125 TSTEP = 0.95*TSTEP*CON(6)/TCGM
      GO TO 105
130 STORE THE MAXIMUM DIFFUSION TEMPERATURE CHANGE
C      CHECK TO SEE IF THERE ARE ANY ARITHMETIC NODES
      IF(MINALE.0) GO TO 185
C      COMPUTE ARITHMETIC TEMPERATURES BY SUCCESSIVE POINT OVER-RELAX
      CON(15) = TCGM
      DO 170 I = 1,LAX
      J1 = J1
      J2 = J2
      TCGM = 0.0
      KON(20) = I
      DO 165 L = 1,NNC
      SUMCV = 0.0
      IF(I.GT.1) GO TO 6000
      INCLUDE VRG2,LIST
      NTYPE = FLD(0,5,NSG2(JJ2))
      LA = FLD(5,17,NSG2(JJ2))
      LK = FLD(22,14,NSG2(JJ2))
      GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
5005 Q(L) = XK(LK)+0(L)
      GO TO 5999
5010 Q1 = 0.0
5012 CALL DID1WM(T(L),A(LA),XK(LK),Q2)
      GO TO 5998
5015 Q1 = 0.0
5017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
      GO TO 5998
5020 CALL DID1WM(COII(14),A(LA),XK(LK),Q1)
5022 JJ2 = JJ2+1
      LA = FLD(5,17,NSG2(JJ2))
      LK = FLD(22,14,NSG2(JJ2))
      GO TO 5017
5025 Q1 = XK(LK)*XK(LA)
      GO TO 5022
5030 CALL DID1WM(COII(14),A(LA),XK(LK),Q1)
      JJ2 = JJ2+1
      LA = FLD(5,17,NSG2(JJ2))

```

*NEW
*NEW
*--3

*NEW
*NEW
*--2

*NEW

CNQUIK

00655 178* LK = FLD(22,14,NSQ2(JJ2))
 00656 178* Q2 = XK(LK)*XK(LA)
 00657 178* GO TO 3998
 00660 178* 5035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
 00661 178* 5037 JJ2 = JJ2+1
 00662 178* LA = FLD(5,17,NSQ2(JJ2))
 00663 178* LK = FLD(22,14,NSQ2(JJ2))
 00664 178* GO TO 5012
 00665 178* 5040 Q1 = XK(LK)*XK(LA)
 00666 178* GO TO 5037
 00667 178* 5998 Q(L) = 01+02*Q(L)
 00670 178* 5999 JJ2 = JJ2+1
 00671 178* 6000 CONTINUE
 00672 178* END
 00673 179* 135 JJ1 = JJ1+1
 00674 180* LG = FLD(5,16,NSQ1(JJ1))
 00675 181* LIA = FLD(22,14,NSQ1(JJ1))
 00676 182* IF(1.GT.1) GO TO 4000
 00700 183* INCLUDE VRG2,LIST
 00700 184* CHECK FOR RADIATION CONDUCTOR
 00701 184* IF(FLD(2,1,NSQ1(JJ1)).EQ.0) GO TO 4000
 00703 184* NTYPE = FLD(10,5,NSQ2(JJ2))
 00704 184* LA = FLD(5,17,NSQ2(JJ2))
 00705 184* LK = FLD(22,14,NSQ2(JJ2))
 00706 184* GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,
 00707 184* 3060,3065), NTYPE
 00710 184* 3005 TM = (T(L)+T(LTA))/2.0
 00711 184* 3007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
 00712 184* GO TO 3999
 00713 184* 3010 TM = T(L)
 00714 184* GO TO 3007
 00715 184* 3015 CALL DID1WM(T(L),A(LA),XK(LK),G1)
 00716 184* 3017 JJ2 = JJ2+1
 00717 184* LA = FLD(5,17,NSQ2(JJ2))
 00720 184* LK = FLD(22,14,NSQ2(JJ2))
 00721 184* CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
 00722 184* GO TO 3998
 00723 184* 3020 G1 = XK(LK)*XK(LA)
 00724 184* GO TO 3017
 00725 184* 3025 CALL DID1WM(T(L),A(LA),XK(LK),G1)
 00726 184* JJ2 = JJ2+1
 00727 184* LA = FLD(5,17,NSQ2(JJ2))
 00730 184* LK = FLD(22,14,NSQ2(JJ2))
 00731 184* G2 = XK(LK)*XK(LA)
 00732 184* GO TO 3998
 00733 184* 3030 TM = (T(L)+T(LTA))/2.0
 00734 184* 3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
 00735 184* GO TO 3999
 00736 184* 3035 TM = T(L)
 00737 184* GO TO 3032
 00740 184* 3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)
 00741 184* 3042 JJ2 = JJ2+1
 00742 184* LA = FLD(5,17,NSQ2(JJ2))
 00743 184* LK = FLD(22,14,NSQ2(JJ2))
 00744 184* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
 00745 184* GO TO 3998
 00746 184* 3045 G1 = XK(LK)*XK(LA)
 00747 184* GO TO 3042
 00750 184* 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)
 JJ2 = JJ2+1

*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***2

CHRUJK

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00751 184* LA = FLD(5,17,NSG2(JJ2))
00752 184* LK = FLD(22,14,NSG2(JJ2))
00753 184* G2 = XK(LK)*XK(LA)
00754 184* GO TO 3998
00755 184* 3055 TM = (T(L)+T(LTA))/2.0
00756 184* CALL D2DIMM(TM,CON(14),A(LA),XK(LK),G(LG))
00757 184* GO TO 3999
00760 184* 3060 TM = T(LTA)
00761 184* GO TO 3007
00762 184* 3065 TM = T(LTA)
00763 184* GO TO 3032
00764 184* 3998 G(LG) = 1./(1./G1+1./G2)
00765 184* IF(FLD(3,1,NSG1(JJ1)).EQ.1) G(LG) = G1*G2
00767 184* 3999 JJ2 = JJ2+1
00770 184* 4000 CONTINUE
00771 184* END
00772 185* IF(FLD(3,1,NSG1(JJ1)).EQ.0) GO TO 140
00774 186* T1 = T(L)+460.0
00775 187* T2 = T(LTA)+460.0
00776 188* GV = G(LG)*(T1+T2)*T2*(T1+T2)
00777 189* GO TO 145
01000 190* 140 GV = G(LG)
01001 191* 145 SUNC = SUNC+GV
01002 192* SUNCV = SUNCV+GV*T(LTA)
01003 193* CHECK FOR LAST CONDUCTOR
01005 195* IF(NSQ(JJ1).GT.0) GO TO 135
01006 196* T2 = DD*(L)+DN*(SUNCV+Q(L))/SUNC
01007 197* T1 = ABS(T(L)-T2)
01008 198* STORE THE NEW TEMPERATURE
01009 199* T(L) = T2
01010 200* 165 SAVE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
01011 201* IF(TCGM.GE.T1) GO TO 165
01012 202* TCGM = T1
01013 203* KON(17) = L
01014 204* 165 CONTINUE
01016 206* C SEE IF RELAXATION CRITERIA WAS MET
01020 208* IF(TCGM.LE.CON(19)) GO TO 175
01022 209* 170 CONTINUE
01023 210* STORE THE MAXIMUM ARITHMETIC RELAXATION CHANGE
01024 211* TCGM = 0.0
01027 213* DO 180 I = 1,INC
01030 214* LE = LE+1
01031 215* T1 = ABS(T(I)-X(LE))
01033 216* IF(T1.LY.TCGM) GO TO 180
01035 218* TCGM = T1
01037 220* KON(18) = I
01041 221* 180 CONTINUE
01042 222* SEE IF ATMPCA WAS SATISFIED
01043 223* IF(TCGM.GT.CON(11)) GO TO 125
01044 224* CON(16) = TCGM
01046 226* CALL VARBL2
01047 227* 185 KON(12) = 0
01048 228* CHECK THE BACKUP SWITCH
01049 229* IF(KON(12).NE.0) GO TO 105
01050 230* ADVANCE TIME
01051 231* CON(13) = CON(11)
01052 232* TSUM = TSUM+TSTEP
01053 233*
01054 234*
01055 235*
01056 236*
01057 237*
01058 238*

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01050 229*
01051 230*
01054 231*
01055 232*
01056 233*
01060 234*
01060 235*
01061 236*
01061 237*
01063 238*
01065 239*
01066 240*
01066 241*
01067 242*
01070 243*
01070 244*
01071 245*
01073 246*
01074 247*
01075 248*
01076 249*
01100 250*
01101 251*
01103 252*
01104 253*
01106 254*
01107 255*
01112 256*
01113 257*
01115 258*
01116 259*
01117 260*
01120 261*
01121 262*
01122 263*
01123 264*
01124 265*

CNGUIK
TSTEP0 = TSTEP
DO 200 I = 1,NND
LE = IE+1
LEH = IEH+I
200 X(LEH) = X(LE)
TSTEP = DELTA*0.95
C CHECK FOR TIME TO PRINT
IF (TSUM_GE_CON(18)) GO TO 190
C CHECK FOR PRINT EVERY ITERATION
IF (KOR(7).EQ.0) GO TO 10
CALL OUTCAL
GO TO 10
C TRY TO EVEN THE OUTPUT INTERVALS
190 TPRINT = TPRINT+TSUM
195 CALL OUTCAL
C IS TIME GREATER THAN END COMPUTE TIME
IF (CON(1)*1.000001.LT.CON(3)) GO TO 5
NTH = IE
NDIM = NLA
RETURN
995 WRITE(6,885)
GO TO 1000
996 WRITE(6,886)
GO TO 1000
997 WRITE(6,887)
GO TO 1000
998 WRITE(6,888) I
999 WRITE(6,889)
1000 CALL OUTCAL
CALL EXIT
885 FORMAT(46H CNGUIK REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
886 FORMAT(24H CS6:IN ZERO OR NEGATIVE)
887 FORMAT(20H TIME STEP TOO SMALL)
888 FORMAT(18H 20H LOCATIONS AVAILABLE)
889 FORMAT(19H NO OUTPUT INTERVAL)
END

END OF UNIVAC 1103 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(15)
CNGUIK CODE RELOCATABLE

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B. COMPUTER LISTINGS OF SINDA IMPLICIT SOLUTION ROUTINES

	Page
CNBACK	B-2
CNFWBK	B-14
CNVARB	B-26

QIW FOR,* CNBACK,CNBACK
UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-100 CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:11

21 FEB 71

SUBROUTINE CNBACK ENTRY POINT 004236

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	004253
0000	*CONST+TEMP	000125
0002	*SIMPLE VAR	000067
0004	*ARRAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COHD	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARL1
0023	OUTCAL
0024	D1D1WM
0025	PLYAHM
0026	D2D1WM
0027	TOPLIN
0030	VARL2
0031	EXIT
0032	NERK2\$
0033	NWOU\$
0034	NI02\$
0035	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0002	R	000017	AA
0002	R	000034	C1	0002	R	000015	DD
0002	R	000044	GSUM	0002	R	000050	G1
0002	I	000025	I	0002	I	000005	IE1
0002	I	000030	JJ	0002	I	000061	JJ2
0015	I	000000	K	0002	I	000063	KK2
0002	I	000062	L	0002	I	000013	LAX
0002	I	000055	LE2	0002	I	000045	LG
0021	I	000007	LS02	0021	I	000005	NAT
0021	I	000003	NGT	0002	I	000001	FIN
0021	I	000000	INND	0013	I	000000	NS01
0002	I	000031	HTYPE	0002	R	000000	PASS
0020	I	000002	MX				
0010	R	000000	C	0002	R	000014	DN
0002	R	000014	DN	0002	R	000040	G2
0002	I	000006	IE3	0002	I	000021	J1
0002	I	000020	K1	0002	I	000020	K1
0002	I	000026	LE1	0002	I	000066	LE
0021	I	000006	LS01	0002	I	000033	LK
0020	I	000000	NDIM	0021	I	000004	NCT
0002	I	000004	NMC	0021	I	000001	NNA
0020	I	000001	NTH	0014	I	000000	NS02
0002	R	000043	GSUM	0011	R	000000	Q
0017	R	000000	CON				
0012	R	000000	G				
0006	R	000000	H				
0002	I	000007	J				
0002	I	000024	J2				
0002	I	000020	K1				
0002	I	000026	LE1				
0021	I	000006	LS01				
0020	I	000000	NDIM				
0002	I	000004	NMC				
0020	I	000001	NTH				
0002	R	000043	GSUM				

CNBACK,CNBACK

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0002 R 000041 Q1      0002 R 000042 Q2      0002 R 000027 RC      0002 R 000022 RLXA      0002 R 000023 RLXD
0002 R 000036 R1      0002 R 000054 R2      0002 R 000037 S      0007 R 000000 T      0002 R 000065 TC6A
0002 R 000064 TC6D    0002 R 000047 TM      0002 R 000010 TPRINT  0002 R 000012 TSTEPN  0002 R 000011 TSUM
0002 R 000051 T1      0002 R 000052 T2      0002 R 000002 X      0015 R 000000 XK      0001 000162 10L
0001 002074 100L      0001 004212 1000L    0001 000420 1005L    0001 000441 1010L    000457 1012L
0001 000512 1015L    0001 000520 1020L    0001 000560 1025L    0001 000604 1030L    0001 000625 1032L
0001 000663 1035L    0001 000671 1040L    0001 000734 1045L    0001 002121 105L    0001 003623 10676
0001 002154 110L     0001 003674 11146    0001 003741 11336    0001 003764 11446    0001 002172 115L
0001 004027 11636    0001 002231 120L     0001 002242 125L     0001 002276 145L     0001 000171 15L
0001 002322 150L     0001 002345 155L     0001 002354 160L     0001 002367 165L     0001 002402 170L
0001 002446 175L     0001 002450 180L     0001 002767 185L     0001 003546 190L     0001 003551 195L
0001 000756 1998L    0001 000761 1999L    0001 000174 20L      0001 0001764 2000L    0001 001335 2005L
0001 001342 2007L    0001 001362 2010L    0001 001365 2015L    0001 001403 2017L    0001 001437 2020L
0001 001445 2025L    0001 001505 2030L    0001 001512 2032L    0001 001535 2035L    0001 001540 2040L
0001 001561 2042L    0001 001620 2045L    0001 001626 2050L    0001 001671 2055L    0001 001717 2060L
0001 001723 2065L    0001 003615 215L    0001 000266 220G    0001 003634 220L    0001 003641 225L
0001 000310 2326     0001 000314 2376     0001 003645 240L     0001 003666 245L     0001 003721 250L
0001 003735 255L     0001 000352 2606     0001 003754 265L     0001 004006 270L     0001 004023 275L
0001 004052 285L     0001 004102 290L     0001 001727 2998L    0001 001755 2999L    0001 000206 30L
0001 001760 3100L    0001 003055 3005L    0001 003663 3007L    0001 003103 3010L    0001 003107 3015L
0001 003126 3017L    0001 003162 3020L    0001 003170 3025L    0001 003231 3030L    0001 003237 3032L
0001 003262 3035L    0001 003266 3040L    0001 003310 3042L    0001 003347 3045L    0001 003355 3050L
0001 003421 3055L    0001 003450 3060L    0001 003454 3065L    0001 000221 35L      0001 003460 3098L
0001 003506 3099L    0001 000233 40L     0001 003451 4000L    0001 001035 4005L    0001 001042 4010L
0001 001043 4012L    0001 001062 4015L    0001 001063 4017L    0001 001100 4020L    0001 001114 4022L
0001 001131 4025L    0001 001137 4030L    0001 001175 4035L    0001 001211 4037L    0001 001226 4040L
0001 000237 45L      0001 001234 4998L    0001 001240 4999L    0001 000150 5L      0001 001243 5000L
0001 002554 5005L    0001 002562 5010L    0001 002563 5012L    0001 002603 5015L    0001 002604 5017L
0001 002621 5020L    0001 002635 5022L    0001 002652 5025L    0001 002660 5030L    0001 002716 5035L
0001 002732 5037L    0001 002747 5040L    0001 002127 5476    0001 002163 5636    0001 002755 5998L
0001 002762 5999L    0001 000343 60L      0001 002765 6000L    0001 002407 6406    0001 002473 6606
0001 002501 6556     0001 001253 70L     0001 001254 75L      0001 002040 80L     0000 000000 800F
0000 000006 801F     0000 000017 802F    0000 000025 803F    0000 000032 804F    0000 000037 805F
0000 000042 806F     0000 000045 807F    0000 000050 808F    0000 000054 809F    0001 004124 90CL
0001 004132 901L    0001 004140 903L    0001 004146 904L    0001 004155 905L    0001 004163 906L
0001 004171 907L    0001 004177 908L    0001 004205 909L

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00101 1* SUBROUTINE CNB/CK
00101 2* C IMPLICIT BACKWARD DIFFERENCING EXECUTION SURROUTINE
00101 3* C THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED, SINDA FORTRAN V
00101 4* C ALL NODES RECEIVE A SUCCESSIVE POINT ITERATION
00101 5* C RELAXATION CRITERIA MUST BE SPECIFIED
00101 6* C OVER-RELAXATION IS ALLOWED, THE DAMPENING FACTORS ARE ADDRESSABLE
00101 7* C INCLUDE COM-MLIST
00103 7* COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
00105 7* COMMON /PC1/NS01(1) /PC2/NS02(1) /KONST/K(1) /ARRAY/A(1)
00106 7* COMMON /FIXCON/KON(1) /XSPACE/NDIM,MTM,X(1)
00107 7* COMMON /DIMENS/ NND,NNA,NHT,NGT,NCT,NAT,LS01,LS02
00110 7* DIMENSION COH(1),XK(1),NK(1)
00111 7* EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NK(1))
00112 7* END
00113 8* INCLUDE DEFF,LIST
00113 8* C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00113 8* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEH)
00113 8* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00113 8* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
00113 8* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR*EXPLICIT (CS0FAC)

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CNRACK.CNBACK

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00113 8* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00113 8* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00113 8* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
00113 8* C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
00113 8* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPF)
00113 8* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
00113 8* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
00113 8* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
00113 8* C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
00113 8* C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEH)
00113 8* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPC)
00113 8* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (CSGMTH)
00113 8* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (OUTPUT)
00113 8* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 8* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 8* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 8* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
00113 8* C CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
00113 8* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
00113 8* C CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
00113 8* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 8* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 8* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 8* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 8* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 8* C CC31 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 8* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENIGBAL)
00113 8* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 8* C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 8* C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 8* C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 8* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 8* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPC
00113 8* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 8* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 8* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINOSM (LAXFAC)
00113 8* C CC50 IS NOT USED AT PRESENT
00114 8* C ENID
00115 9* IF (KOR(5).LE.0) GO TO 999
00117 10* IF (CON(6).LE.0) CON(6) = 1.E+8
00121 11* IF (CON(8).LE.0) CON(8) = 1.E+8
00123 12* IF (CON(9).LE.0) CON(9) = 1.0
00125 13* IF (CON(10).LE.0) CON(10) = 1.0
00127 14* IF (CON(11).LE.0) CON(11) = 1.E+8
00131 15* IF (CON(13).LE.0) CON(13) GO TO 990
00133 16* IF (CON(18).LE.0) GO TO 998
00135 17* IF (NNA-GT.0.AND.CON(19).LE.0.) GO TO 997
00137 18* IF (CON(22).LE.0.) GO TO 996
00141 19* IF (IND-GT.0.AND.CON(26).LE.0.) GO TO 995
00143 20* IF (KOR(31).NE.1) GO TO 991
00145 21* PASS = -1.0
00146 22* NN = NND+1
00147 23* NLA = NDIM
00150 24* IEL = NTH
00151 25* NPI = NND+NNA
00152 26* IEL2 = NTH+NNT
00153 27* IEL3 = IEL2+NND
00154 28* J = 2*NND+NNT
00155 29* NTH = NTH+J

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CNBACK,CNBACK

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30* 00156 NDIM = NDIM-J
31* 00156 CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
32* 00157 IF (NDIM.LT.0) GO TO 994
33* 00161 IPRINT = CON(13)
34* 00161 INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
35* 00162 TSUM = 0.0
36* 00162 DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
37* 00163 IF (CON(13)+CON(18).GT.CON(3)) CON(18) = CON(3)-CON(13)
38* 00163 DONT EXCEED IT
39* 00165 TSTEPN = CON(22)
40* 00166 IF (TSTEPN.LE.CON(8)) GO TO 20
41* 00170 TSTEPN = CON(8)
42* 00171 GO TO 35
43* 00171 DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
44* 00172 IF (TSUM+TSTEPN-CON(18)) 30,35,25
45* 00172 DONT EXCEED IT
46* 00175 TSTEPN = CON(18)-TSUM
47* 00176 GO TO 35
48* 00176 DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
49* 00177 IF (TSUM+2.0*TSTEPN.LE.CON(18)) GO TO 35
50* 00177 APPROACH THE OUTPUT INTERVAL GRADUALLY
51* 00201 TSTEPN = (CON(18)-TSUM)/2.0
52* 00201 STORE DELTA TIME STEP IN THE CONSTANTS
53* 00202 CON(2) = TSTEPN
54* 00202 CALCULATE THE NEW TIME
55* 00203 IF (PASS.GT.0.) GO TO 40
56* 00205 CON(1) = IPRINT
57* 00206 CON(2) = 0.0
58* 00207 GO TO 45
59* 00210 CON(1) = IPRINT+TSUM+TSTEPN
60* 00210 COMPUTE THE MEAN TIME BETWEEN ITERATIONS
61* 00211 CON(14) = (CON(1)+CON(13))/2.0
62* 00212 LAX = KON(5)
63* 00213 DN = CON(10)
64* 00214 DD = 1.0-DN
65* 00215 AN = CON(9)
66* 00216 AA = 1.0-AN
67* 00216 DO THE RELAXATION LOOP
68* 00217 DO 240 K1 = 1,LAX
69* 00222 KON(20) = K1
70* 00223 J1 = 0
71* 00224 RLXA = 0.0
72* 00225 RLXD = 0.0
73* 00226 IF (K1.GT.1) GO TO 110
74* 00230 J2 = 1
75* 00230 ZERO OUT ALL SOURCE LOCATIONS AND SHIFT TEMPERATURES
76* 00231 DO 50 I = 1,NNC
77* 00234 Q(I) = 0.0
78* 00236 DO 55 I = 1,NNT
79* 00241 LE1 = IE1+I
80* 00242 X(LE1) = Y(I)
81* 00244 KON(12) = 0
82* 00245 CALL VARBL1
83* 00245 CHECK THE BACKUP SWITCH
84* 00246 IF (KON(12).NE.0) GO TO 15
85* 00246 CHECK FOR FIRST PASS
86* 00250 IF (PASS.GE.0.) GO TO 60
87* 00252 CALL OUTCAL
88* 00253 PASS = 1.0
89* 00254 GO TO 10

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CNBACK,CNBACK

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00255 90*
00256 91*
00257 92*
00258 93*
00259 94*
00260 95*
00261 95*
00262 95*
00263 95*
00264 95*
00265 95*
00266 95*
00267 95*
00268 95*
00269 95*
00270 95*
00271 95*
00272 95*
00273 95*
00274 95*
00275 95*
00276 95*
00277 95*
00300 95*
00301 95*
00302 95*
00303 95*
00304 95*
00305 95*
00306 95*
00307 95*
00310 95*
00311 95*
00312 95*
00313 95*
00314 95*
00315 95*
00316 95*
00317 95*
00320 95*
00321 95*
00322 95*
00323 95*
00324 95*
00325 95*
00326 95*
00327 95*
00330 95*
00331 95*
00332 95*
00333 95*
00334 95*
00335 95*
00336 95*
00337 96*
00340 97*
00341 98*
00342 99*
00343 100*
00344 100*
00345 100*
00346 100*
00347 100*
00350 100*
00351 100*
00352 100*

60 RC = 1.E+8
JJ = 0
C CALCULATE FIRST PASS TEMPERATURES AND CSGMIN
DO 105 I = 1,NMP
INCLUDE VARC,LIST
FOLD DELTAT INTO THE CAPACITANCES*
IF (FLD(1,1,NSQ(1,J1+1)),EQ,0) GO TO 2000
NTYPE = FLD(0,5,NSQ2(J2))
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
1005 CALL DID1WN(T(I),A(LA),XK(LK),C(I))
GO TO 1999
1010 CALL DID1WN(T(I),A(LA),XK(LK),C1)
1012 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
CALL DID1WN(T(I),A(LA),XK(LK),C2)
GO TO 1998
1015 C1 = XK(LK)*XK(LA)
GO TO 1012
1020 CALL DID1WN(T(I),A(LA),XK(LK),C1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
C2 = XK(LK)*XK(LA)
GO TO 1998
1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
GO TO 1999
1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
1032 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
GO TO 1998
1035 C1 = XK(LK)*XK(LA)
GO TO 1032
1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
C2 = XK(LK)*XK(LA)
GO TO 1998
1045 CALL D2D1WN(T(I),CON(14),A(LA),XK(LK),C(I))
GO TO 1999
1996 C(I) = C1+C2
1999 J2 = J2+1
2000 CONTINUE
END
C(I) = C(I)/TSTEPN
R1 = 0.0
S = 0.0
G2 = 0.0
INCLUDE VARC,LIST
IF (FLD(4,1,NSQ(1,J1+1)),EQ,0) GO TO 5000
NTYPE = FLD(0,5,NSQ2(J2))
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4045), NTYPE
4005 G(I) = XK(LK)+G(I)

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*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-3

CNBACK,CNBACK

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00353 100*
00354 100*
00355 100*
00356 100*
00357 100*
00360 100*
00361 100*
00363 100*
00364 100*
00365 100*
00366 100*
00367 100*
00370 100*
00371 100*
00372 100*
00373 100*
00374 100*
00375 100*
00376 100*
00377 100*
00400 100*
00401 100*
00402 100*
00403 100*
00404 100*
00405 100*
00406 100*
00407 100*
00410 100*
00411 100*
00412 101*
00413 102*
00414 103*
00415 104*
00416 105*
00417 106*
00420 107*
00421 107*
00423 107*
00424 107*
00425 107*
00426 107*
00426 107*
00427 107*
00430 107*
00431 107*
00432 107*
00433 107*
00434 107*
00435 107*
00436 107*
00437 107*
00440 107*
00441 107*
00442 107*
00443 107*
00444 107*
00445 107*
00446 107*

60 TO 4999
4010 Q1 = 0-0
4012 CALL DID1WM(T(I),A(LA),XK(LK),Q2)
GO TO 4998
4015 Q1 = 0-0
4017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
GO TO 4998
4020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GO TO 4017
4025 Q1 = XK(LK)*XK(LA)
GO TO 4022
4030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
Q2 = XK(LK)*XK(LA)
GO TO 4998
4035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
4037 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GO TO 4012
4040 Q1 = XK(LK)*XK(LA)
GO TO 4037
4998 Q(I) = Q1+Q2+Q(I)
4999 J2 = J2+1
5000 CONTINUE
END
Q(I) = Q(I)+C(I)*(T(I)+460.0)
QSUM = Q(I)
QSUM = C(I)
65UM = C(I)
70 J1 = J1+1
LG = FLD(5,16,NSQ1(J1))
LTA = FLD(22,14,NSQ1(J1))
INCLUDE VARG,LIST
IF(FLD(2,1,NSQ1(J1)),EQ,0) GO TO 3000
NTYPE = FLD(0,5,NSQ2(J2))
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GOTO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
2060,2065), NTYPE
2005 TM = (T(I)+LTA)/2.0
2007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
GO TO 2999
2010 TM = T(I)
GO TO 2007
2015 CALL DID1WM(T(I),A(LA),XK(LK),G1)
2017 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
GO TO 2998
2020 G1 = XK(LK)*XK(LA)
GO TO 2017
2025 CALL DID1WM(T(I),A(LA),XK(LK),G1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))

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*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW

*NEW
***2

*NEW
*NEW
***2

*NEW
*NEW
***2

CNBACK,CNBACK

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00447 107* LK = FLD(22,14,NSQ2(J2))
00450 107* G2 = XK(LK)*XK(LA)
00451 107* GO TO 2998
00452 107* 2030 TM = (T(I)+T(LTA))/2.0
00453 107* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00454 107* GO TO 2999
00455 107* 2035 TM = T(I)
00456 107* GO TO 2032
00457 107* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00460 107* 2042 J2 = J2+1
00461 107* LA = FLD(5,17,NSQ2(J2))
00462 107* LK = FLD(22,14,NSQ2(J2))
00463 107* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00464 107* GO TO 2998
00465 107* 2045 G1 = XK(LK)*XK(LA)
00466 107* GO TO 2042
00467 107* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00470 107* J2 = J2+1
00471 107* LA = FLD(5,17,NSQ2(J2))
00472 107* LK = FLD(22,14,NSQ2(J2))
00473 107* G2 = XK(LK)*XK(LA)
00474 107* GO TO 2998
00475 107* 2055 TM = (T(I)+T(LTA))/2.0
00476 107* CALL D2DIRM(TM,CON(14),A(LA),XK(LK),G(LG))
00477 107* GO TO 2999
00500 107* 2060 TM = T(LTA)
00501 107* GO TO 2007
00502 107* 2065 TM = T(LTA)
00503 107* GO TO 2032
00504 107* 2998 G(LG) = 1./(1./G1+1./G2)
00505 107* IF(FLD(3,1,NSQ1(J1)).EQ.0) G(LG) = G1*G2
00507 107* 2999 J2 = J2+1
00510 107* 3000 CONTINUE
00511 107* ENO
00512 108* T1 = T(I)+460.0
00513 109* T2 = T(LTA)+460.0
00514 110* C CHECK FOR RADIATION CONDUCTOR
00515 111* IF(FLD(3,1,NSQ1(J1)).EQ.0) GO TO 75
00516 112* R1 = R1+G(LG)
00517 113* GSUM = GSUM+G(LG)+T2**4
00520 114* G2 = G2+G(LG)*(T1+T1+T2*T2)*(T1+T2)
00521 115* GO TO 80
00522 116* 75 GV = G(LG)
00523 117* G2 = G2+GV
00524 118* GSUM = GSUM+GV
00525 119* GSUM = GSUM+GV*T2
00525 120* C CHECK FOR LAST CONDUCTOR
00526 121* IF(INSQ1(J1).GT.0) GO TO 70
00526 122* DAMPEN RADIATION ON THIS NODE IF PRESENT
00530 123* IF(R1.LE.0.) GO TO 100
00532 124* R2 = R1+T1**4
00533 125* T2 = (GSUM-R2)/GSUM
00534 126* R1 = R1+T2**4
00535 127* S = (R1+R2)/2.*n
00535 128* C OBTAIN THE NEW TEMPERATURE
00536 129* 100 T(I) = (DN*((GSUM-S)/GSUM)+DD*T1)+460.0
00537 130* R1 = C(I)/G2
00540 131* IF(R1.GE.RC) GO TO 105
00542 132* RC = R1
00543 133* KOT(35) = I

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CNBACK-CNBACK

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00544 134*
00544 135*
00546 136*
00551 137*
00552 138*
00553 139*
00555 140*
00556 141*
00560 143*
00561 144*
00562 145*
00565 146*
00566 147*
00567 148*
00570 149*
00571 150*
00572 151*
00573 152*
00574 153*
00574 154*
00576 155*
00577 156*
00600 157*
00601 158*
00602 159*
00602 160*
00603 161*
00603 162*
00605 163*
00607 164*
00610 165*
00611 166*
00612 167*
00612 168*
00613 169*
00613 170*
00614 171*
00614 172*
00615 173*
00616 174*
00617 175*
00620 176*
00621 177*
00622 178*
00623 179*
00624 180*
00625 181*
00626 182*
00627 183*
00631 184*
00632 185*
00633 186*
00635 187*
00635 188*
00636 189*
00637 190*
00642 191*
00642 192*
00643 193*

C 105 CONTINUE
C CONVERT TEMPERATURES TO RANKINE
DO 65 I = 1,NNT
LE1 = IE1+I
T(I) = T(I)+460.
65 X(LE1) = X(LE1)+460.
CON(17) = RC+TSTEPN
IF(RC.LE.0.) GO TO 993
GO TO 225

C 110 JJ = J+1
DO 165 I = 1,NHD
R1 = 0.0
S = 0.0
GSUM = G(I)
GSUM = C(I)
115 J1 = J+1
LG = FLD(5,16,NSQ1(J1))
LTA = FLD(22,14,NSQ1(J1))
CHECK FOR RADIATION CONDUCTOR
IF(FLD(3,1,NSG1(J1)).EQ.0) GO TO 120
R1 = R1+G(LG)
GSUM = GSUM+G(LG)*T(LTA)**4
GO TO 125

120 GSUM = GSUM+G(LG)
GSUM = GSUM+G(LG)*T(LTA)
CHECK FOR LAST CONDUCTOR

C 125 IF(NSQ1(J1).GT.0) GO TO 115
IF(R1.LE.0.) GO TO 145
R2 = R1*T(I)**4
T2 = (GSUM-R2)/GSUM
R1 = R1+T2**4
S = (R1+R2)/2.0
OBTAIN THE NEW TEMPERATURE

C 145 T2 = DN*((GSUM-S)/GSUM)+DO*T(I)
OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
T1 = ABS(T(I)-T2)
STORE THE NEW AND OLD TEMPERATURES
GO TO (160,155,150), JJ
LE2 = IE2+1
LE3 = IE3+1
R1 = T2-T(I)
X(LE2) = T(I)
X(LE3) = R1/(R1-X(LE3))
GO TO 160

155 LE3 = IE3+1
X(LE3) = T2-T(I)
160 T(I) = T2
IF(RLXD.GE.T1) GO TO 165
RLXD = T1
KK1 = I

165 CONTINUE
GO TO (180,180,170), JJ
PERFORM LINEAR EXTRAPOLATION ON THE ERROR FUNCTION CURVE

C 170 JJ = 0
DO 175 I = 1,NHD
LE3 = IE3+1
SEE IF THE EXTRAPOLATION IS ALLOWABLE
IF(X(LE3).GE.0.) GO TO 175

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CNBACK,CNBACK

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00643 194* LIMIT THE EXTRAPOLATION
00645 195* IF(X(LE3).LT.-10.) X(LE3) = -10.
00647 196* LE2 = IZ2+1
00650 197* T(I) = X(LE3)*X(LE2)+(1.0-X(LE3))*T(I)
00651 198*
00653 199* 175 CONTINUE
160 IF(NMIA.LE.0) GO TO 220
    JJI = J1
    JJ2 = J2
    DO 230 I = 1,NNT
230 T(I) = T(I)-460.0
    DO 215 I = NN,NIRC
    L = I
    GSUM = 0.0
    IF(K1.GT.2) GO TO 6000
    INCLUDE VRG2*LIST
    IF (FLD(4,1,NSQ1(JJ1+1)),EQ.0) GO TO 6000
    NTYPE = FLD(10,5,NSQ2(JJ2))
    LA = FLD(5,17,NSQ2(JJ2))
    LK = FLD(22,14,NSQ2(JJ2))
    GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5050), NTYPE
5005 Q(L) = XK(LK)+Q(L)
    GO TO 5999
5010 Q1 = 0.0
5012 CALL DID1WM(T(L),A(LA),XK(LK),Q2)
    GO TO 5998
5015 Q1 = 0.0
5017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
    GO TO 5998
5020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
5022 JJ2 = JJ2+1
    LA = FLD(5,17,NSQ2(JJ2))
    LK = FLD(22,14,NSQ2(JJ2))
    GO TO 5017
5025 Q1 = XK(LK)*XK(LA)
    GO TO 5022
5030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
    JJ2 = JJ2+1
    LA = FLD(5,17,NSQ2(JJ2))
    LK = FLD(22,14,NSQ2(JJ2))
    Q2 = XK(LK)*XK(LA)
    GO TO 5998
5035 CALL DID1WM(CO1(14),A(LA),XK(LK),G1)
5037 JJ2 = JJ2+1
    LA = FLD(5,17,NSQ2(JJ2))
    LK = FLD(22,14,NSQ2(JJ2))
    GO TO 5012
5040 Q1 = XK(LK)*XK(LA)
    GO TO 5037
5998 Q(L) = Q1+Q2+Q(L)
5999 JJ2 = JJ2+1
6000 CONTINUE
    ENDO
    GSUM = Q(I)
185 JJI = JJI+1
    LG = FLD(5,16,NSQ1(JJ1))
    LIA = FLD(22,14,NSQ1(JJ1))
    IF(K1.GT.2) GO TO 4000
    INCLUDE VRG2*LIST
    IF(FLD(12,1,NSQ1(JJ1)),EQ.0) GO TO 4000
    NTYPE = FLD(10,5,NSQ2(JJ2))

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*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW

CNRACK, CNRACK

00754	214*	LA = FLD(5,17,NS02(JJ2))	*NEW
00755	214*	LK = FLD(22,14,NS02(JJ2))	*NEW
00756	214*	GO TO (3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,	*-3
00757	214*	3060,3065), NTYPE	
00760	214*	3005 TM = (T(L)+T(LTA))/2.0	
00761	214*	3007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))	
00762	214*	GO TO 3999	
00763	214*	3010 TM = T(L)	
00764	214*	GO TO 3007	
00765	214*	3015 CALL DID1WM(T(L),A(LA),XK(LK),61)	
00766	214*	3017 JJ2 = JJ2+1	
00767	214*	LA = FLD(5,17,NS02(JJ2))	
00770	214*	LK = FLD(22,14,NS02(JJ2))	
00771	214*	CALL DID1WM(T(LTA),A(LA),XK(LK),62)	
00772	214*	GO TO 3998	
00773	214*	3020 G1 = XK(LK)*XK(LA)	
00774	214*	GO TO 3017	
00775	214*	3025 CALL DID1WM(T(L),A(LA),XK(LK),61)	
00776	214*	JJ2 = JJ2+1	
00777	214*	LA = FLD(5,17,NS02(JJ2))	
01000	214*	LK = FLD(22,14,NS02(JJ2))	
01001	214*	G2 = XK(LK)*XK(LA)	
01002	214*	GO TO 3998	
01003	214*	3030 TM = (T(L)+T(LTA))/2.0	
01004	214*	3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))	
01005	214*	GO TO 3999	
01006	214*	3035 TM = T(L)	
01007	214*	GO TO 3032	
01010	214*	3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
01011	214*	3042 JJ2 = JJ2+1	
01012	214*	LA = FLD(5,17,NS02(JJ2))	
01013	214*	LK = FLD(22,14,NS02(JJ2))	
01014	214*	CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)	
01015	214*	GO TO 3998	
01016	214*	3045 G1 = XK(LK)*XK(LA)	
01017	214*	GO TO 3042	
01020	214*	3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
01021	214*	JJ2 = JJ2+1	
01022	214*	LA = FLD(5,17,NS02(JJ2))	
01023	214*	LK = FLD(22,14,NS02(JJ2))	
01024	214*	G2 = XK(LK)*XK(LA)	
01025	214*	GO TO 3998	
01026	214*	3055 TM = (T(L)+T(LTA))/2.0	
01027	214*	CALL P2D1WM(TM,CON(14),A(LA),XK(LK),6(LG))	
01030	214*	GO TO 3999	
01031	214*	3060 TM = T(LTA)	
01032	214*	GO TO 3007	
01033	214*	3065 TM = T(LTA)	
01034	214*	GO TO 3032	
01035	214*	3998 G(LG) = 1./(1./G1+1./G2)	
01037	214*	IF(FLD(3,1,NS01(JJ1)).EQ.1) G(LG) = G1*G2	
01040	214*	3999 JJ2 = JJ2+1	
01041	214*	4000 CONTINUE	
01042	215*	END	
01043	216*	T1 = T(I)+460.0	
01044	217*	T2 = T(LTA)+460.0	
01046	219*	C CHECK FOR RADIATION CONDUCTOR	
01047	220*	IF (FLD(3,1,NS01(JJ1)).EQ.0) GO TO 190	
		GV = G(LG)*(T1+T2+T2)*T1+T2	
		GO TO 195	

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01050 221*
01051 222*
01052 223*
01053 224*
01054 225*
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01098 269*
01099 270*
01100 271*
01101 272*
01102 273*
01103 274*
01104 275*
01105 276*
01106 277*
01107 278*
01108 279*
01109 280*

190 GV = G(LG)
195 GSUM = GSUM+GV
    GSUM = GSUM+GV*T2
C CHECK FOR LAST CONDUCTOR
  IF(NSG(I,JJ1).GT.0) GO TO 185
C CALCULATE THE NEW TEMPERATURE
  T2 = AN*GSUM/GSUM+AA*T1
  T1 = ABS(T2-T1)
  IF(RLXA.GE.T1) GO TO 215
  RLXA = T1
  KK2 = I
215 CONTINUE
  DO 235 I = 1,N*NT
  T(I) = T(I)+460.0
C SEE IF THE ARITHMETIC RELAXATION CRITERIA WAS MET
  IF(RLXA.GT.CON(19)) GO TO 225
C SEE IF THE DIFFUSION RELAXATION CRITERIA WAS MET
  IF(RLXD.LE.CON(26)) GO TO 245
  IF(KON(7).EQ.0) GO TO 240
  CALL OUTCAL
240 CONTINUE
  IF(KON(28).GE.65) CALL TOPLIN
  WRITE(6,982)
  KON(28) = KON(28)+2
C SEE IF THE TEMPERATURE CHANGES WERE TOO LARGE
  TCGD = 0.0
  TCGA = 0.0
  DO 250 I = 1,N*ND
  LE = IEI+I
  C(I) = C(I)*TSTEPN
  T1 = ABS(T(I)-X(LE))
  IF(TCGD.GT.T1) GO TO 250
  TCGD = T1
  KON(36) = I
250 CONTINUE
  IF(TCGD.LE.CON(6)) GO TO 265
  TSTEPN = 0.95*TSTEPN*CON(6)/TCGD
255 DO 260 I = 1,N*NT
  LE = IEI+I
  T(I) = X(LE)-460.0
  GO TO 30
265 IF(NHA.LE.0) GO TO 275
  DO 270 I = NN,NNC
  LE = IEI+I
  T1 = ARS(T(I)-X(LE))
  IF(TCGA.GT.T1) GO TO 270
  TCGA = T1
  KON(38) = I
270 CONTINUE
  IF(TCGA.LE.CON(11)) GO TO 275
  TSTEPN = 0.95*TSTEPN*CON(11)/TCGA
  GO TO 255
C CONVERT TEMPERATURES BACK TO FARENHEIT
275 DO 280 I = 1,N*NT
280 T(I) = T(I)-460.0
C STORE THE TEMPERATURE AND RELAXATION CHANGES
  CON(15) = TCGD
  CON(16) = TCGA
  CON(27) = RLXD

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CNBACK*CNBACK

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01172 281* IF(RLXA.GT.RLXD) GO TO 285
01174 282* KK2 = KK1
01175 283* RLXA = RLXD
01176 284* 285 KON(37) = KK2
01177 285* KON(30) = RLXA
01200 286* KON(12) = 0
01201 287* CALL VARBL2
01201 288* CHECK THE BACKUP SWITCH
01202 289* IF(KON(12).NE.0) GO TO 255
01202 290* C ADVANCE TIME
01204 291* C CON(13) = CON(1)
01205 292* TSUM = TSUM+TSTEPN
01205 293* C CHECK FOR TIME TO PRINT
01206 294* IF(TSUM.GE.CON(18)) GO TO 290
01206 295* C CHECK FOR PRINT EVERY ITERATION
01210 296* IF(KON(7).NE.0) CALL OUTCAL
01212 297* GO TO 10
01212 298* C TRY TO EVEN THE OUTPUT INTERVALS
01213 299* C 290 TPRINT = TPRINT+TSUM
01214 300* CALL OUTCAL
01214 301* C IS TIME GREATER THAN END COMPUTE TIME
01215 302* IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
01217 303* NTH = IE1
01220 304* NDIM = NLA
01221 305* RETURN
01222 306* 990 WRITE(6,880)
01224 307* GO TO 1000
01225 308* 981 WRITE(6,881)
01227 309* GO TO 1000
01230 310* 993 WRITE(6,883)
01232 311* GO TO 1000
01233 312* 994 WRITE(6,884) NDIM
01236 313* 995 WRITE(6,885)
01241 315* GO TO 1000
01242 316* 996 WRITE(6,886)
01244 317* GO TO 1000
01245 318* 997 WRITE(6,887)
01247 319* GO TO 1000
01250 320* 998 WRITE(6,888)
01252 321* GO TO 1000
01253 322* 999 WRITE(6,889)
01255 323* 1000 CALL OUTCAL
01256 324* CALL EXIT
01257 325* 880 FORMAT(29H TRANSIENT TIME NOT SPECIFIED)
01260 326* 881 FORMAT(45H CNBACK REQUIRES LONG PSEUDO-COMPUTE SEQUENCE)
01261 327* 882 FORMAT(29H RELAXATION CRITERIA NOT MET)
01263 328* 883 FORMAT(19,20H CG-MIN ZERO OR NEGATIVE)
01264 329* 884 FORMAT(19,20H LOCATIONS AVAILABLE)
01264 330* 885 FORMAT(10H NO (RLXCA))
01265 331* 886 FORMAT(10H NO (TIMEI))
01266 332* 887 FORMAT(10H NO (ARLXCA))
01267 333* 888 FORMAT(19H NO OUTPUT INTERVAL)
01270 334* 889 FORMAT(9H NO NLOOP)
01271 335* END

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END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S) 12 JAN 70 23:06:44 0 42603742 14 335 (DELETED)

CNBACK SYMBOLIC 12 JAN 70 23:06:44 0 42615064 84 1 (DELETED)

CNBACK CODE HELCCATABLE 12 JAN 70 23:06:44 0 42615210 14 226

DIW FOR.** CNFWBK.CNFWBK
UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-100 CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:43

21 FEB 71

SUBROUTINE CNFWBK ENTRY POINT 004271

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	004307
0000	*CONST+TEMP	000127
0002	*SIMPLE VAR	000071
0004	*ARAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	CONC	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000301
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARBL1
0023	OUTCAL
0024	DID1WM
0025	PLYAWM
0026	D2D1WM
0027	TOPLIN
0030	VARBL2
0031	EXIT
0032	NERR2\$
0033	NWDUS
0034	NI02\$
0035	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0002	R	000017	AA
0002	R	000035	C1	0002	R	000015	DD
0002	R	000045	GSUM	0002	R	000052	G1
0002	I	000026	I	0002	I	000005	IE2
0002	I	000031	JJ	0002	I	000063	JJ2
0015	I	000000	K	0002	I	000065	KK2
0002	I	000064	L	0002	I	000013	LAX
0002	I	000057	LE2	0002	I	000046	LG
0021	I	000007	LS02	0002	I	000050	LTAE
0020	I	000000	NJIM	0002	I	000002	NLA
0002	I	000004	RNC	0021	I	000002	RHT
0020	I	000001	NTH	0002	I	000032	NTYPE
0017	R	000030	CON	0002	R	000016	AN
0012	R	000000	G	0002	R	000015	DD
0006	R	000000	H	0002	R	000052	G1
0002	I	000007	J	0002	I	000005	IE2
0002	I	000025	J2	0002	I	000063	JJ2
0002	I	000021	K1	0002	I	000065	KK2
0002	I	000027	LE1	0002	I	000013	LAX
0021	I	000006	LS01	0002	I	000046	LG
0021	I	000004	NCT	0002	I	000050	LTAE
0021	I	000001	NNA	0002	I	000002	NLA
0014	I	000000	N502	0021	I	000002	RHT
0011	R	000000	G	0002	R	000032	NTYPE

CNF, WBK, CNF, BK

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0002 R 000044 QSUM          0002 R 000042 Q1          0002 R 000043 Q2          0002 R 000030 RC          0002 R 000023 RLXA
0002 R 000024 RLXD          0002 R 000037 R1          0002 R 000056 R2          0002 R 000040 S          0007 R 000000 Y
0002 R 000067 TGA          0002 R 000066 TCGD          0002 R 000051 TM          0002 R 000010 TSTEP          0002 R 000020 YSTEP
0002 R 000012 TSTEPN          0002 R 000011 TSGM          0002 R 000053 T1          0002 R 000054 T2          0002 R 000002 X
0015 R 000000 XK          0001 000162 10L          0001 000126 100L          0001 004244 1000L          0002 R 000023 1005L
0001 000444 1010L          0001 000462 1012L          0001 000515 1015L          0001 000523 1020L          0001 000563 1025L
0001 000607 1030L          0001 000630 1032L          0001 000666 1035L          0001 000674 1040L          0001 000737 1045L
0001 002153 105L          0001 003655 1073G          0001 002206 110L          0001 003226 110L          0001 003773 1137G
0001 002224 115L          0001 004016 1150G          0001 004061 1167G          0001 002263 120L          0001 002274 125L
0001 002330 145L          0001 000171 15L          0001 002377 155L          0001 002406 160L          0001 002406 160L
0001 002421 165L          0001 002434 170L          0001 002500 175L          0001 003021 185L          0001 003021 185L
0001 003600 190L          0001 003603 195L          0001 000761 1998L          0001 000764 1999L          0001 000174 20L
0001 000767 2000L          0001 001344 2005L          0001 001351 2007L          0001 001371 2010L          0001 001374 2015L
0001 001412 2017L          0001 001446 2020L          0001 001454 2025L          0001 001514 2030L          0001 001521 2032L
0001 001544 2035L          0001 001547 2040L          0001 001570 2042L          0001 001627 2045L          0001 001635 2050L
0001 001700 2055L          0001 001726 2060L          0001 001732 2065L          0001 003647 215L          0001 003666 220L
0001 000271 221G          0001 003673 225L          0001 000313 233G          0001 000317 240G          0001 003677 240L
0001 003720 245L          0001 003753 250L          0001 003767 255L          0001 000355 261G          0001 004006 265L
0001 004040 270L          0001 004055 275L          0001 004104 285L          0001 004134 290L          0001 001736 2998L
0001 001764 2999L          0001 000206 30L          0001 001767 3000L          0001 003107 3005L          0001 003115 3007L
0001 003135 3010L          0001 003141 3015L          0001 003160 3017L          0001 003214 3020L          0001 003222 3025L
0001 003263 3030L          0001 003271 3032L          0001 003314 3035L          0001 003320 3040L          0001 003342 3042L
0001 003401 3045L          0001 003407 3050L          0001 003453 3055L          0001 003502 3060L          0001 003506 3065L
0001 000221 30L          0001 003512 3098L          0001 003540 3099L          0001 000233 40L          0001 003543 4000L
0001 001040 4005L          0001 001045 4010L          0001 001046 4012L          0001 001065 4015L          0001 001066 4017L
0001 001103 4020L          0001 001117 4022L          0001 001134 4025L          0001 001142 4030L          0001 001200 4035L
0001 001214 4037L          0001 001231 4040L          0001 001231 4040L          0001 001237 4998L          0001 001243 4999L
0001 000150 5L          0001 001246 5000L          0001 002605 5005L          0001 002613 5010L          0001 002614 5012L
0001 002634 5015L          0001 002635 5017L          0001 002652 5020L          0001 002666 5022L          0001 002703 5025L
0001 002711 5030L          0001 002747 5030L          0001 002763 5037L          0001 003046 60L          0001 002161 553G
0001 002215 507G          0001 003006 5098L          0001 003013 5099L          0001 000346 60L          0001 003016 6000L
0001 002441 644G          0001 002525 664G          0001 002533 671G          0001 001260 70L          0001 002047 75L
0001 002072 80L          0000 000000 80F          0000 000006 80F          0000 000017 802F          0000 000025 803F
0000 000032 804F          0000 000037 805F          0000 000042 806F          0000 000045 807F          0000 000050 808F
0000 000054 809F          0001 004156 900L          0001 004164 901L          0001 004172 903L          0001 004200 904L
0001 004207 905L          0001 004215 906L          0001 004223 907L          0001 004231 909L          0001 004237 909L

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00101 SURROUTINE CNF, BK
00101 1* C
00101 2* C IMPLICIT FORWARD-BACKWARD DIFFERENCING EXECUTION SUBROUTINE
00101 3* C THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED, SINDA FORTRAN V
00101 4* C ALL NODES RECEIVE A SUCCESSIVE POINT ITERATION
00101 5* C RELAXATION CRITERIA MUST BE SPECIFIED
00101 6* C OVER-RELAXATION IS ALLOWED, THE DAMPENING FACTORS ARE ADDRESSABLE
00101 7* C INCLUDE COMMLIST
00101 8* C COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
00101 9* C COMMON /PC1/NSQ(1) /PC2/NSR2(1) /KONST/K(1) /ARRAY/A(1)
00101 10* C COMMON /FIXCON/KON(1) /XSPACE/NDIM, NTHX(1)
00101 11* C COMMON /DIMENS/ NMD, NNA, NITNGT, NCT, NAT, L501, L502
00101 12* C DIMENSION CON(1), XK(1), NK(1)
00101 13* C EQUIVALENCE (KON(1), CON(1)), (K(1), XK(1)), (X(1), NK(1))
00101 14* C END
00101 15* C ***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00101 16* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
00101 17* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00101 18* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
00101 19* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR EXPLICIT (CSGFAC)

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00113 8* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
 00113 8* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
 00113 8* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPETRH)
 00113 8* C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
 00113 8* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
 00113 8* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
 00113 8* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
 00113 8* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
 00113 8* C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
 00113 8* C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
 00113 8* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPC)
 00113 8* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPC)
 00113 8* C CC17 CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSGMIN)
 00113 8* C CC18 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
 00113 8* C CC19 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
 00113 8* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
 00113 8* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (CSGMAX)
 00113 8* C CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
 00113 8* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
 00113 8* C CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
 00113 8* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
 00113 8* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
 00113 8* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
 00113 8* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
 00113 8* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
 00113 8* C CC31 IS INDICATOR, 0-THERMAL SPCS, 1-THERMAL LPCS, 2-GENERAL (LSPCS)
 00113 8* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGRAL)
 00113 8* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
 00113 8* C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
 00113 8* C CC35 CONTAINS RELATIVE NODE NUMBER OF -CSGMIN
 00113 8* C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
 00113 8* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
 00113 8* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPC
 00113 8* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
 00113 8* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
 00113 8* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
 00113 8* C CC50 IS NOT USED AT PRESENT
 00114 8* C
 00115 9* IF (CON(5).LE.0) GO TO 999
 00117 10* IF (CON(6).LE.0) CON(6) = 1.E+8
 00121 11* IF (CON(8).LE.0) CON(8) = 1.E+8
 00123 12* IF (CON(9).LE.0) CON(9) = 1.0
 00125 13* IF (CON(10).LE.0) CON(10) = 1.0
 00127 14* IF (CON(11).LE.0) CON(11) = 1.E+8
 00131 15* IF (CON(3).LE.0) CON(13) GO TO 990
 00133 16* IF (CON(18).LE.0) GO TO 998
 00135 17* IF (NNA.GT.0.AND.CON(19).LE.0) GO TO 997
 00137 18* IF (CON(22).LE.0) GO TO 996
 00141 19* IF (NND.GT.0.AND.CON(26).LE.0) GO TO 995
 00143 20* IF (CON(31).NE.1) GO TO 991
 90145 21* PASS = -1.0
 90146 22* NND = NND+1
 90147 23* NLA = NNDIM
 90150 24* NTH = NTH
 90151 25* NNC = NND+NNA
 90152 26* NTH = NTH+NNT
 90153 27* NLS = 1E2+NND
 90154 28* J = 2*NND+NNT
 90155 29* NTH = NTH+J

CNFWBK,CNF*BK

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00156 NDIM = NDIM-J
00157 CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00158 IF(INDIM.LT.0) GO TO 994
00159 TPRINT = CON(13)
00160 INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00161 TSUM = 0.0
00162 DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00163 IF(CON(13)+CON(18).GT.CON(3)) CON(18) = CON(3)-CON(13)
00164 DONT EXCEED IT
00165 TSTEPN = CON(22)
00166 IF(TSTEPN.LE.CON(8)) GO TO 20
00167 15 TSTEPN = CON(8)
00168 GO TO 35
00169 20 DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
00170 25 TSTEPN = CON(18)-TSUM
00171 GO TO 35
00172 DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
00173 IF(TSUM+2.0*TSTEPN.LE.CON(18)) GO TO 35
00174 APPROACH THE OUTPUT INTERVAL GRADUALLY
00175 TSTEPN = (CON(18)-TSUM)/2.0
00176 STARE DELTA TIME STEP IN THE CONSTANTS
00177 CON(2) = TSTEPN
00178 CALCULATE THE NEW TIME
00179 IF(PASS.GT.0.) GO TO 40
00180 CON(1) = TPRINT
00181 CON(2) = 0.0
00182 GO TO 45
00183 40 CON(1) = TPRINT+TSUM+TSTEPN
00184 COMPUTE THE MEAN TIME BETWEEN ITERATIONS
00185 CON(14) = (CON(1)+CON(13))/2.0
00186 LAX = KON(5)
00187 DN = CON(10)
00188 DD = 1.0-DN
00189 AN = CON(9)
00190 AA = 1.0-AN
00191 TSTEP = TSTEPN/2.0
00192 DO THE RELAXATION LOOP
00193 DO 240 K1 = 1,LAX
00194 KON(20) = K1
00195 J1 = 0
00196 RLXA = 0.0
00197 RLXD = 0.0
00198 IF(K1.GT.1) GO TO 110
00199 J2 = 1
00200 ZERO OUT ALL SOURCE LOCATIONS AND SHIFT TEMPERATURES
00201 50 0(I) = 0.0
00202 DO 50 I = 1,MNC
00203 DO 55 I = 1,MNT
00204 LE1 = IE1+I
00205 X(LE1) = T(I)
00206 KON(12) = 0
00207 CALL VARBL1
00208 CHECK THE BACKUP SWITCH
00209 IF(KON(12).NE.0) GO TO 15
00210 CHECK FOR FIRST PASS
00211 IF(PASS.GE.0.) GO TO 60
00212 CALL OUTCAL
00213 PASS = 1.0
00214 89*
00215 88*
00216 87*
00217 86*
00218 85*
00219 84*
00220 83*
00221 82*
00222 81*
00223 80*
00224 79*
00225 78*
00226 77*
00227 76*
00228 75*
00229 74*
00230 73*
00231 72*
00232 71*
00233 70*
00234 69*
00235 68*
00236 67*
00237 66*
00238 65*
00239 64*
00240 63*
00241 62*
00242 61*
00243 60*
00244 59*
00245 58*
00246 57*
00247 56*
00248 55*
00249 54*
00250 53*
00251 52*
00252 51*
00253 50*
00254 49*
00255 48*
00256 47*
00257 46*
00258 45*
00259 44*
00260 43*
00261 42*
00262 41*
00263 40*
00264 39*
00265 38*
00266 37*
00267 36*
00268 35*
00269 34*
00270 33*
00271 32*
00272 31*
00273 30*
```

```

00255 90* GO TO 10
00256 91* RC = 1.E+8
00257 92* JJ = 0
00258 93* CALCULATE FIRST PASS TEMPERATURES AND CSGHIN
00259 94* DO 105 I = 1,NND
00260 95* INCLUDE VARQ,LIST
00261 96* FOLD DELTAT INTO THE CAPACITANCES
00262 97* IF(FLD(1,1,NSQ1(J1+1)),EQ.0) GO TO 2000
00263 98* NTYPE = FLD(0,5,NSQ2(J2))
00264 99* LA = FLD(5,17,NSQ2(J2))
00265 00* LK = FLD(22,14,NSQ2(J2))
00266 01* GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
00267 02* CALL D2DIWM(T(I),A(LA),XK(LK),C(I))
00268 03* GO TO 1999
00269 04* 1010 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00270 05* 1012 J2 = J2+1
00271 06* LA = FLD(5,17,NSQ2(J2))
00272 07* LK = FLD(22,14,NSQ2(J2))
00273 08* CALL D1DIWM(T(I),A(LA),XK(LK),C2)
00274 09* GO TO 1998
00275 10* 1015 C1 = XK(LK)*XK(LA)
00276 11* GO TO 1012
00277 12* 1020 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00278 13* J2 = J2+1
00279 14* LA = FLD(5,17,NSQ2(J2))
00280 15* LK = FLD(22,14,NSQ2(J2))
00281 16* C2 = XK(LK)*XK(LA)
00282 17* GO TO 1998
00283 18* 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
00284 19* GO TO 1999
00285 20* 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00286 21* 1032 J2 = J2+1
00287 22* LA = FLD(5,17,NSQ2(J2))
00288 23* LK = FLD(22,14,NSQ2(J2))
00289 24* CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00290 25* GO TO 1998
00291 26* 1035 C1 = XK(LK)*XK(LA)
00292 27* GO TO 1032
00293 28* 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00294 29* J2 = J2+1
00295 30* LA = FLD(5,17,NSQ2(J2))
00296 31* LK = FLD(22,14,NSQ2(J2))
00297 32* C2 = XK(LK)*XK(LA)
00298 33* GO TO 1948
00299 34* 1045 CALL D2DIWM(T(I),CON(14),A(LA),XK(LK),C(I))
00300 35* GO TO 1999
00301 36* 1990 C(I) = C1+C2
00302 37* 1999 J2 = J2+1
00303 38* 2000 CONTINUE
00304 39* END
00305 40* C(I) = C(I)/TSTEP
00306 41* R1 = 0.0
00307 42* S = 0.0
00308 43* G2 = 0.0
00309 44* INCLUDE VARQ,LIST
00310 45* IF(FLD(4,1,NSQ1(J1+1)),EQ.0) GO TO 5000
00311 46* NTYPE = FLD(0,5,NSQ2(J2))
00312 47* LA = FLD(5,17,NSQ2(J2))
00313 48* LK = FLD(22,14,NSQ2(J2))
00314 49* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4045), NTYPE
00315 50* CALL D2DIWM(T(I),A(LA),XK(LK),C(I))
00316 51* GO TO 1999
00317 52* 4005 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00318 53* 4010 J2 = J2+1
00319 54* 4015 LA = FLD(5,17,NSQ2(J2))
00320 55* 4020 LK = FLD(22,14,NSQ2(J2))
00321 56* 4025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00322 57* 4030 GO TO 1998
00323 58* 4035 C1 = XK(LK)*XK(LA)
00324 59* 4040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00325 60* 4045 J2 = J2+1
00326 61* 4050 LA = FLD(5,17,NSQ2(J2))
00327 62* 4055 LK = FLD(22,14,NSQ2(J2))
00328 63* 4060 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00329 64* 4065 GO TO 1998
00330 65* 4070 C1 = XK(LK)*XK(LA)
00331 66* 4075 GO TO 1072
00332 67* 4080 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00333 68* 4085 J2 = J2+1
00334 69* 4090 LA = FLD(5,17,NSQ2(J2))
00335 70* 4095 LK = FLD(22,14,NSQ2(J2))
00336 71* 4100 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00337 72* 4105 GO TO 1998
00338 73* 4110 C1 = XK(LK)*XK(LA)
00339 74* 4115 GO TO 1072
00340 75* 4120 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00341 76* 4125 J2 = J2+1
00342 77* 4130 LA = FLD(5,17,NSQ2(J2))
00343 78* 4135 LK = FLD(22,14,NSQ2(J2))
00344 79* 4140 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00345 80* 4145 GO TO 1998
00346 81* 4150 C1 = XK(LK)*XK(LA)
00347 82* 4155 GO TO 1072
00348 83* 4160 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00349 84* 4165 J2 = J2+1
00350 85* 4170 LA = FLD(5,17,NSQ2(J2))
00351 86* 4175 LK = FLD(22,14,NSQ2(J2))
00352 87* 4180 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00353 88* 4185 GO TO 1998
00354 89* 4190 C1 = XK(LK)*XK(LA)
00355 90* 4195 GO TO 1072
00356 91* 4200 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00357 92* 4205 J2 = J2+1
00358 93* 4210 LA = FLD(5,17,NSQ2(J2))
00359 94* 4215 LK = FLD(22,14,NSQ2(J2))
00360 95* 4220 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00361 96* 4225 GO TO 1998
00362 97* 4230 C1 = XK(LK)*XK(LA)
00363 98* 4235 GO TO 1072
00364 99* 4240 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00365 00* 4245 J2 = J2+1
00366 01* 4250 LA = FLD(5,17,NSQ2(J2))
00367 02* 4255 LK = FLD(22,14,NSQ2(J2))
00368 03* 4260 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00369 04* 4265 GO TO 1998
00370 05* 4270 C1 = XK(LK)*XK(LA)
00371 06* 4275 GO TO 1072
00372 07* 4280 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00373 08* 4285 J2 = J2+1
00374 09* 4290 LA = FLD(5,17,NSQ2(J2))
00375 10* 4295 LK = FLD(22,14,NSQ2(J2))
00376 11* 4300 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00377 12* 4305 GO TO 1998
00378 13* 4310 C1 = XK(LK)*XK(LA)
00379 14* 4315 GO TO 1072
00380 15* 4320 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00381 16* 4325 J2 = J2+1
00382 17* 4330 LA = FLD(5,17,NSQ2(J2))
00383 18* 4335 LK = FLD(22,14,NSQ2(J2))
00384 19* 4340 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00385 20* 4345 GO TO 1998
00386 21* 4350 C1 = XK(LK)*XK(LA)
00387 22* 4355 GO TO 1072
00388 23* 4360 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00389 24* 4365 J2 = J2+1
00390 25* 4370 LA = FLD(5,17,NSQ2(J2))
00391 26* 4375 LK = FLD(22,14,NSQ2(J2))
00392 27* 4380 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00393 28* 4385 GO TO 1998
00394 29* 4390 C1 = XK(LK)*XK(LA)
00395 30* 4395 GO TO 1072
00396 31* 4400 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00397 32* 4405 J2 = J2+1
00398 33* 4410 LA = FLD(5,17,NSQ2(J2))
00399 34* 4415 LK = FLD(22,14,NSQ2(J2))
00400 35* 4420 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00401 36* 4425 GO TO 1998
00402 37* 4430 C1 = XK(LK)*XK(LA)
00403 38* 4435 GO TO 1072
00404 39* 4440 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00405 40* 4445 J2 = J2+1
00406 41* 4450 LA = FLD(5,17,NSQ2(J2))
00407 42* 4455 LK = FLD(22,14,NSQ2(J2))
00408 43* 4460 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00409 44* 4465 GO TO 1998
00410 45* 4470 C1 = XK(LK)*XK(LA)
00411 46* 4475 GO TO 1072
00412 47* 4480 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00413 48* 4485 J2 = J2+1
00414 49* 4490 LA = FLD(5,17,NSQ2(J2))
00415 50* 4495 LK = FLD(22,14,NSQ2(J2))
00416 51* 4500 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00417 52* 4505 GO TO 1998
00418 53* 4510 C1 = XK(LK)*XK(LA)
00419 54* 4515 GO TO 1072
00420 55* 4520 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00421 56* 4525 J2 = J2+1
00422 57* 4530 LA = FLD(5,17,NSQ2(J2))
00423 58* 4535 LK = FLD(22,14,NSQ2(J2))
00424 59* 4540 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00425 60* 4545 GO TO 1998
00426 61* 4550 C1 = XK(LK)*XK(LA)
00427 62* 4555 GO TO 1072
00428 63* 4560 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00429 64* 4565 J2 = J2+1
00430 65* 4570 LA = FLD(5,17,NSQ2(J2))
00431 66* 4575 LK = FLD(22,14,NSQ2(J2))
00432 67* 4580 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00433 68* 4585 GO TO 1998
00434 69* 4590 C1 = XK(LK)*XK(LA)
00435 70* 4595 GO TO 1072
00436 71* 4600 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00437 72* 4605 J2 = J2+1
00438 73* 4610 LA = FLD(5,17,NSQ2(J2))
00439 74* 4615 LK = FLD(22,14,NSQ2(J2))
00440 75* 4620 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00441 76* 4625 GO TO 1998
00442 77* 4630 C1 = XK(LK)*XK(LA)
00443 78* 4635 GO TO 1072
00444 79* 4640 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00445 80* 4645 J2 = J2+1
00446 81* 4650 LA = FLD(5,17,NSQ2(J2))
00447 82* 4655 LK = FLD(22,14,NSQ2(J2))
00448 83* 4660 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00449 84* 4665 GO TO 1998
00450 85* 4670 C1 = XK(LK)*XK(LA)
00451 86* 4675 GO TO 1072
00452 87* 4680 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00453 88* 4685 J2 = J2+1
00454 89* 4690 LA = FLD(5,17,NSQ2(J2))
00455 90* 4695 LK = FLD(22,14,NSQ2(J2))
00456 91* 4700 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00457 92* 4705 GO TO 1998
00458 93* 4710 C1 = XK(LK)*XK(LA)
00459 94* 4715 GO TO 1072
00460 95* 4720 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00461 96* 4725 J2 = J2+1
00462 97* 4730 LA = FLD(5,17,NSQ2(J2))
00463 98* 4735 LK = FLD(22,14,NSQ2(J2))
00464 99* 4740 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00465 00* 4745 GO TO 1998
00466 01* 4750 C1 = XK(LK)*XK(LA)
00467 02* 4755 GO TO 1072
00468 03* 4760 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00469 04* 4765 J2 = J2+1
00470 05* 4770 LA = FLD(5,17,NSQ2(J2))
00471 06* 4775 LK = FLD(22,14,NSQ2(J2))
00472 07* 4780 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00473 08* 4785 GO TO 1998
00474 09* 4790 C1 = XK(LK)*XK(LA)
00475 10* 4795 GO TO 1072
00476 11* 4800 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00477 12* 4805 J2 = J2+1
00478 13* 4810 LA = FLD(5,17,NSQ2(J2))
00479 14* 4815 LK = FLD(22,14,NSQ2(J2))
00480 15* 4820 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00481 16* 4825 GO TO 1998
00482 17* 4830 C1 = XK(LK)*XK(LA)
00483 18* 4835 GO TO 1072
00484 19* 4840 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00485 20* 4845 J2 = J2+1
00486 21* 4850 LA = FLD(5,17,NSQ2(J2))
00487 22* 4855 LK = FLD(22,14,NSQ2(J2))
00488 23* 4860 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00489 24* 4865 GO TO 1998
00490 25* 4870 C1 = XK(LK)*XK(LA)
00491 26* 4875 GO TO 1072
00492 27* 4880 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00493 28* 4885 J2 = J2+1
00494 29* 4890 LA = FLD(5,17,NSQ2(J2))
00495 30* 4895 LK = FLD(22,14,NSQ2(J2))
00496 31* 4900 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00497 32* 4905 GO TO 1998
00498 33* 4910 C1 = XK(LK)*XK(LA)
00499 34* 4915 GO TO 1072
00500 35* 4920 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00501 36* 4925 J2 = J2+1
00502 37* 4930 LA = FLD(5,17,NSQ2(J2))
00503 38* 4935 LK = FLD(22,14,NSQ2(J2))
00504 39* 4940 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00505 40* 4945 GO TO 1998
00506 41* 4950 C1 = XK(LK)*XK(LA)
00507 42* 4955 GO TO 1072
00508 43* 4960 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00509 44* 4965 J2 = J2+1
00510 45* 4970 LA = FLD(5,17,NSQ2(J2))
00511 46* 4975 LK = FLD(22,14,NSQ2(J2))
00512 47* 4980 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00513 48* 4985 GO TO 1998
00514 49* 4990 C1 = XK(LK)*XK(LA)
00515 50* 4995 GO TO 1072
00516 51* 5000 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00517 52* 5005 J2 = J2+1
00518 53* 5010 LA = FLD(5,17,NSQ2(J2))
00519 54* 5015 LK = FLD(22,14,NSQ2(J2))
00520 55* 5020 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00521 56* 5025 GO TO 1998
00522 57* 5030 C1 = XK(LK)*XK(LA)
00523 58* 5035 GO TO 1072
00524 59* 5040 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00525 60* 5045 J2 = J2+1
00526 61* 5050 LA = FLD(5,17,NSQ2(J2))
00527 62* 5055 LK = FLD(22,14,NSQ2(J2))
00528 63* 5060 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00529 64* 5065 GO TO 1998
00530 65* 5070 C1 = XK(LK)*XK(LA)
00531 66* 5075 GO TO 1072
00532 67* 5080 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00533 68* 5085 J2 = J2+1
00534 69* 5090 LA = FLD(5,17,NSQ2(J2))
00535 70* 5095 LK = FLD(22,14,NSQ2(J2))
00536 71* 5100 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00537 72* 5105 GO TO 1998
00538 73* 5110 C1 = XK(LK)*XK(LA)
00539 74* 5115 GO TO 1072
00540 75* 5120 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00541 76* 5125 J2 = J2+1
00542 77* 5130 LA = FLD(5,17,NSQ2(J2))
00543 78* 5135 LK = FLD(22,14,NSQ2(J2))
00544 79* 5140 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00545 80* 5145 GO TO 1998
00546 81* 5150 C1 = XK(LK)*XK(LA)
00547 82* 5155 GO TO 1072
00548 83* 5160 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00549 84* 5165 J2 = J2+1
00550 85* 5170 LA = FLD(5,17,NSQ2(J2))
00551 86* 5175 LK = FLD(22,14,NSQ2(J2))
00552 87* 5180 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00553 88* 5185 GO TO 1998
00554 89* 5190 C1 = XK(LK)*XK(LA)
00555 90* 5195 GO TO 1072
00556 91* 5200 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00557 92* 5205 J2 = J2+1
00558 93* 5210 LA = FLD(5,17,NSQ2(J2))
00559 94* 5215 LK = FLD(22,14,NSQ2(J2))
00560 95* 5220 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00561 96* 5225 GO TO 1998
00562 97* 5230 C1 = XK(LK)*XK(LA)
00563 98* 5235 GO TO 1072
00564 99* 5240 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00565 00* 5245 J2 = J2+1
00566 01* 5250 LA = FLD(5,17,NSQ2(J2))
00567 02* 5255 LK = FLD(22,14,NSQ2(J2))
00568 03* 5260 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00569 04* 5265 GO TO 1998
00570 05* 5270 C1 = XK(LK)*XK(LA)
00571 06* 5275 GO TO 1072
00572 07* 5280 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00573 08* 5285 J2 = J2+1
00574 09* 5290 LA = FLD(5,17,NSQ2(J2))
00575 10* 5295 LK = FLD(22,14,NSQ2(J2))
00576 11* 5300 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00577 12* 5305 GO TO 1998
00578 13* 5310 C1 = XK(LK)*XK(LA)
00579 14* 5315 GO TO 1072
00580 15* 5320 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00581 16* 5325 J2 = J2+1
00582 17* 5330 LA = FLD(5,17,NSQ2(J2))
00583 18* 5335 LK = FLD(22,14,NSQ2(J2))
00584 19* 5340 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00585 20* 5345 GO TO 1998
00586 21* 5350 C1 = XK(LK)*XK(LA)
00587 22* 5355 GO TO 1072
00588 23* 5360 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00589 24* 5365 J2 = J2+1
00590 25* 5370 LA = FLD(5,17,NSQ2(J2))
00591 26* 5375 LK = FLD(22,14,NSQ2(J2))
00592 27* 5380 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00593 28* 5385 GO TO 1998
00594 29* 5390 C1 = XK(LK)*XK(LA)
00595 30* 5395 GO TO 1072
00596 31* 5400 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00597 32* 5405 J2 = J2+1
00598 33* 5410 LA = FLD(5,17,NSQ2(J2))
00599 34* 5415 LK = FLD(22,14,NSQ2(J2))
00600 35* 5420 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00601 36* 5425 GO TO 1998
00602 37* 5430 C1 = XK(LK)*XK(LA)
00603 38* 5435 GO TO 1072
00604 39* 5440 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00605 40* 5445 J2 = J2+1
00606 41* 5450 LA = FLD(5,17,NSQ2(J2))
00607 42* 5455 LK = FLD(22,14,NSQ2(J2))
00608 43* 5460 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00609 44* 5465 GO TO 1998
00610 45* 5470 C1 = XK(LK)*XK(LA)
00611 46* 5475 GO TO 1072
00612 47* 5480 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00613 48* 5485 J2 = J2+1
00614 49* 5490 LA = FLD(5,17,NSQ2(J2))
00615 50* 5495 LK = FLD(22,14,NSQ2(J2))
00616 51* 5500 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00617 52* 5505 GO TO 1998
00618 53* 5510 C1 = XK(LK)*XK(LA)
00619 54* 5515 GO TO 1072
00620 55* 5520 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00621 56* 5525 J2 = J2+1
00622 57* 5530 LA = FLD(5,17,NSQ2(J2))
00623 58* 5535 LK = FLD(22,14,NSQ2(J2))
00624 59* 5540 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00625 60* 5545 GO TO 1998
00626 61* 5550 C1 = XK(LK)*XK(LA)
00627 62* 5555 GO TO 1072
00628 63* 5560 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00629 64* 5565 J2 = J2+1
00630 65* 5570 LA = FLD(5,17,NSQ2(J2))
00631 66* 5575 LK = FLD(22,14,NSQ2(J2))
00632 67* 5580 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00633 68* 5585 GO TO 1998
00634 69* 5590 C1 = XK(LK)*XK(LA)
00635 70* 5595 GO TO 1072
00636 71* 5600 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00637 72* 5605 J2 = J2+1
00638 73* 5610 LA = FLD(5,17,NSQ2(J2))
00639 74* 5615 LK = FLD(22,14,NSQ2(J2))
00640 75* 5620 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00641 76* 5625 GO TO 1998
00642 77* 5630 C1 = XK(LK)*XK(LA)
00643 78* 5635 GO TO 1072
00644 79* 5640 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00645 80* 5645 J2 = J2+1
00646 81* 5650 LA = FLD(5,17,NSQ2(J2))
00647 82* 5655 LK = FLD(22,14,NSQ2(J2))
00648 83* 5660 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00649 84* 5665 GO TO 1998
00650 85* 5670 C1 = XK(LK)*XK(LA)
00651 86* 5675 GO TO 1072
00652 87* 5680 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00653 88* 5685 J2 = J2+1
00654 89* 5690 LA = FLD(5,17,NSQ2(J2))
00655 90* 5695 LK = FLD(22,14,NSQ2(J2))
00656 91* 5700 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00657 92* 5705 GO TO 1998
00658 93* 5710 C1 = XK(LK)*XK(LA)
00659 94* 5715 GO TO 1072
00660 95* 5720 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00661 96* 5725 J2 = J2+1
00662 97* 5730 LA = FLD(5,17,NSQ2(J2))
00663 98* 5735 LK = FLD(22,14,NSQ2(J2))
00664 99* 5740 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00665 00* 5745 GO TO 1998
00666 01* 5750 C1 = XK(LK)*XK(LA)
00667 02* 5755 GO TO 1072
00668 03* 5760 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00669 04* 5765 J2 = J2+1
00670 05* 5770 LA = FLD(5,17,NSQ2(J2))
00671 06* 5775 LK = FLD(22,14,NSQ2(J2))
00672 07* 5780 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00673 08* 5785 GO TO 1998
00674 09* 5790 C1 = XK(LK)*XK(LA)
00675 10* 5795 GO TO 1072
00676 11* 5800 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00677 12* 5805 J2 = J2+1
00678 13* 5810 LA = FLD(5,17,NSQ2(J2))
00679 14* 5815 LK = FLD(22,14,NSQ2(J2))
00680 15* 5820 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00681 16* 5825 GO TO 1998
00682 17* 5830 C1 = XK(LK)*XK(LA)
00683 18* 5835 GO TO 1072
00684 19* 5840 CALL D1DIWM(T(I),A(LA),XK(LK),C1)
00685 20* 58
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CNF*WBK,CNF*WBK

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00353 101* 4005 Q(I) = XK(LK)+Q(I)
00354 101* GO TO 4999
00355 101* 4010 Q1 = 0.0
00356 101* 4012 CALL DID1WM(T(I),A(LA),XK(LK),Q2)
00357 101* GO TO 4998
00360 101* 4015 Q1 = 0.0
00361 101* 4017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
00362 101* GO TO 4998
00363 101* 4020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00364 101* 4022 J2 = J2+1
00365 101* LA = FLD(5,17,NSQ2(J2))
00366 101* LK = FLD(22,14,NSQ2(J2))
00367 101* GO TO 4017
00370 101* 4025 Q1 = XK(LK)*XK(LA)
00371 101* GO TO 4022
00372 101* 4030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00373 101* J2 = J2+1
00374 101* LA = FLD(5,17,NSQ2(J2))
00375 101* LK = FLD(22,14,NSQ2(J2))
00376 101* Q2 = XK(LK)*XK(LA)
00377 101* GO TO 4998
00400 101* 4035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00401 101* 4037 J2 = J2+1
00402 101* LA = FLD(5,17,NSQ2(J2))
00403 101* LK = FLD(22,14,NSQ2(J2))
00404 101* GO TO 4012
00405 101* 4040 Q1 = XK(LK)*XK(LA)
00406 101* GO TO 4037
00407 101* 4998 Q(I) = Q1+Q2+Q(I)
00410 101* 4999 J2 = J2+1
00411 101* 5000 CONTINUE
00412 101* END
00413 102* Q(I) = 2.0*Q(I)+C(I)*(T(I)+460.0)
00414 103* GSUM = G(I)
00415 104* GSUM = C(I)
00416 105* 70 J1 = J1+1
00417 106* LG = FLD(5,16,NSQ1(J1))
00420 107* LTA = FLD(22,14,NSQ1(J1))
00421 108* LTAE = LTA+IE1
00422 109* INCLUDE VARG,LIST
00423 109* IF(FLD(2,1,NSQ1(J1)),EQ,0) GO TO 3000
00425 109* NTYPE = FLD(0,5,NSQ2(J2))
00426 109* LA = FLD(5,17,NSQ2(J2))
00427 109* LK = FLD(22,14,NSQ2(J2))
00430 109* 60TO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00431 109* 2060,2065), NTYPE
00432 109* 2005 TM = (T(I)+T(LTA))/2.0
00433 109* 2007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
00434 109* GO TO 2999
00435 109* 2010 TM = T(I)
00436 109* 2015 CALL DID1WM(T(I),A(LA),XK(LK),G1)
00437 109* 2017 J2 = J2+1
00440 109* LA = FLD(5,17,NSQ2(J2))
00441 109* LK = FLD(22,14,NSQ2(J2))
00442 109* CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
00443 109* GO TO 2998
00445 109* 2020 G1 = XK(LK)*XK(LA)
00446 109* GO TO 2017
00447 109* 2025 CALL DID1WM(T(I),A(LA),XK(LK),G1)

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*NEW
*NEW
**-2

*NEW
*NEW
**-2

*NEW
*NEW
**-2

*NEW
*NEW
**-3

*NEW
*NEW
**-2

*NEW
*NEW
**=2

*NEW
*NEW
**=2

*NEW
*NEW
**=2

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00447 109* J2 = J2+1
00450 109* LA = FLD(15,17,NSQ2(J2))
00451 109* LK = FLD(22,14,NSQ2(J2))
00452 109* G2 = XK(LK)*XK(LA)
00453 109* GO TO 2998
00454 109* 2030 TM = (T(I)+T(LTA))/2.0
00455 109* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00456 109* GO TO 2999
00457 109* 2035 TM = T(I)
00460 109* GO TO 2032
00461 109* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00462 109* 2042 J2 = J2+1
00463 109* LA = FLD(15,17,NSQ2(J2))
00464 109* LK = FLD(22,14,NSQ2(J2))
00465 109* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00466 109* GO TO 2998
00467 109* 2045 G1 = XK(LK)*XK(LA)
00470 109* GO TO 2042
00471 109* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00472 109* J2 = J2+1
00473 109* LA = FLD(15,17,NSQ2(J2))
00474 109* LK = FLD(22,14,NSQ2(J2))
00475 109* G2 = XK(LK)*XK(LA)
00476 109* GO TO 2998
00477 109* 2055 TM = (T(I)+T(LTA))/2.0
00500 109* CALL D2D1WM(TM,CON(14),A(LA),XK(LK),G(LG))
00501 109* GO TO 2999
00502 109* 2060 TM = T(LTA)
00503 109* GO TO 2007
00504 109* 2065 TM = T(LTA)
00505 109* GO TO 2032
00506 109* 2998 G(LG) = 1./(1./G1+1./G2)
00507 109* IF(FLD(3,1,NSQ1(J1)).EQ.1) G(LG) = G1*G2
00511 109* 2999 J2 = J2+1
00512 109* 3000 CONTINUE
00513 109* END
00514 110* T1 = T(I)+460.0
00515 111* T2 = T(LTA)+460.0
00515 112* CHECK FOR RADIATION CONDUCTOR
00516 113* IF(FLD(3,1,NSQ1(J1)).EQ.0) GO TO 75
00520 114* R1 = R1+G(LG)
00521 115* OSUM = OSUM+G(LG)*T2**4
00522 116* G2 = G2+G(LG)*(T1*T1+T2*T2)*(T1+T2)
00523 117* Q(I) = Q(I)+G(LG)*(X(LTAE)+460.0)**4-T1**4
00524 118* GO TO 80
00525 119* 75 GV = G(LG)
00526 120* Q(I) = Q(I)+GV*(X(LTAE)-T(I))
00527 121* G2 = G2+GV
00530 122* OSUM = OSUM+GV
00531 123* OSUM = OSUM+GV*T2
00531 124* CHECK FOR LAST CONDUCTOR
00532 125* IF(NSQ1(J1).GT.0) GO TO 70
00532 126* DAMPEN RADIATION ON THIS NODE IF PRESENT
00534 127* IF(R1.LE.0.) GO TO 100
00536 128* R2 = R1+T1**4
00537 129* T2 = (OSUM-R2)/OSUM
00540 130* R1 = R1+T2**4
00541 131* S = (R1+R2)/2.0
00541 132* OBTAIN THE NEW TEMPERATURE
00542 133* T(I) = (DN*((OSUM-S)/OSUM)+DD+T1)-460.0

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CNFWBK*CNFWRK

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00543 134* R1 = C(I)/62
00544 135* IF(R1.GE.RC) GO TO 105
00546 136* RC = R1
00547 137* KON(35) = I
00550 138* C 105 CONTINUE
00550 139* CONVERT TEMPERATURES TO RANKINE
00552 140* DO 65 I = 1,NNT
00555 141* LE1 = IE1+I
00556 142* T(I) = T(I)+60.
00557 143* 65 X(LE1) = X(LE1)+460.
00561 144* CON(17) = RC*TSTEP
00562 145* IF(RC.LE.0.) GO TO 993
00564 146* GO TO 225
C 110 NOW RELAX THE NETWORK BY SUCCESSIVE POINT AND EXTRAPOLATION
00564 147* JJ = JJ+1
00565 148* DO 165 I = 1,NND
00566 149* R1 = 0.0
00571 150* S = 0.0
00572 151* OSUM = 0(I)
00573 152* GSUM = C(I)
00574 153* 115 J1 = J1+1
00575 154* LG = FLD(5,16,NSQ1(J1))
00576 155* LTA = FLD(22,14,NSQ1(J1))
00577 156* CHECK FOR RADIATION CONDUCTOR
00600 158* IF(FLD(3,1,NSQ1(J1)).EQ.0) GO TO 120
00602 159* R1 = R1+G(LG)
00603 160* OSUM = OSUM+G(LG)*T(LTA)**4
00604 161* GO TO 125
00605 162* 120 GSUM = GSUM+G(LG)
00606 163* OSUM = OSUM+G(LG)*T(LTA)
00607 164* CHECK FOR LAST CONDUCTOR
00607 165* C 125 IF(NSQ1(J1).GT.0) GO TO 115
00607 166* DAMPEN RADIATION ON THIS NODE IF PRESENT
00611 167* IF(R1.LE.0.) GO TO 145
00613 168* R2 = R1*T(I)**4
00614 169* T2 = (OSUM-R2)/GSUM
00615 170* R1 = R1*T2**4
00616 171* S = (R1+R2)/2.0
00616 172* C 145 OBTAIN THE NEW TEMPERATURE
00617 173* T2 = DN*((OSUM-S)/GSUM)+DD*T(I)
00617 174* C OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
00620 175* T1 = ABS(T(I)-T2)
00621 176* C STORE THE NEW AND OLD TEMPERATURES
00621 177* GO TO (160,155,150), JJ
00622 178* 150 LE2 = IE2+I
00623 179* LE3 = IE3+I
00624 180* R1 = T2-T(I)
00625 181* X(LE2) = T(I)
00626 182* X(LE3) = R1/(R1-X(LE3))
00627 183* GO TO 160
00630 184* 155 LE3 = IE3+I
00631 185* X(LE3) = T2-T(I)
00632 186* T(I) = T2
00633 187* 160 IF(RLX(16E,11) GO TO 165
00635 188* RLXD = T1
00636 189* KK1 = 1
00637 190* 165 CONTINUE
00641 191* GO TO (180,180,170), JJ
00641 192* C PERFORM LINEAR EXTRAPOLATION ON THE ERROR FUNCTION CURVE
00642 193* JJ = 0

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CNFWBK,CNFWBK

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00643 194* DO 175 I = 1,NND
00646 195* LE3 = IE3+1
00646 196* SEE IF THE EXTRAPOLATION IS ALLOWABLE
00647 197* IF(X(LE3).GE.0.) GO TO 175
00647 198* LIMIT THE EXTRAPOLATION
00651 199* IF(X(LE3).LT.-10.) X(LE3) = -10.
00653 200* LE2 = IE2+1
00654 201* T(I) = X(LE3)*X(LE2)+(1.0-X(LE3))*T(I)
00655 202* 175 CONTINUE
00657 203* 180 IF(NNA.LE.0) GO TO 220
00661 204* JJ1 = J1
00662 205* JJ2 = J2
00663 206* DO 230 I = 1,NNT
00666 207* T(I) = T(I)-460.0
00670 208* DO 215 I = NN,MNC
00673 209* L = 1
00674 210* IF(KI.GT.2) GO TO 6000
00676 211* INCLUDE VR02,LIST
00677 211* IF(FLD(4,1,NS01(JJ1+1),EQ.0) GO TO 6000
00701 211* NTYPE = FLD(0,5,NS02(JJ2))
00702 211* LA = FLD(5,17,NS02(JJ2))
00703 211* LK = FLD(22,14,NS02(JJ2))
00704 211* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
00705 211* 5005 0(L) = XK(LK)*0(L)
00706 211* GO TO 5999
00707 211* 5010 01 = 0.0
00710 211* 5012 CALL DID1WM(T(L),A(LA),XK(LK),02)
00711 211* GO TO 5998
00712 211* 5015 01 = 0.0
00713 211* 5017 CALL DID1WM(CON(14),A(LA),XK(LK),02)
00714 211* GO TO 5998
00715 211* 5020 CALL DIO1WM(CON(14),A(LA),XK(LK),01)
00716 211* 5022 JJ2 = JJ2+1
00720 211* LA = FLD(5,17,NS02(JJ2))
00721 211* LK = FLD(22,14,NS02(JJ2))
00722 211* GO TO 5017
00723 211* 5025 01 = XK(LK)*XK(LA)
00724 211* GO TO 5022
00725 211* 5030 CALL DID1WM(CON(14),A(LA),XK(LK),01)
00726 211* JJ2 = JJ2+1
00727 211* LA = FLD(5,17,NS02(JJ2))
00730 211* LK = FLD(22,14,NS02(JJ2))
00731 211* 02 = XK(LK)*XK(LA)
00732 211* GO TO 5998
00733 211* 5035 CALL DID1WM(CON(14),A(LA),XK(LK),01)
00734 211* 5037 JJ2 = JJ2+1
00735 211* LA = FLD(5,17,NS02(JJ2))
00736 211* LK = FLD(22,14,NS02(JJ2))
00737 211* GO TO 5012
00740 211* 5040 01 = XK(LK)*XK(LA)
00741 211* GO TO 5037
00742 211* 5998 0(L) = 01+02+0(L)
00743 211* 5999 JJ2 = JJ2+1
00744 211* 6000 CONTINUE
00745 212* EHD
00746 213* GSUM = 0.0
00747 214* OSUM = 0(I)
00750 215* JJ1 = JJ1+1
00751 216* L6 = FLD(5,16,NS01(JJ1))
LTA = FLD(22,14,NS01(JJ1))

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*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

CNFWBK,CNFWRK

00752	217*	IF(K1,GT.2) GO TO 4000			
00754	218*	INCLUDE VRG2,LIST			
00755	218*	IF(FLD(2,1,NS01(JJ1)).EQ.0) GO TO 4000			*NEW
00757	218*	NTPY = FLD(0,5,NS02(JJ2))			*NEW
00760	218*	LA = FLD(5,17,NS02(JJ2))			*NEW
00761	218*	LK = FLD(22,14,NS02(JJ2))			**--3
00762	218*	GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,			
00763	218*	3060,3065), NTPY			
00764	218*	TM = (T(L)+T(LTA))/2.0			
00765	218*	CALL DIDIWM(TM,A(LA),XK(LK),G(LG))			
00766	218*	GO TO 3999			
00767	218*	TM = T(L)			
00770	218*	GO TO 3007			
00771	218*	CALL DIDIWM(T(L),A(LA),XK(LK),G1)			
00772	218*	LA = FLD(5,17,NS02(JJ2))			*NEW
00773	218*	LK = FLD(22,14,NS02(JJ2))			*NEW
00774	218*	CALL DIDIWM(T(LTA),A(LA),XK(LK),G2)			**--2
00775	218*	GO TO 3998			
00776	218*	G1 = XK(LK)*XK(LA)			
00777	218*	GO TO 3017			
01000	218*	CALL DIDIWM(T(L),A(LA),XK(LK),G1)			
01001	218*	JJ2 = JJ2+1			
01002	218*	LA = FLD(5,17,NS02(JJ2))			*NEW
01003	218*	LK = FLD(22,14,NS02(JJ2))			*NEW
01004	218*	G2 = XK(LK)*XK(LA)			**--2
01005	218*	GO TO 3998			
01006	218*	TM = (T(L)+T(LTA))/2.0			
01007	218*	CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))			
01010	218*	GO TO 3999			
01011	218*	TM = T(L)			
01012	218*	GO TO 3032			
01013	218*	CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)			
01014	218*	LA = FLD(5,17,NS02(JJ2))			*NEW
01015	218*	LK = FLD(22,14,NS02(JJ2))			*NEW
01016	218*	CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)			**--2
01017	218*	GO TO 3998			
01020	218*	G1 = XK(LK)*XK(LA)			
01021	218*	GO TO 3042			
01022	218*	CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)			
01023	218*	JJ2 = JJ2+1			
01024	218*	LA = FLD(5,17,NS02(JJ2))			*NEW
01025	218*	LK = FLD(22,14,NS02(JJ2))			*NEW
01026	218*	G2 = XK(LK)*XK(LA)			**--2
01027	218*	GO TO 3998			
01030	218*	TM = (T(L)+T(LTA))/2.0			
01031	218*	CALL DIDIWM(TM,CON(14),A(LA),XK(LK),G(LG))			
01032	218*	GO TO 3999			
01033	218*	TM = T(LTA)			
01034	218*	GO TO 3007			
01035	218*	TM = T(LTA)			
01036	218*	GO TO 3032			
01037	218*	G(LG) = 1./(1/G1+1/G2)			
01040	218*	IF(FLD(3,1,NS01(JJ1)).EQ.1) G(LG) = G1*G2			
01041	218*	JJ2 = JJ2+1			
01043	218*	4000 CONTINUE			
01044	218*	END			
01045	218*	T1 = T(L)+460.0			
01046	219*	T2 = T(LTA)+460.0			
01047	220*				

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01047 221* CHECK FOR RADIATION CONDUCTOR
01050 222* IF(FLD(3,1,NSQ1(JJ1)).EQ.0) GO TO 190
01052 223* GV = G(LG)*(T1+T2)*T1+T2*(T1+T2)
01053 224* GO TO 195
01054 225* 190 GV = G(LG)
01055 226* 195 GSUM = GSUM+GV
01056 227* GSUM = GSUM+GV*T2
01057 228* CHECK FOR LAST CONDUCTOR
01057 229* IF(NSQ1(JJ1).GT.0) GO TO 185
01057 230* CALCULATE THE NEW TEMPERATURE
01061 231* T2 = AN*GSUM/GSUM+AA*T1
01062 232* T1 = ABS(T2-T1)
01063 233* T(I) = T2-460.0
01064 234* IF(RLXA.GE.T1) GO TO 215
01066 235* RLXA = T1
01067 236* KK2 = I
01070 237* 215 CONTINUE
01072 238* DO 235 I = 1,NNT
01075 239* T(I) = T(I)+60.0
01075 240* 235 SEE IF THE ARITHMETIC RELAXATION CRITERIA WAS MET
01077 241* IF(RLXA.GT.CON(19)) GO TO 225
01077 242* SEE IF THE DIFFUSION RELAXATION CRITERIA WAS MET
01101 243* 220 IF(RLXD.LE.CON(26)) GO TO 245
01103 244* 225 IF(KON(17).EQ.0) GO TO 240
01105 245* CALL OUTCAL
01106 246* 240 CONTINUE
01110 247* IF(KON(28).GE.65) CALL TOPLIN
01112 248* WRITE(G*882)
01114 249* KON(28) = KON(28)+2
01114 250* SEE IF THE TEMPERATURE CHANGES WERE TOO LARGE
01115 251* 245 TCGD = 0.0
01116 252* TCGA = 0.0
01117 253* DO 250 I = 1,N1D
01122 254* LE = IEI+I
01123 255* C(I) = C(I)*TSTEP
01124 256* T1 = ABS(T(I)-X(LE))
01125 257* IF(TCGD.GT.T1) GO TO 250
01127 258* TCGD = T1
01130 259* KON(36) = I
01131 260* 250 CONTINUE
01133 261* IF(TCGD.LE.CON(6)) GO TO 265
01135 262* TSTEPN = 0.95*TSTEPN+CON(6)/TCGD
01136 263* 255 DO 260 I = 1,NNT
01141 264* LE = IEI+I
01142 265* 260 T(I) = X(LE)-460.0
01144 266* GO TO 30
01145 267* 265 IF(MNA.LE.0) GO TO 275
01147 268* DO 270 I = NN,NNC
01152 269* LE = IEI+I
01153 270* T1 = ABS(T(I)-X(LE))
01154 271* IF(TCGA.GT.T1) GO TO 270
01156 272* TCGA = T1
01157 273* KON(38) = I
01160 274* 270 CONTINUE
01162 275* IF(TCGA.LE.CON(11)) GO TO 275
01164 276* TSTEPN = 0.95*TSTEPN+CON(11)/TCGA
01165 277* GO TO 255
01165 278* C CONVERT TEMPERATURES BACK TO FARENHEIT
01166 279* DO 280 I = 1,NNT
01171 280* 280 T(I) = T(I)-460.0

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CNF.WBK.CNF.WBK

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01171 281*
01172 282*
01173 283*
01174 284*
01175 284*
01176 285*
01200 286*
01201 287*
01202 288*
01203 289*
01204 290*
01205 291*
01205 292*
01205 293*
01206 294*
01210 295*
01211 296*
01211 297*
01212 298*
01212 299*
01214 300*
01216 301*
01216 302*
01217 303*
01220 304*
01220 305*
01221 306*
01223 307*
01225 308*
01226 310*
01230 311*
01231 312*
01233 313*
01234 314*
01236 315*
01237 316*
01242 317*
01243 318*
01245 319*
01246 320*
01250 321*
01251 322*
01253 323*
01254 324*
01256 325*
01257 326*
01261 327*
01263 328*
01263 329*
01264 330*
01265 331*
01266 332*
01267 333*
01270 334*
01271 335*
01272 336*
01273 337*
01274 338*
01275 339*

C STORE THE TEMPERATURE AND RELAXATION CHANGES
CON(15) = TCGD
CON(16) = TCGA
CON(27) = RLXD
IF(RLXA.GT.RLXD) GO TO 285
KK2 = KK1
RLXA = RLXD
285 KON(37) = KK2
CON(30) = RLXA
KON(12) = 0
CALL VARBL2
CHECK THE RACKUP SWITCH
IF(KON(12).NE.0) GO TO 255
ADVANCE TIME
CON(13) = CON(1)
TSUM = TSUM+TSTEPN
CHECK FOR TIME TO PRINT
IF(TSUM.GE.CON(18)) GO TO 290
CHECK FOR PRINT EVERY ITERATION
IF(KON(7).NE.0) CALL OUTCAL
GO TO 10
C TRY TO EVEN THE OUTPUT INTERVALS
TPRINT = TPRINT+TSUM
CALL OUTCAL
C IS TIME GREATER THAN END COMPUTE TIME
IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
NTH = IE1
NDIM = NLA
RETURN
990 WRITE(6,880)
GO TO 1000
991 WRITE(6,881)
GO TO 1000
993 WRITE(6,883)
GO TO 1000
994 WRITE(6,884) NDIM
GO TO 1000
995 WRITE(6,885)
GO TO 1000
996 WRITE(6,886)
GO TO 1000
997 WRITE(6,887)
GO TO 1000
998 WRITE(6,888)
GO TO 1000
999 WRITE(6,889)
1000 CALL OUTCAL
CALL EXIT
880 FORMAT(29H TRANSIENT TIME NOT SPECIFIED)
881 FORMAT(45H CNF.WBK REQUIRES LONG PSEUDO-COMPUTE SEQUENCE)
882 FORMAT(28H RELAXATION CRITERIA NOT MET)
883 FORMAT(24H CSGMIN ZERO OR NEGATIVE)
884 FORMAT(18,20H LOCATIONS AVAILABLE)
885 FORMAT(10H NO DRLXCA)
886 FORMAT(10H NO DTIMEI)
887 FORMAT(10H NO APLXCA)
888 FORMAT(19H NO OUTPUT INTERVAL)
889 FORMAT(19H NO NLOOP)
END
```

QIM FOR** CNVARB,CNVARB
UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-100 CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:51

SUBROUTINE CNVARB ENTRY POINT 004272

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	004311
0000	*CONST+TEMP	000131
0002	*SIMPLE VAR	000072
0004	*ARRAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	CONU	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARB1
0023	OUTCAL
0024	D10IWM
0025	PLYAWM
0026	D2DIWM
0027	TOPLIN
0030	VAR:L2
0031	EXIT
0032	NERN2\$
0033	NWOU\$
0034	NIO2\$
0035	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0002	R	000017	AA
0010	R	000000	C	0017	R	000000	CON
0002	R	000014	DN	0012	R	000000	G
0002	I	000042	G2	0006	R	000000	H
0002	I	000006	IE3	0002	I	000007	J
0002	I	000022	J1	0002	I	000025	J2
0017	I	000000	KON	0002	I	000021	K1
0002	I	000071	LE	0002	I	000027	LE1
0002	I	000034	LA	0021	I	000006	LS01
0021	I	000004	MCY	0020	I	000000	MDIM
0021	I	000001	NIA	0002	I	000003	NMC
0014	I	000000	NS02	0020	I	000001	NTH
0002	R	000053	BETA0	0002	R	000016	AN
0002	R	000015	DD	0002	R	000035	C1
0002	R	000046	G1	0002	R	000057	GSUM
0002	I	000005	IE2	0002	I	000026	I
0002	I	000054	JJ2	0002	I	000031	JJ
0002	I	000056	KK2	0015	I	000000	K
0002	I	000013	LG	0002	I	000065	L
0021	I	000005	MAT	0002	I	000037	LC2
0002	I	000002	NN	0021	I	000007	LS02
0013	I	000000	NS01	0021	I	000003	NGT
0002	R	000000	PASS	0020	I	000002	NNT
				0020	I	000002	NX
				0002	R	000054	BETAN
				0002	R	000036	C2
				0002	R	000047	GV
				0002	I	000004	IE1
				0002	I	000063	JJ1
				0002	I	000062	KK1
				0002	I	000033	LA
				0002	I	000061	LE3
				0021	I	000044	LTA
				0002	I	000001	NLA
				0021	I	000002	NNT
				0020	I	000002	NX

CVARB,CNVARB

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0011 R 000000 Q      0002 R 000056 QSUM      0002 R 000040 Q1      0002 R 000041 Q2
0012 R 000023 RLXA   0002 R 000024 RLXD      0002 R 000052 R1      0002 R 000060 R2
0017 R 000000 T      0002 R 000070 TCGA      0002 R 000067 TCGD      0002 R 000045 TM
0020 R 000020 TSTEP  0002 R 000012 TSTEPN      0002 R 000011 T5UM      0002 R 000050 T1
0020 R 000002 X      0015 R 000000 XK      0001 000162 10L      0001 002075 100L
0001 000427 1005L   0001 000450 1010L   0001 000466 1012L   0001 000521 1015L
0001 000567 1025L   0001 000613 1030L   0001 000534 1032L   0001 000672 1035L
0001 000743 1045L   0001 002150 105L    0001 003655 1073G   0001 002201 110L
0001 003773 1137G   0001 002205 115L    0001 004016 1150G   0001 004061 1167G
0001 002256 125L    0001 002325 145L    0001 002171 15L     0001 002351 150L
0001 002403 160L   0001 002416 165L    0001 002431 170L   0001 002500 175L
0001 003021 185L   0001 003600 190L    0001 003603 195L   0001 000770 199L
0001 000174 20L    0001 000773 2000L   0001 001346 2005L   0001 001353 2007L
0001 001376 2015L  0001 001414 2017L   0001 001450 2020L   0001 001456 2025L
0001 001523 2032L  0001 001546 2035L   0001 001451 2040L   0001 001572 2042L
0001 001637 2050L  0001 001702 2055L   0001 001730 2060L   0001 001734 2065L
0001 003666 220L   0001 000274 221G   0001 003673 225L   0001 000316 233G
0001 003677 240L   0001 003720 245L   0001 003753 250L   0001 003767 255L
0001 004006 265L   0001 004040 270L   0001 004055 275L   0001 004104 285L
0001 001740 2998L   0001 001766 2999L   0001 002006 30L    0001 001771 3000L
0001 003115 3007L  0001 003135 3010L   0001 003141 3015L   0001 003160 3017L
0001 003222 3025L   0001 003263 3030L   0001 003271 3032L   0001 003314 3035L
0001 003342 3042L  0001 003401 3045L   0001 003407 3050L   0001 003453 3055L
0001 003506 3065L  0001 000221 35L    0001 003512 3998L   0001 003540 3999L
0001 003543 4000L  0001 001046 4005L   0001 001053 4010L   0001 001054 4012L
0001 001074 4017L  0001 001111 4020L   0001 001125 4022L   0001 001142 4025L
0001 001206 4035L  0001 001222 4037L   0001 001237 4040L   0001 001237 45L
0001 001251 4999L  0001 000150 5L    0001 001254 5000L   0001 002604 5005L
0001 002613 5012L  0001 002633 5015L   0001 002634 5017L   0001 002651 5020L
0001 002702 5025L  0001 002710 5030L   0001 002746 5035L   0001 002762 5037L
0001 002132 544G   0001 002166 555G   0001 003005 5998L   0001 003012 5999L
0001 003015 6000L  0001 002441 644G   0001 002525 664G   0001 002533 671G
0001 002041 75L    0001 002052 80L   0001 000300 800F   0000 000006 886F
0000 000025 833F   0000 000032 843F   0000 000037 895F   0000 000042 886F
0000 000050 888F   0000 000054 895F   0001 004156 990L   0001 004164 991L
0001 004200 994L   0001 004207 995L   0001 004215 996L   0001 004223 997L

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00101 1* SUBROUTINE CNVARB
00101 2* IMPLICIT FORWARD-BACKWARD DIFFERENCING EXECUTION SUBROUTINE
00101 3* C WITH AN INTERNALLY CALCULATED BETA WEIGHTING FACTOR
00101 4* C THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED, SINDA FORTRAN V
00101 5* C ALL NODES RECEIVE A SUCCESSIVE POINT ITERATION
00101 6* C RELAXATION CRITERIA MUST BE SPECIFIED
00101 7* C OVER-RELAXATION IS ALLOWED, THE DAMPENING FACTORS ARE ADDRESSABLE
00101 8* INCLUDE COMM,LIST
00103 8* COMMON /TITLE/HH(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
00104 8* COMMON /PCI/NS01(1) /PC2/NS02(1) /KONST/K(1) /ARRAY/A(1)
00105 8* COMMON /FIXCON/KON(1) /XSPACE/NDIM/NTHX(1)
00106 8* COMMON /DIMENS/ NND,NNA,NNT,NGT,NCT,NAT,LS01,LS02
00110 8* DIMENSION CON(1),XK(1),NX(1)
00111 8* EQUIVALENCE (KON(1),CON(1)),K(1),XK(1)),X(1),RX(1)
00112 8* END
00113 9* INCLUDE DEFF,LIST
00113 9* C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00113 9* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
00113 9* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTINEU)

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00113 9* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
00113 9* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR EXPLICIT (CSGFAC)
00113 9* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00113 9* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00113 9* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPCTR)
00113 9* C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
00113 9* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
00113 9* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
00113 9* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
00113 9* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
00113 9* C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEH)
00113 9* C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
00113 9* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DIMPCC)
00113 9* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
00113 9* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSGMIN)
00113 9* C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPT)
00113 9* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 9* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (DLOOPCT)
00113 9* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
00113 9* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (CSGMAX)
00113 9* C CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
00113 9* C CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
00113 9* C CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
00113 9* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 9* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 9* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 9* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 9* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 9* C CC31 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 9* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGBAL)
00113 9* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 9* C CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS (NOCOPY)
00113 9* C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 9* C CC36 CONTAINS RELATIVE NODE NUMBER OF DIMPCC
00113 9* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 9* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 9* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 9* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 9* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 9* C CC50 IS NOT USED AT PRESENT
00114 9* C END
00115 10* IF(KON(5),LE.0) GO TO 999
00117 11* IF(CON(6),LE.0) CON(6) = 1.E+8
00121 12* IF(CON(8),LE.0) CON(8) = 1.E+8
00123 13* IF(CON(9),LE.0) CON(9) = 1.0
00125 14* IF(CON(10),LE.0) CON(10) = 1.0
00127 15* IF(CON(11),LE.0) CON(11) = 1.E+8
00131 16* IF(CON(3),LE.0) CON(13) GO TO 990
00133 17* IF(CON(19),LE.0) GO TO 998
00135 18* IF(NNIA.GT.0.AND.CON(19),LE.0) GO TO 997
00137 19* IF(CON(22),LE.0) GO TO 996
00141 20* IF(NND.GT.0.AND.CON(26),LE.0) GO TO 995
00143 21* IF(KON(31),NE.1) GO TO 991
00145 22* PASS = -1.0
00146 23* NLA = NDIM
00147 24* NLA = NND+1
00150 25* NRC = NND+MNA
00151 26* IE1 = NTH
00152 27* IE2 = NTH+HINT
00153 28* IE3 = IE2+NND

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CNVARB,CNVARB

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00154 29* J = 2*NND+NNI
00155 30* NTH = NTH+J
00156 31* NDIM = NDIM-J
00157 32* CHECK FOR EXTRA LOCATIONS FOR CALCULATED NODES
00158 33* IF(NDIM.LT.0) GO TO 994
00159 34* TPRINT = CON(13)
00160 35* INITIALIZE TIME SUM BETWEEN OUTPUT INTERVALS
00161 36* TSUM = 0.0
00162 37* DOES OLD TIME PLUS THE OUTPUT INTERVAL EXCEED THE STOP TIME
00163 38* IF(CON(13)+CON(18).GT.CON(3)) CON(18) = CON(3)-CON(13)
00164 39* DONT EXCEED IT
00165 40* TSTEPN = CON(22)
00166 41* IF(TSTEPN.LE.CON(8)) GO TO 20
00167 42* TSTEPN = CON(8)
00168 43* GO TO 35
00169 44* DOES THE TIME SUM PLUS THE TIME STEP EXCEED OUTPUT INTERVAL
00170 45* IF(TSUM+TSTEPN-CON(18)) 30,35,25
00171 46* DONT EXCEED IT
00172 47* TSTEPN = CON(18)-TSUM
00173 48* GO TO 35
00174 49* DOES TIME SUM PLUS TWO TIME STEPS EXCEED OUTPUT INTERVAL
00175 50* IF(TSUM+2.0*TSTEPN.LE.CON(18)) GO TO 35
00176 51* APPROACH THE OUTPUT INTERVAL GRADUALLY
00177 52* TSTEPN = (CON(18)-TSUM)/2.0
00178 53* STORE DELTA TIME STEP IN THE CONSTANTS
00179 54* CON(2) = TSTEPN
00180 55* CALCULATE THE NEW TIME
00181 56* IF(PASS.GT.0.) GO TO 40
00182 57* CON(1) = TPRINT
00183 58* CON(2) = 0.0
00184 59* GO TO 45
00185 60* CON(1) = TPRINT+TSUM+TSTEPN
00186 61* COMPUTE THE MEAN TIME BETWEEN ITERATIONS
00187 62* CON(14) = (CON(1)+CON(13))/2.0
00188 63* LAX = CON(5)
00189 64* DN = CON(10)
00190 65* DD = 1.0-DN
00191 66* AN = CON(9)
00192 67* AA = 1.0-AN
00193 68* TSTEP = TSTEPN/2.0
00194 69* DO THE RELAXATION LOOP
00195 70* DO 240 K1 = 1,LAX
00196 71* KON(20) = K1
00197 72* J1 = 0
00198 73* RLXA = 0.0
00199 74* RLXD = 0.0
00200 75* IF(K1.GT.1) GO TO 105
00201 76* J2 = 1
00202 77* ZERO OUT ALL SOURCE LOCATIONS AND SHIFT TEMPERATURES
00203 78* DO 50 I = 1,NNC
00204 79* Q(I) = 0.0
00205 80* DO 55 I = 1,NNI
00206 81* LEI = IEI+I
00207 82* X(LEI) = T(I)
00208 83* CON(12) = 0.0
00209 84* CALL VARBLI
00210 85* CHECK THE BACKUP SWITCH
00211 86* IF(KON(12).NE.0) GO TO 15
00212 87* CHECK FOR FIRST PASS
00213 88* IF(PASS.GE.0.) GO TO 60
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00253 89* CALL OUTCAL
00254 90* PASS = 1.0
00255 91* GO TO 10
00256 92* 60 RC = 1.E+8
00257 93* JJ = 0
00257 94* C CALCULATE THE OLD CONTRIBUTION AND THE CSGMIN
00260 95* DO 100 I = 1,NND
00263 95* INCLUDE VARG,LIST
00263 96* FOLD DELTAY INTO THE CAPACITANCES
00264 97* IF(FLD(1,1,NSQ1(J1+1)),EQ.0) GO TO 2000
00266 97* NTYPE = FLD(0,5,NSQ2(J2))
00267 97* LA = FLD(5,17,NSQ2(J2))
00270 97* LK = FLD(22,14,NSQ2(J2))
00271 97* GO TO (1005,1010,1015,1020,1025,1030,1035,1040,1045), NTYPE
00272 97* CALL DID1WM(T(I),A(LA),XK(LK),C(I))
00273 97* GO TO 1999
00274 97* 1010 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00275 97* 1012 J2 = J2+1
00276 97* LA = FLD(5,17,NSQ2(J2))
00277 97* LK = FLD(22,14,NSQ2(J2))
00300 97* CALL DID1WM(T(I),A(LA),XK(LK),C2)
00301 97* GO TO 1998
00302 97* 1015 C1 = XK(LK)*XK(LA)
00303 97* GO TO 1012
00304 97* 1020 CALL DID1WM(T(I),A(LA),XK(LK),C1)
00305 97* J2 = J2+1
00306 97* LA = FLD(5,17,NSQ2(J2))
00307 97* LK = FLD(22,14,NSQ2(J2))
00310 97* C2 = XK(LK)*XK(LA)
00311 97* GO TO 1998
00312 97* 1025 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C(I))
00313 97* GO TO 1999
00314 97* 1030 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00315 97* 1032 J2 = J2+1
00316 97* LA = FLD(5,17,NSQ2(J2))
00317 97* LK = FLD(22,14,NSQ2(J2))
00320 97* CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C2)
00321 97* GO TO 1998
00322 97* 1035 C1 = XK(LK)*XK(LA)
00323 97* GO TO 1032
00324 97* 1040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),C1)
00325 97* J2 = J2+1
00326 97* LA = FLD(5,17,NSQ2(J2))
00327 97* LK = FLD(22,14,NSQ2(J2))
00330 97* C2 = XK(LK)*XK(LA)
00331 97* GO TO 1998
00332 97* 1045 CALL D2DIWM(T(I),CON(14),A(LA),XK(LK),C(I))
00333 97* GO TO 1999
00334 97* 1998 C(I) = C1+C2
00335 97* 1999 J2 = J2+1
00336 97* 2000 CONTINUE
00337 97* END
00340 98* C(I) = C(I)/TSTEP
00341 99* LE2 = IE2+1
00342 100* X(LE2) = 0.0
00343 101* INCLUDE VARG,LIST
00344 101* IF(FLD(4,1,NSQ1(J1+1)),EQ.0) GO TO 5000
00346 101* NTYPE = FLD(0,5,NSQ2(J2))
00347 101* LA = FLD(5,17,NSQ2(J2))
00350 101* LK = FLD(22,14,NSQ2(J2))

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*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

CNVARB,CNVARB

00351 101* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE
 00352 101* Q(I) = XK(LK)+Q(I)
 00353 101* GO TO 4999
 00354 101* 4010 Q1 = 0.0
 00355 101* 4012 CALL DID1WM(T(I),A(LA),XK(LK),62)
 00356 101* GO TO 4998
 00357 101* 4015 Q1 = 0.0
 00360 101* 4017 CALL DID1WM(CON(14),A(LA),XK(LK),62)
 00361 101* GO TO 4998
 00362 101* 4020 CALL DID1WM(CON(14),A(LA),XK(LK),61)
 00363 101* 4022 J2 = J2+1
 00364 101* LA = FLD(5,17,NSQ2(J2))
 00365 101* LK = FLD(22,14,NSQ2(J2))
 00366 101* GO TO 4017
 00367 101* 4025 Q1 = XK(LK)*XK(LA)
 00370 101* GO TO 4022
 00371 101* 4030 CALL DID1WM(CON(14),A(LA),XK(LK),61)
 00372 101* J2 = J2+1
 00373 101* LA = FLD(5,17,NSQ2(J2))
 00374 101* LK = FLD(22,14,NSQ2(J2))
 00375 101* Q2 = XK(LK)*XK(LA)
 00376 101* GO TO 4998
 00377 101* 4035 CALL DID1WM(CON(14),A(LA),XK(LK),61)
 00400 101* 4037 J2 = J2+1
 00401 101* LA = FLD(5,17,NSQ2(J2))
 00402 101* LK = FLD(22,14,NSQ2(J2))
 00403 101* GO TO 4012
 00404 101* 4040 Q1 = XK(LK)*XK(LA)
 00405 101* GO TO 4037
 00406 101* 4998 Q(I) = Q1+Q2+Q(I)
 00407 101* 4999 J2 = J2+1
 00410 101* 5000 CONTINUE
 00411 101* ENC
 00412 102* Q(I) = 2.0*Q(I)+C(I)*(I)+(I)+460.0
 00413 103* G2 = 0.0
 00414 104* 70 J1 = J1+1
 00415 105* LG = FLD(5,16,NSQ1(J1))
 00416 106* LTA = FLD(22,14,NSQ1(J1))
 00417 107* INCLUD VARG,LIST
 00420 107* IF(FLD(2,1,NSQ1(J1)),EQ,0) GO TO 3000
 00422 107* NTYPE = FLD(10,8,NSQ2(J2))
 00423 107* LA = FLD(5,17,NSQ2(J2))
 00424 107* LK = FLD(22,14,NSQ2(J2))
 00425 107* GOTO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
 \$ 2060,2065), NTYPE
 00426 107* 2005 TM = (Y(I)+I(LTA))/2.0
 00427 107* 2007 CALL DID1WM(TM,A(LA),XK(LK),6(LG))
 00430 107* GO TO 2999
 00431 107* 2010 TM = I(I)
 00432 107* GO TO 2007
 00433 107* 2015 CALL DID1WM(T(I),A(LA),XK(LK),61)
 00434 107* 2017 J2 = J2+1
 00435 107* LA = FLD(5,17,NSQ2(J2))
 00436 107* LK = FLD(22,14,NSQ2(J2))
 00437 107* CALL DID1WM(LTA),A(LA),XK(LK),62)
 00440 107* GO TO 2003
 00441 107* 2020 G1 = XK(LK)*XK(LA)
 00442 107* GO TO 2017
 00443 107* 2025 CALL DID1WM(T(I),A(LA),XK(LK),61)
 00444 107* J2 = J2+1

00-3

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CNVARB,CNVARB

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00445 107* LA = FLD(5,17,NSQ2(J2))
00446 107* LK = FLD(22,14,NSQ2(J2))
00447 107* G2 = XK(LK)*XK(LA)
00450 107* GO TO 2998
00451 107* 2030 TM = (T(I)+T(LTA))/2.0
00452 107* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00453 107* GO TO 2999
00454 107* 2035 TM = T(I)
00455 107* GO TO 2032
00456 107* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00457 107* 2042 J2 = J2+1
00460 107* LA = FLD(5,17,NSQ2(J2))
00461 107* LK = FLD(22,14,NSQ2(J2))
00462 107* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00463 107* GO TO 2998
00464 107* 2045 G1 = XK(LK)*XK(LA)
00465 107* GO TO 2042
00466 107* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00467 107* J2 = J2+1
00470 107* LA = FLD(5,17,NSQ2(J2))
00471 107* LK = FLD(22,14,NSQ2(J2))
00472 107* G2 = XK(LK)*XK(LA)
00473 107* GO TO 2998
00474 107* 2055 TM = (T(I)+T(LTA))/2.0
00475 107* CALL D2DIM(TM,CON(14),A(LA),XK(LK),G(LG))
00476 107* GO TO 2999
00477 107* 2060 TM = T(LTA)
00500 107* GO TO 2007
00501 107* 2065 TM = T(LTA)
00502 107* GO TO 2032
00503 107* 2998 G(LG) = 1./(1./G1+1./G2)
00504 107* IF(FLD(3,1,NSQ1(J1)).EQ.1) G(LG) = G1*G2
00506 107* 2999 J2 = J2+1
00507 107* 3000 CONTINUE
00510 107* END
00511 108* GV = G(LG)
00512 109* T1 = X(IE1+I)+460.0
00513 110* T2 = X(IE1+LTA)+460.0
00514 111* CHECK FOR RADIATION CONDUCTOR
00516 113* IF(FLD(3,1,NSQ1(J1)).EQ.0) GO TO 75
00517 114* G2 = G2*GV*(T1+T1+T2)*(T1+T2)
00520 115* X(LE2) = X(LE2)+GV*(T2**4-T1**4)
00521 116* GO TO 80
00522 117* 75 X(LE2) = X(LE2)+GV*(T2-T1)
00523 118* G2 = G2*GV
00523 118* 80 IF(NSQ1(J1).GT.0) GO TO 70
00523 119* C OBTAIN THE MINIMUM STABILITY CRITERIA
00525 120* RI = C(I)/G2
00526 121* IF(R1.GE.RC) GO TO 100
00530 122* RC = R1
00531 123* KON(35) = I
00532 124* 100 CONTINUE
00533 125* CON(17) = RC*TSSTEP
00535 126* IF(ARC.LE.0.) GO TO 993
00537 127* BETA0 = 2.0*CON(17)/CON(2)
00540 128* IF(INF.TAO.GI.1.0) BETA0 = 1.0
00542 129* BETA0 = 2.0-BETA0
00542 130* C CONVERT TEMPERATURES TO RANKINE
00543 131* DO 65 I = 1,NNT
00546 132* LEI = IEI+I

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CNVARB,CNVARB

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00547 133*      T(I) = T(I)+460.
00550 134*      X(LE1) = X(LE1)+460.
00550 135*      C    NOW RELAX THE NETWORK BY SUCCESSIVE POINT AND EXTRAPOLATION
00552 136*      J1 = 0
00553 137*      JJ = JJ+1
00554 138*      DO 165 I = 1,AND
00557 139*      IF(K1.GT.1) GO TO 110
00561 140*      LE2 = IE2+1
00562 141*      G(I) = G(I)+BETA0*X(LE2)
00563 142*      110 R1 = 0.0
00564 143*      S = 0.0
00565 144*      OSUM = 0.0
00566 145*      GSUM = 0.0
00567 146*      115 J1 = J1+1
00570 147*      LG = FLD(5,16,NS01(J1))
00571 148*      LTA = FLD(22,14,NS01(J1))
00572 149*      GV = G(LG)
00573 150*      C    CHECK FOR RADIATION CONDUCTOR
00575 151*      IF(FLD(3,1,NS01(J1)).EQ.0) GO TO 120
00576 152*      R1 = R1+GV
00577 153*      OSUM = OSUM+GV*T(LTA)**4
00577 154*      GO TO 125
00600 155*      120 GSUM = GSUM+GV
00601 156*      OSUM = OSUM+GV*T(LTA)
00602 157*      125 IF(NS01(J1).GT.0) GO TO 115
00604 158*      GSUM = G(I)+BETAN*OSUM
00605 159*      OSUM = C(I)+BETAN*GSUM
00605 160*      C    DAMPEN RADIATION ON THIS NODE IF PRESENT
00606 161*      IF(R1.LE.0.) GO TO 145
00610 162*      R1 = R1*BETAN
00611 163*      R2 = R1*T(I)**4
00612 164*      T2 = (OSUM-R2)/GSUM
00613 165*      S = (R1+R2)/2.0
00614 166*      C    OBTAIN THE NEW TEMPERATURE
00614 167*      T2 = DN*((OSUM-S)/GSUM)+DD*T(I)
00615 168*      C    OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
00615 169*      T1 = ABS(T(I)-T2)
00616 170*      C    STORE THE NEW AND OLD TEMPERATURES
00617 171*      GO TO (160,155,150), JJ
00620 172*      150 LE2 = IE2+1
00621 173*      LE3 = IE3+1
00622 174*      R1 = T2-T(I)
00622 175*      X(LE2) = T(I)
00623 176*      X(LE3) = R1/(R1-X(LE3))
00624 177*      GO TO 160
00625 178*      155 LE3 = IE3+1
00626 179*      X(LE3) = T2-T(I)
00627 180*      T(I) = T2
00630 181*      160 IF(RLXD.GE.11) GO TO 165
00631 182*      RLXD = 11
00633 183*      KK1 = 1
00634 184*      165 CONTINUE
00635 185*      GO TO (160,180,170), JJ
00637 186*      C    PERFORM LINEAR EXTRAPOLATION ON THE ERROR FUNCTION CURVE
00637 187*      JJ = 0
00640 188*      170 IF(KOH(34).NE.0) GO TO 180
00641 189*      DO 175 I = 1,NH0
00643 190*      LE3 = 1E3+1
00646 191*      C    SEE IF THE EXTRAPOLATION IS ALLOWABLE
00646 192*

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CNVARB,CNVARB

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00647 193* IF(X(LE3).GE.0.) GO TO 175
00647 194* LIMIT THE EXTRAPOLATION
00651 195* IF(X(LE3).LT.-10.) X(LE3) = -10.
00653 196* LE2 = IE2+1
00654 197* T(I) = X(LE3)*X(LE2)+(1.0-X(LE3))*T(I)
00655 198*
00657 199* 175 CONTINUE
00661 200* 180 IF(NNA.LE.0) GO TO 220
00662 201* JJI = JI
00663 202* JJJ = JJ
00663 203* DO 230 I = 1,NNT
00666 204* T(I) = T(I)-460.
00670 205* DO 215 I = NN,NNC
00673 206* GSUM = 0.0
00674 207* IF(K1.GT.1) GO TO 6000
00676 208* INCLUDE VR02,LIST
00677 209* IF(FLD(4,1,NSQ1(JJI+1)),EQ.0) GO TO 6000
00701 207* NTYP = FLD(10,5,NSQ2(JJ2))
00702 207* LA = FLD(5,17,NSQ2(JJ2))
00703 207* LK = FLD(22,14,NSQ2(JJ2))
00704 207* GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYP
00705 207* 5005 Q(L) = XK(LK)+Q(L)
00706 207* GO TO 5999
00707 207* 5010 Q1 = 0.0
00710 207* 5012 CALL DID1WM(T(L),A(LA),XK(LK),Q2)
00711 207* GO TO 5998
00712 207* 5015 Q1 = 0.0
00713 207* 5017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
00714 207* GO TO 5998
00715 207* 5020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00716 207* 5022 JJ2 = JJ2+1
00717 207* LA = FLD(5,17,NSQ2(JJ2))
00720 207* LK = FLD(22,14,NSQ2(JJ2))
00721 207* GO TO 5017
00722 207* 5025 Q1 = XK(LK)*XK(LA)
00723 207* GO TO 5022
00724 207* 5030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00725 207* JJ2 = JJ2+1
00726 207* LA = FLD(5,17,NSQ2(JJ2))
00727 207* LK = FLD(22,14,NSQ2(JJ2))
00730 207* Q2 = XK(LK)*XK(LA)
00731 207* GO TO 5998
00732 207* 5035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00733 207* 5037 JJ2 = JJ2+1
00734 207* LA = FLD(5,17,NSQ2(JJ2))
00735 207* LK = FLD(22,14,NSQ2(JJ2))
00736 207* GO TO 5012
00737 207* 5040 Q1 = XK(LK)*XK(LA)
00740 207* GO TO 5037
00741 207* 5998 Q(L) = Q1+Q2+Q(L)
00742 207* 5999 JJ2 = JJ2+1
00743 207* 6000 CONTINUE
00744 207* END
00745 208* OSUM = Q(I)
00746 209* L = I
00747 210* 185 JJI = JJI+1
00750 211* LG = FLD(5,16,NSQ1(JJI))
00751 212* LTA = FLD(22,14,NSQ1(JJI))
00752 213* IF(K1.GT.2) GO TO 4000
00754 214* INCLUDE VR62,LIST
00754 215* CHECK FOR RADIATION CONDUCTOR

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CNVARB,CNVARB

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01053 220*      GO TO 195
01054 221*      190 GV = G(LG)
01055 222*      195 GSUM = GSUM+GV
01056 223*      QSUM = GSUM+GV*T2
01056 224*      C CHECK FOR NEGATIVE CONDUCTOR
01057 225*      IF(NSG(JJI).GT.0) GO TO 185
01057 226*      C CALCULATE THE NEW TEMPERATURE
01061 227*      T2 = AN*OSUM/GSUM+AA*T1
01062 228*      T1 = ABS(T2-T1)
01063 229*      T(I) = T2-460.0
01064 230*      IF(RLXA.GE.T1) GO TO 215
01066 231*      RLXA = T1
01067 232*      KK2 = I
01070 233*      215 CONTINUE
01072 234*      DO 235 I = 1,N/IT
01075 235*      235 T(I) = T(I)+460.
01075 236*      C SEE IF THE ARITHMETIC RELAXATION CRITERIA WAS MET
01077 237*      IF(PLXA.GT.CON(19)) GO TO 225
01077 238*      C SEE IF THE DIFFUSION RELAXATION CRITERIA WAS MET
01101 239*      220 IF(RLX).LE.CON(26)) GO TO 245
01103 240*      225 IF(KON(7).EQ.0) GO TO 240
01105 241*      CALL OUTCAL
01106 242*      240 CONTINUE
01110 243*      IF(KON(20).GE.65) CALL TOPLIN
01112 244*      WRITE(6,B82)
01114 245*      KON(28) = KON(28)+2
01114 246*      C SEE IF THE TEMPERATURE CHANGES WERE TOO LARGE
01115 247*      245 TCGD = 0.0
01116 248*      TCGA = 0.0
01117 249*      DO 250 I = 1,N/ND
01122 250*      LE = IE1+I
01123 251*      C(I) = C(I)*TSTEP
01124 252*      T1 = ABS(T(I)-X(LE))
01125 253*      IF(TCGD.GT.T1) GO TO 250
01127 254*      TCGD = T1
01130 255*      KON(36) = I
01131 256*      250 CONTINUE
01133 257*      IF(TCGD.LE.CON(6)) GO TO 265
01135 258*      TSTEPN = 0.95*TSTEPN+CON(16)/TCGD
01136 259*      255 DO 260 I = 1,N/NT
01141 260*      LE = IE1+I
01142 261*      260 T(I) = X(LE)-460.
01144 262*      GO TO 30
01145 263*      265 IF(HNA.LE.0) GO TO 275
01147 264*      DO 270 I = NN/NNC
01152 265*      LE = IE1+I
01153 266*      T1 = ABS(T(I)-X(LE))
01154 267*      IF(TCGA.GT.T1) GO TO 270
01156 268*      TCGA = T1
01157 269*      KON(38) = I
01160 270*      270 CONTINUE
01162 271*      IF(TCGA.LE.CON(11)) GO TO 275
01164 272*      TSTEPN = 0.95*TSTEPN+CON(11)/TCGA
01165 273*      GO TO 255
01165 274*      C CONVERT TEMPERATURES BACK TO FARENHEIT
01156 275*      275 DO 280 I = 1,N/IT
01171 276*      280 T(I) = T(I)-460.
01171 277*      C STORE THE TEMPERATURE AND RELAXATION CHANGES
01173 278*      CON(15) = TCGD
01174 279*      CON(16) = TCGA

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CNVARB,CNVARB

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01175 280*      CON(27) = RLXD
01176 281*      IF(ARLXA.GT.RLXD) GO TO 285
01200 282*      KK2 = KK1
01201 283*      RLXA = RLXD
01202 284*      KON(37) = KK2
01203 285*      KON(30) = RLXA
01204 286*      CON(12) = 0.0
01205 287*      CALL VARBL2
01206 288*      CHECK THE BACKUP SWITCH
01207 289*      IF(KON(12).NE.0) GO TO 255
01210 291*      ADVANCE TIME
01211 292*      CON(13) = CON(11)
01211 293*      TSUM = TSUM+TSTEPN
01212 294*      CHECK FOR TIME TO PRINT
01212 295*      IF(TSUM.GE.CON(18)) GO TO 290
01214 296*      CHECK FOR PRINT EVERY ITERATION
01216 297*      IF(KON(7).NE.0) CALL OUTCAL
01216 298*      GO TO 10
01217 299*      TRY TO EVEN THE OUTPUT INTERVALS
01220 300*      CALL OUTCAL
01220 301*      IS TIME GREATER THAN END COMPUTE TIME
01221 302*      IF(CON(1)*1.000001.LT.CON(3)) GO TO 5
01223 303*      NTH = IE1
01224 304*      NDIM = MLA
01225 305*      RETURN
01226 306*      990 WRITE(6,880)
01230 307*      GO TO 1000
01231 308*      991 WRITE(6,881)
01233 309*      GO TO 1000
01234 310*      993 WRITE(6,883)
01236 311*      GO TO 1000
01242 312*      994 WRITE(6,884) NDIM
01243 314*      GO TO 1090
01245 315*      995 WRITE(6,885)
01246 316*      GO TO 1000
01250 317*      996 WRITE(6,886)
01251 318*      GO TO 1000
01253 319*      997 WRITE(6,887)
01254 320*      GO TO 1000
01256 321*      998 WRITE(6,888)
01257 322*      GO TO 1000
01261 323*      999 WRITE(6,889)
01262 324*      1000 CALL OUTCAL
01263 325*      CALL EXIT
01264 326*      880 FORMAT(29H TRANSIENT TIME NOT SPECIFIED)
01265 327*      881 FORMAT(45H CNVARB REQUIRES LONG PSEUDO-COMPUTE SEQUENCE)
01266 328*      882 FORMAT(28H RELAXATION CRITERIA NOT MET)
01267 329*      883 FORMAT(24H CSGMIN ZERO OR NEGATIVE)
01270 330*      884 FORMAT(18,20H LOCATIONS AVAILABLE)
01271 331*      885 FORMAT(10H NO DRLXCA)
01272 332*      886 FORMAT(10H NO DTIMEI)
01273 333*      887 FORMAT(10H NO ARLXCA)
01274 334*      888 FORMAT(19H NO OUTPUT INTERVAL)
01275 335*      889 FORMAT(9H NO NLOOP)
                                END

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END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(5)
CNVARB SYMBOLIC
CNVARB CODE RELOCATABLE

C. **COMPUTER LISTINGS OF SINDA STEADY STATE SOLUTION ROUTINES**

	Page
CINDSS	C-2
CINDSL	C-11
CINDSM	C-20

QIW FOR.* CINDSS,CINDSS
UNIVAC 1108 FORTRAN V ATHENA VERSION: 131K-100 CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 14:00:06

: 1 FEB 71

SUBROUTINE CINDSS ENTRY POINT 003062

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	003077
0000	*CONST+TEMP	000106
0002	*SIMPLE VAR	000045
0004	*ARRAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COND	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARL1
0023	OUTCAL
0024	DIDIM
0025	PLYAWM
0026	D2DIM
0027	NONLIN
0030	VARL2
0031	EXIT
0032	NERR25
0033	NWDUS
0034	NI025
0035	NER105

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C	0017	R	000000	CON
0012	R	000000	G	0002	R	000040	GSUM	0002	R	000033	GV
0006	R	000000	H	0002	I	000007	I	0002	I	000003	IE
0002	I	000010	J1	0002	I	000011	J2	0015	I	000000	K
0002	I	000041	L	0002	I	000020	LA	0002	I	000005	LAX
0002	I	000024	LG	0002	I	000021	LK	0002	I	000006	LS01
0021	I	000005	NAT	0021	I	000004	NCT	0020	I	000000	NDIM
0002	I	000001	NR1	0021	I	000001	NNA	0002	I	000002	NMC
0013	I	000000	NS01	0014	I	000000	NS02	0020	I	000001	NTH
0002	I	000035	N1	0002	I	000042	N2	0002	R	000000	PASS
0002	R	000043	OUT	0002	R	000022	O1	0002	R	000023	O2
0007	R	000000	T	0002	R	000026	TM	0002	R	000031	T1
0002	R	000014	DN	0002	R	000015	DD	0002	R	000017	NTYPE
0002	R	000030	G2	0002	R	000027	G1	0002	R	000000	Q
0002	I	000037	JJ2	0017	I	000036	JJ1	0011	R	000000	Q
0002	I	000006	K1	0002	I	000000	KON	0002	R	000013	RLXA
0002	I	000034	LF1	0002	I	000016	LE	0002	R	000032	T2
0002	I	000025	LTA	0021	I	000007	LS02	0021	I	000003	NGT
0002	I	000004	NLA	0021	I	000003	NGT	0021	I	000000	NND
0002	I	000002	NJT	0021	I	000000	NND	0021	I	000002	NPT
0020	R	000002	NP	0002	I	000017	NTYPE	0020	R	000002	NP
0002	R	000044	QTN	0011	R	000000	Q	0002	R	000044	QTN
0002	R	000012	RLXD	0002	R	000013	RLXA	0002	R	000012	RLXD
0020	P	000002	X	0002	R	000032	T2	0020	P	000002	X

CINROSS, CINROSS

0015	R	000000	XK	0001	000104	10L	0001	000036	1030L	0001	002537	110L	002543	115L
0001		002605	125L	0001	000125	15L	0001	000140	155G	0001	002620	155L	000145	161G
0001		002634	165L	0001	002712	170L	0001	002721	175L	0001	000256	1998L	000261	1999L
0001		000264	2000L	0001	000624	2005L	0001	000631	2007L	0001	000651	2010L	000654	2015L
0001		000572	2017L	0001	000726	2020L	0001	000734	2025L	0001	000203	203G	000774	2030L
0001		001001	2032L	0001	001024	2035L	0001	001027	2040L	0001	001050	2042L	001107	2045L
0001		001115	2050L	0001	001160	2055L	0001	001206	2060L	0001	001212	2065L	000215	211G
0001		000121	25L	0001	001216	2998L	0001	001244	2999L	0001	001247	3000L	001777	3005L
0001		002005	3007L	0001	002025	3010L	0001	002031	3015L	0001	002050	3017L	002104	3020L
0001		002112	3025L	0001	002153	3030L	0001	002161	3032L	0001	002204	3035L	002210	3040L
0001		002232	3042L	0001	002271	3045L	0001	002277	3050L	0001	002343	3055L	002372	3060L
0001		002376	3065L	0001	000535	35L	0001	002402	3998L	0001	002430	3999L	001300	40L
0001		002433	4000L	0001	000327	4005L	0001	000334	4010L	0001	000335	4012L	000354	4015L
0001		000355	4017L	0001	000372	4020L	0001	000406	4022L	0001	000423	4025L	000431	4030L
0001		00047	4035L	0001	000593	4037L	0001	000520	4040L	0001	001431	434G	001303	45L
0001		000526	4998L	0001	000532	4999L	0001	001351	50L	0001	000535	5000L	001502	5005L
0001		001510	5010L	0001	001511	5012L	0001	001531	5015L	0001	001532	5017L	001547	5020L
0001		001563	5022L	0001	001600	5025L	0001	001606	5030L	0001	001644	5035L	001660	5037L
0001		001675	5040L	0001	001703	5998L	0001	001710	5999L	0001	001713	6000L	002631	657G
0001		001406	70L	0001	001416	75L	0001	001713	80L	0001	002470	85L	000000	862F
0000		000001	823F	0000	000007	844F	0000	000020	885F	0000	000030	886F	000035	887F
0000		000040	885F	0000	000043	889F	0001	002473	90L	0001	003000	994L	000035	887F
0001		003015	997L	0001	003023	998L	0001	003031	999L	0001	003000	994L	0003006	996L

1* SUBROUTINE CINROSS
2* STEADY STATE EXECUTION SUBROUTINE FOR SINDA FORTRAN V
3* THE SHORT PSEUDO-COMPUTE SEQUENCE IS REQUIRED
4* DIFFUSION NODES RECEIVE A BLOCK ITERATION
5* ARITHMETIC NODES RECEIVE A SUCCESSIVE POINT ITERATION
6* OVER-RELAXATION IS ALLOWED, THE DAMPING FACTORS ARE ADDRESSABLE
7* INCLUDE COMM,LIST
8* COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
9* COMMON /PCI/NSQ(1) /PC2/NSQ2(1) /KONST/K(1) /ARRAY/A(1)
10* COMMON /FIXCON/KON(1) /XSPACE/NDIM,NTH,X(1)
11* COMMON /DIMENS/ NRD,NNA,NRT,NGT,NCI,MAT,LS01,LS02
12* DIMENSION CON(1),XK(1),NK(1)
13* EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NK(1))
14* END
15* INCLUDE DEFF,LIST
16* C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
17* C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
18* C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
19* C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
20* C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR*EXPLICIT (CSGFAC)
21* C CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
22* C CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
23* C CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
24* C CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
25* C CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMPA)
26* C CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPO)
27* C CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
28* C CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
29* C CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
30* C CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
31* C CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)
32* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
33* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSSMIN)

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00113 6* CONTROL CONSTANT IS CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 8* CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 8* CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED, INTEGER (LOOPCT)
00113 8* CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEL)
00113 8* CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
00113 8* CC23 CONTAINS THE C/56 MAXIMUM (CSGMAX)
00113 8* CC24 CONTAINS THE C/56 RANGE CALCULATED (CSGRAL)
00113 8* CC25 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (CSGRCL)
00113 8* CC26 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
00113 8* CC27 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 8* CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 8* CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
00113 8* CC31 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 8* CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGRAL)
00113 8* CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 8* CC34 CONTAINS THE HOCOPY SWITCH FOR MATRIX USERS (HOCOPY)
00113 8* CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 8* CC36 CONTAINS RELATIVE NODE NUMBER OF DIMPCC
00113 8* CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 8* CC38 CONTAINS RELATIVE NODE NUMBER OF ATPCC
00113 8* CC39-40-41-42-45 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 8* CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 8* CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSS (LAXFAC)
00113 8* CC50 IS NOT USED AT PRESENT
00114 8* END
00115 9* IF (KON(5).LE.0) GO TO 999
00117 10* IF (CON(9).LE.0.) CON(9) = 1.0
00121 11* IF (CON(10).LE.0.) CON(10) = 1.0
00123 12* IF (NNA.GT.0.AND.CON(19).LE.0.) GO TO 998
00125 13* IF (NND.GT.0.AND.CON(26).LE.0.) GO TO 997
00127 14* IF (KON(31).NE.0) GO TO 994
00131 15* PASS = -1.0
00132 16* NN = N/ID+1
00133 17* NNC=NNA+NNND
00134 18* IE=NTN
00135 19* NLA = I/IDIM
00136 20* NTH=NTN+NNND
00137 21* NDIM=NDIM+NNND
00140 22* IF (NDIM.LI.0) GO TO 996
00142 23* CON(1) = CON(13)
00143 24* CON(2) = 0.0
00144 25* CON(14) = CON(13)
00145 26* GO TO 15
00146 27* CON(1) = CON(13)+CON(18)
00147 28* IF (CON(1)-CON(3).GT.0.) CON(1) = CON(3)
00151 29* CON(14) = (CON(1)+CON(13))/2.0
00152 30* CON(2) = CON(1)-CON(13)
00152 31* COMPUTE STEADY STATE TEMPERATURES
00153 32* LAX = KON(5)
00154 33* DO 120 K1 = 1,LAX
00157 34* KON(20) = K1
00157 35* ZERO OUT ALL SOURCE LOCATIONS
00160 36* DO 20 I = 1,NNC
00163 37* O(I)=0.
00165 38* CALL VAPOLI
00166 39* IF (PASS.GE.0.) GO TO 25
00170 40* CALL OUTCAL
00171 41* PASS = 1.0
00172 42* 25 J1 = 0

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00173 43* J2 = 1
00174 44* RLXD = 0.0
00175 45* RLXA = 0.0
00176 46* IF(NND.LE.0) GO TO 75
00200 47* DN = CON(10)
00201 48* DD = 1.0-DN
00201 49* ZERO OUT EXTRA LOCATIONS
00202 50* DO 30 I = 1,NND
00205 51* LE=IE+1
00206 52* 30 X(LE)=0.0
00206 53* DO A BLOCK ITERATION ON THE DIFFUSION NODES
00210 54* DO 70 I = 1,NND
00213 55* LE=IE+1
00214 56* INCLUDE DUMC,LIST
00215 57* IF(FLD(1,1),NSG1(JI+1),EQ.0) GO TO 2000
00217 56* NTYPE = FLD(0,5,NSQ2(J2))
00220 56* GO TO (1999,1998,1998,1998,1999,1998,1998,1998,1999), NTYPE
00221 56* 1998 J2 = J2+1
00222 56* 1999 J2 = J2+1
00223 56* 2000 CONTINUE
00224 56* END
00225 57* INCLUDE VARO,LIST
00226 57* IF(FLD(4,1),NSQ1(JI+1),EQ.0) GO TO 5000
00230 57* NTYPE = FLD(0,5,NSQ2(J2))
00231 57* LA = FLD(5,17,NSQ2(J2))
00232 57* LK = FLD(22,14,NSQ2(J2))
00233 57* GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE
00234 57* 4005 G(I) = XK(LK)+G(I)
00235 57* GO TO 4999
00236 57* 4010 G1 = 0.0
00237 57* 4012 CALL DIDIWM(T(I),A(LA),XK(LK),G2)
00240 57* GO TO 4998
00241 57* 4015 G1 = 0.0
00242 57* 4017 CALL DIDIWM(CON(14),A(LA),XK(LK),G2)
00243 57* GO TO 4998
00244 57* 4020 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00245 57* 4022 J2 = J2+1
00246 57* LA = FLD(5,17,NSQ2(J2))
00247 57* LK = FLD(22,14,NSQ2(J2))
00250 57* GO TO 4017
00251 57* 4025 G1 = XK(LK)*XK(LA)
00252 57* GO TO 4022
00253 57* 4030 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00254 57* J2 = J2+1
00255 57* LA = FLD(5,17,NSQ2(J2))
00256 57* LK = FLD(22,14,NSQ2(J2))
00257 57* G2 = XK(LK)*XK(LA)
00260 57* GO TO 4998
00261 57* 4035 CALL DIDIWM(CON(14),A(LA),XK(LK),G1)
00262 57* 4037 J2 = J2+1
00263 57* LA = FLD(5,17,NSQ2(J2))
00264 57* LK = FLD(22,14,NSQ2(J2))
00265 57* GO TO 4012
00266 57* 4040 G1 = XK(LK)*XK(LA)
00267 57* GO TO 4037
00270 57* 4998 G(I) = G1+G2+G(I)
00271 57* 4999 J2 = J2+1
00272 57* 5000 CONTINUE
00273 57* END
00274 58* J5 J1 = J1+1

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CINDSS,CINDSS

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00275 59* LG = FLD(5,16,NSQ1(J1))
00276 60* IF(LG.EQ.0) GO TO 50
00300 61* LTA = FLD(22,14,NSQ1(J1))
00301 62* INCLUDE VARG,LIST
00301 63* CHECK FOR RADIATION CONDUCTOR
00302 63* IF(FLD(2,1,NSQ1(J1)).EQ.0) GO TO 3000
00304 63* NTYPE = FLD(0,5,NSQ2(J2))
00305 63* LA = FLD(5,17,NSQ2(J2))
00306 63* LK = FLD(22,14,NSQ2(J2))
00307 63* GOTO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
00310 63* 2060,2065), NTYPE
00311 63* 2005 TM = (T(I)+T(LTA))/2.0
00312 63* 2007 CALL D2DIMM(TM,A(LA),XK(LK),G(LG))
00313 63* GO TO 2999
00314 63* 2010 TM = T(I)
00315 63* GO TO 2007
00316 63* 2015 CALL D1DIMM(T(I),A(LA),XK(LK),G1)
00317 63* 2017 J2 = J2+1
00320 63* LA = FLD(5,17,NSQ2(J2))
00321 63* LK = FLD(22,14,NSQ2(J2))
00322 63* CALL D1DIMM(T(LTA),A(LA),XK(LK),G2)
00323 63* GO TO 2998
00324 63* 2020 G1 = XK(LK)*XK(LA)
00325 63* GO TO 2017
00326 63* 2025 CALL D1DIMM(T(I),A(LA),XK(LK),G1)
00327 63* J2 = J2+1
00330 63* LA = FLD(5,17,NSQ2(J2))
00331 63* LK = FLD(22,14,NSQ2(J2))
00332 63* G2 = XK(LK)*XK(LA)
00333 63* GO TO 2998
00334 63* 2030 TM = (T(I)+T(LTA))/2.0
00335 63* 2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
00336 63* GO TO 2999
00337 63* 2035 TM = T(I)
00340 63* GO TO 2032
00341 63* 2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00342 63* 2042 J2 = J2+1
00343 63* LA = FLD(5,17,NSQ2(J2))
00344 63* LK = FLD(22,14,NSQ2(J2))
00345 63* CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
00346 63* GO TO 2998
00347 63* 2045 G1 = XK(LK)*XK(LA)
00350 63* GO TO 2042
00351 63* 2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
00352 63* J2 = J2+1
00353 63* LA = FLD(5,17,NSQ2(J2))
00354 63* LK = FLD(22,14,NSQ2(J2))
00355 63* G2 = XK(LK)*XK(LA)
00356 63* GO TO 2998
00357 63* 2055 TM = (T(I)+T(LTA))/2.0
00360 63* CALL D2DIMM(TM,CON(14),A(LA),XK(LK),G(LG))
00361 63* GO TO 2999
00362 63* 2060 TM = T(LTA)
00363 63* GO TO 2007
00364 63* 2065 TM = T(LTA)
00365 63* GO TO 2032
00366 63* 2994 G(LG) = 1./(1./G1+1./G2)
00370 63* IF(FLD(3,1,NSQ1(J1)).EQ.1) G(LG) = G1*G2
00371 63* 2999 J2 = J2+1
3000 CONTINUE

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CINDSS,CINDSS

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00372 63*
00373 64*
00375 65*
00376 66*
00377 67*
00400 68*
00401 69*
00402 70*
00403 71*
00403 72*
00404 73*
00406 74*
00407 76*
00410 76*
00410 77*
00411 78*
00413 79*
00413 80*
00414 81*
00414 82*
00415 83*
00415 84*
00416 85*
00420 86*
00421 87*
00422 88*
00424 89*
00425 90*
00427 91*
00430 92*
00431 93*
00432 94*
00432 95*
00433 96*
00436 97*
00437 98*
00440 99*
00441 99*
00443 99*
00444 99*
00445 99*
00446 99*
00447 99*
00450 99*
00451 99*
00452 99*
00453 99*
00454 99*
00455 99*
00456 99*
00457 99*
00460 99*
00461 99*
00462 99*
00463 99*
00464 99*
00465 99*
00466 99*
00467 99*
00470 99*

END
IF(FLD(3,1,NS01(J1)),EQ,0) GO TO 40
T1 = T(I)+460.0
T2 = T(LTA)+460.0
GV = G(LG)*(T1+T1+T2+T2)*(T1+T2)
GO TO 45
40 GV = G(LG)
45 X(LE) = X(LE)+GV
G(I) = G(I)+GV*(LTA)
CHECK FOR ADJOINING DIFFUSION NODE, WATCH FOR ONE WAY CONDUCTOR
IF(LTA.GT.NND.OR.FLD(21,1,NS01(J1)),EQ,1) GO TO 50
LE1 = IE+LTA
X(LE1) = X(LE1)+GV
G(LTA) = G(LTA)+GV*(I)
CHECK FOR LAST CONDUCTOR TO THIS NODE
IF(NS01(J1).GT,0) GO TO 35
T2 = DD*(I)+DN*(I)/X(LE)
OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
T1=ABS(T(I)-T2)
STORE THE NEW TEMPERATURE
T(I)=T2
SAVE THE MAXIMUM DIFFUSION RELAXATION CHANGE
IF(RLXD.GT.T1) GO TO 70
RLXD = T1
N1 = I
70 CONTINUE
CON(27) = RLXD
IF(NMA.LE.0) GO TO 115
75 DN = CON(9)
DD = 1.0-DN
JJ1 = J1
JJ2 = J2
DO A SUCCESSIVE POINT ITERATION ON THE ARITHMETIC NODES
DO 110 I = NN,RNC
650% = 0.0
L = I
INCLUDE VR02,LIST
IF(FLD(4,1,NS01(JJ1+1)),EQ,0) GO TO 6000
NTYPE = FLD(10,5,NS02(JJ2))
LA = FLD(5,17,NS02(JJ2))
LK = FLD(22,14,NS02(JJ2))
GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5050), NTYPE
5005 O(L) = XK(LK)+C(L)
GO TO 5999
5010 O1 = 0.0
5012 CALL DID1WM(T(L),A(LA),XK(LK),O2)
GO TO 5998
5015 O1 = 0.0
5017 CALL DID1WM(CON(14),A(LA),XK(LK),O2)
GO TO 5998
5020 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
5022 JJ2 = JJ2+1
LA = FLD(5,17,NS02(JJ2))
LK = FLD(22,14,NS02(JJ2))
GO TO 5017
5025 O1 = XK(LK)*XK(LA)
GO TO 5022
5030 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
JJ2 = JJ2+1
LA = FLD(5,17,NS02(JJ2))

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*NEW

LK = FLD(22,14,NS02(JJ2))
 02 = XK(LK)*XK(LA)
 60 TO 5998
 5035 CALL DID1WM(CON(14),A(LA),XK(LK),01)
 5037 JJ2 = JJ2+1
 LA = FLD(5,17,NS02(JJ2))
 LK = FLD(22,14,NS02(JJ2))
 60 TO 5012
 5040 01 = XK(LK)*XK(LA)
 60 TO 5037
 5998 0(L) = 01+02+G(L)
 5999 JJ2 = JJ2+1
 6000 CONTINUE
 END
 80 JJ1 = JJ1+1
 LG = FLD(5,16,NS01(JJ1))
 LTA = FLD(22,14,NS01(JJ1))
 INCLUDE VRG2*LIST
 CHECK FOR RADIATION CONDUCTOR
 IF(FLD(12,1,NS01(JJ1)).EQ.0) GO TO 4000
 NTYPE = FLD(10,5,NS02(JJ2))
 LA = FLD(5,17,NS02(JJ2))
 LK = FLD(22,14,NS02(JJ2))
 60 TO (3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,
 3060,3065), NTYPE
 3005 TM = (T(L)+T(LTA))/2.0
 3007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
 60 TO 3999
 3010 TM = T(L)
 60 TO 3007
 3015 CALL DID1WM(T(L),A(LA),XK(LK),G1)
 3017 JJ2 = JJ2+1
 LA = FLD(5,17,NS02(JJ2))
 LK = FLD(22,14,NS02(JJ2))
 CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
 60 TO 3998
 3020 G1 = XK(LK)*XK(LA)
 60 TO 3017
 3025 CALL DID1WM(T(L),A(LA),XK(LK),G1)
 JJ2 = JJ2+1
 LA = FLD(5,17,NS02(JJ2))
 LK = FLD(22,14,NS02(JJ2))
 G2 = XK(LK)*XK(LA)
 60 TO 3998
 3030 TM = (T(L)+T(LTA))/2.0
 3032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
 60 TO 3999
 3035 TM = T(L)
 60 TO 3032
 3040 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)
 3042 JJ2 = JJ2+1
 LA = FLD(5,17,NS02(JJ2))
 LK = FLD(22,14,NS02(JJ2))
 CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
 60 TO 3998
 3045 G1 = XK(LK)*XK(LA)
 60 TO 3042
 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)
 JJ2 = JJ2+1
 LA = FLD(5,17,NS02(JJ2))

00471 99*
 00472 99*
 00473 99*
 00474 99*
 00475 99*
 00476 99*
 00477 99*
 00500 99*
 00501 99*
 00502 99*
 00503 99*
 00504 99*
 00505 99*
 00506 99*
 00507 100*
 00510 101*
 00511 102*
 00512 103*
 00512 104*
 00513 104*
 00515 104*
 00516 104*
 00517 104*
 00520 104*
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 00522 104*
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 00560 104*
 00561 104*
 00562 104*
 00563 104*

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CINDSS,CINDSS

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00564 104* LK = FLD(22,14,NSQ2(JJ2))
00565 104* G2 = XK(LK)*XK(LA)
00566 104* GO TO 3998
00567 104* TM = (T(L)+T(LTA))/2.0
00570 104* CALL D2DIM(TM,CON(14),A(LA),XK(LK),G(LG))
00571 104* GO TO 3999
00572 104* TM = T(LTA)
00573 104* GO TO 3007
00574 104* TM = T(LTA)
00575 104* GO TO 3032
00576 104* G(LG) = 1./(1./G1+1./G2)
00577 104* IF(FLD(3,1,NSQ1(JJ1)),EQ,1) G(LG) = G1*G2
00601 104* 3999 JJ2 = JJ2+1
00602 104* 4000 CONTINUE
00603 104* END
00604 105* IF(FLD(3,1,NSQ1(JJ1)),EQ,0) GO TO 85
00606 106* T1 = T(I)+460.0
00607 107* T2 = T(LTA)+460.0
00610 108* GV = G(LG)*(T1+T2)*T1+T2
00611 109* GO TO 90
00612 110* 85 GV = G(LG)
00613 111* 90 G(I) = G(I)+GV*T(LTA)
00614 112* GSUM = GSUM+GV
00614 113* CHECK FOR LAST CONDUCTOR TO THIS NODE
00615 114* IF(NSQ1(JJ1),GT,0) GO TO 80
00617 115* T2 = DD*(I)+DH*G(I)/GSUM
00620 117* T1 = ABS(T(I)-T2)
00621 118* T(I) = T2
00622 119* IF(RLXA,GT,T1) GO TO 110
00624 120* RLXA = T1
00625 121* N2 = I
00630 123* 110 CONTINUE
00630 124* CON(30) = RLXA
00631 125* SEE IF THE RELAXATION CRITERIA ARE MET
00633 126* IF(KON(7),NE,0) CALL OUTCAL
00635 127* 120 CONTINUE
00637 128* WRITE(6,882)
00641 129* WRITE(6,885) LAX
00644 130* KON(28) = KON(2R)+2
00645 131* KON(37) = N2
00646 132* IF(RLXA,GT,RLXD) GO TO 155
00650 133* CON(30) = RLXD
00651 134* KON(37) = N1
00652 135* CHECK THE ENERGY BALANCE OF THE SYSTEM
00653 136* GOUT = 0.0
00654 137* GIN = 0.0
00655 139* J1 = 0
00656 140* DO 195 I = 1,NNC
00661 141* GIN = GIN+Q(I)
00662 142* J1 = J1+1
00663 143* LTA = FLD(22,14,NSQ1(J1))
00664 144* IF(LTA,LE,NNC) GO TO 175
00666 145* LG = FLD(5,16,NSQ1(J1))
00667 146* IF(FLD(3,1,NSQ1(J1)),EQ,0) GO TO 170
00671 147* T1 = T(I)+460.0
00672 148* T2 = T(LTA)+460.0
00673 149* GOUT = GOUT+G(LG)*T1**4-T2**4
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CINDSS,CINDSS

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00674 150*
00675 151*
00676 152*
00677 153*
00678 154*
00679 155*
00680 156*
00681 157*
00682 158*
00683 159*
00684 160*
00685 161*
00686 162*
00687 163*
00688 164*
00689 165*
00690 166*
00691 167*
00692 168*
00693 169*
00694 170*
00695 171*
00696 172*
00697 173*
00698 174*
00699 175*
00700 176*
00701 177*
00702 178*
00703 179*
00704 180*
00705 181*
00706 182*
00707 183*
00708 184*
00709 185*

      GO TO 175
      GOUT = GOUT+G(LG)*(T(I)-T(LTA))
      CHECK FOR LAST CONDUCTOR TO THIS NODE
      IF(NSQ(I,J1).GT.0) GO TO 165
      175 CONTINUE
      CON(32) = ABS(QIN-GOUT)
      CALL VARBL2
      CON(13) = CON(1)
      CALL OUTCAL
      WRITE(6,882)
      WRITE(6,883) KON(20),CON(32)
      KON(28) = KON(28)+2
      IF(CON(3).GT.CON(1)*1.000001) GO TO 10
      NTH=IE
      NDIM = NLA
      RETURN
      994 WRITE(6,884)
      GO TO 1000
      996 WRITE(6,886) NDIM
      GO TO 1000
      997 WRITE(6,887)
      GO TO 1000
      998 WRITE(6,888)
      GO TO 1000
      999 WRITE(6,889)
      1000 CALL OUTCAL
      CALL EXIT
      882 FORMAT(1H )
      883 FORMAT(10H LOOPCT = I6,10H ENGBAL = E12.5)
      884 FORMAT(46H CINDSS REQUIRES SHORT PSEUDO-COMPUTE SEQUENCE)
      885 FORMAT(35H ITERATION COUNT EXCEEDED, NLOOP = , I10)
      886 FORMAT(18,20H LOCATIONS AVAILABLE)
      887 FORMAT(10H NO DRLXCA)
      888 FORMAT(10H NO ARLXCA)
      889 FORMAT(14H NO LOOP COUNT)
      END

```

END OF UNIVAC 1108 FORTRAN V COMPILATION. 0 *DIAGNOSTIC* MESSAGE(S)
 CINDSS SYMBOLIC
 CINDSS CODE RELOCATABLE

CINOSL,CIN:SL

PIV FOR,* CINOSL,CINOSL
UNIVAC 1108 FORTRAN V ATHENA VERSION: 131K-10D CREATED ON 20 AUG 70
THIS COMPILATION WAS DONE ON 09 JUN 70 AT 13:59:58

SUBROUTINE CINOSL ENTRY POINT 003163

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	003200
0000	*CONST+TEMP	000110
0002	*SIMPLE VAR	000050
0004	*ARRAYS	000000
0005	*BLANK	000000
0006	TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	CONU	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	VARBL1
0023	OUTCAL
0024	DID1WM
0025	PLYAWM
0026	DZD1WM
0027	NONLIN
0030	VARHL2
0031	EXIT
0032	NERK25
0033	NWDUS
0034	NI025
0035	NER105

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C
0012	R	000000	G	0002	R	000020	GSUM
0006	R	000000	H	0002	I	000011	I
0002	I	000042	JJ1	0002	I	000043	JJ2
0017	I	000000	KON	0002	I	000010	K1
0002	I	000036	LE1	0002	I	000037	LE2
0021	I	000007	LSQ2	0002	I	000027	LTA
0021	I	000003	NGT	0002	I	000005	NLA
0021	I	000000	NND	0021	I	000002	MNT
0002	I	000021	NTYPE	0020	K	000002	NX
0011	R	000000	Q	0002	K	000047	QIN
0002	R	000015	RLXA	0002	K	000014	RLXD
0017	R	000000	CON	0017	R	000000	CON
0002	R	000035	GV	0002	R	000035	GV
0002	I	000003	IE1	0002	I	000003	IE1
0002	I	000043	J1	0002	I	000012	J1
0002	I	000044	L	0002	I	000044	L
0002	I	000026	LG	0002	I	000026	LG
0021	I	000005	NAT	0021	I	000005	NAT
0002	I	000001	NN	0002	I	000001	NN
0013	I	000000	NS01	0013	I	000000	NS01
0002	I	000046	OUT	0002	I	000046	OUT
0002	R	000040	R1	0002	R	000040	R1
0002	R	000016	DN	0002	R	000016	DN
0002	R	000032	G2	0002	R	000032	G2
0002	I	000006	JJ	0002	I	000006	JJ
0015	I	000000	K	0015	I	000000	K
0002	I	000007	LAX	0002	I	000007	LAX
0021	I	000006	LS01	0021	I	000006	LS01
0020	I	000000	NDIM	0020	I	000000	NDIM
0002	I	000002	NHC	0002	I	000002	NHC
0020	I	000001	NTH	0020	I	000001	NTH
0002	R	000000	PASS	0002	R	000000	PASS
0002	R	000025	G2	0002	R	000025	G2
0002	R	000030	TM	0002	R	000030	TM
0002	R	000017	DD	0002	R	000017	DD
0002	R	000031	G1	0002	R	000031	G1
0002	I	000004	IE2	0002	I	000004	IE2
0002	I	000013	J2	0002	I	000013	J2
0002	I	000022	LA	0002	I	000022	LA
0002	I	000023	LK	0002	I	000023	LK
0021	I	000004	NCT	0021	I	000004	NCT
0021	I	000001	NNA	0021	I	000001	NNA
0014	I	000000	NS02	0014	I	000000	NS02
0002	I	000045	N2	0002	I	000045	N2
0002	R	000024	Q1	0002	R	000024	Q1
0007	R	000000	T	0007	R	000000	T


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0002 R 000033 T1      0002 R 000034 T2      0020 R 000002 X      0015 R 000000 YK      0001 000133 10L
0001 003137 1000L    0001 002522 110L    0001 002545 115L    0001 002554 120L
0001 002573 130L    0001 002660 135L    0001 002662 140L    0001 002706 150L
0001 000147 160G    0001 000157 165G    0001 002735 165L    0001 003013 170L
0001 000253 1998L   0001 000256 1999L   0001 000173 20L    0001 000261 2000L
0001 000621 2007L  0001 000641 2010L  0001 000644 2015L  0001 000562 2017L
0001 000724 2025L  0001 000764 2030L  0001 000771 2032L  0001 001014 2035L
0001 001040 2042L  0001 001077 2045L  0001 001105 2050L  0001 001150 2055L
0001 001202 2065L  0001 000214 207G  0001 000532 25L    0001 001206 2998L
0001 001270 30L    0001 001237 3000L  0001 001770 3005L  0001 001234 2999L
0001 002022 3015L  0001 002041 3017L  0001 002075 3020L  0001 002016 3010L
0001 002152 3032L  0001 002175 3035L  0001 002201 3040L  0001 002144 3030L
0001 002270 3050L  0001 002334 3055L  0001 002363 3060L  0001 002262 3045L
0001 002373 3998L  0001 002421 3998L  0001 002424 4000L  0001 001273 35L
0001 000332 4012L  0001 000351 4015L  0001 000352 4017L  0001 000331 4010L
0001 000420 4025L  0001 000426 4030L  0001 000454 4035L  0001 000403 4022L
0001 001424 434G  0001 000523 4998L  0001 000527 4999L  0001 000515 4040L
0001 001475 5005L  0001 001503 5010L  0001 001504 5012L  0001 000522 5000L
0001 001542 5020L  0001 001556 5022L  0001 001573 5025L  0001 001525 5017L
0001 001653 5037L  0001 001670 5040L  0001 001335 55L    0001 001637 5035L
0001 001360 60L    0001 001706 6000L  0001 001367 65L    0001 001703 5998L
0001 002732 707G  0001 001412 75L    0001 001706 80L    0001 001402 70L
0000 000001 863F   0000 000007 884F   0000 000020 885F   0000 000000 8A2F
0000 000040 808F  0000 000043 884F  0001 002460 90L    0000 000035 8A7F
0001 003116 997L  0001 003124 998L  0001 003132 999L  0001 003107 996L

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00101 1* SUPROUTINE.CIN.SL
00101 2* STEADY STATE EXECUTION SUBROUTINE FOR SINDA
00101 3* THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED
00101 4* DIFFUSION NODES RECEIVE A SUCCESSIVE POINT ITERATION
00101 5* ARITHMETIC NODES RECEIVE A SUCCESSIVE POINT ITERATION
00101 6* OVER-RELAXATION IS ALLOWED, THE DAMPING FACTORS ARE ADDRESSABLE
00103 7* INCLUDE COMM.LIST
00104 7* COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/S(1) /COND/G(1)
00105 7* COMMON /PCI/NSQ(1) /PC2/HSQ2(1) /KONST/K(1) /ARRAY/A(1)
00106 7* COMMON /FIXCON/KON(1) /XSPACE/NDIM,NTH,X(1)
00107 7* COMMON /DIMENS/ NND,NNA,NHT,NGT,NCT,NAT,LSQ1,LSQ2
00110 7* DIMENSION CON(1),XK(1),NX(1)
00111 7* EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
00112 7* END
00113 8* INCLUDE DEFF.LIST
00113 8* ***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
00113 8* CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEN)
00113 8* CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (DTIMEU)
00113 8* CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEND)
00113 8* CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR*EXPLICIT (CSGFAC)
00113 8* CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
00113 8* CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
00113 8* CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
00113 8* CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEN)
00113 8* CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMP)
00113 8* CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
00113 8* CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
00113 8* CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
00113 8* CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
00113 8* CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
00113 8* CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPCC)

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CINDSL* CINDSL

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00113 8* C CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPCC)
00113 8* C CONTROL CONSTANT 17 IS RESERVED FOR THE C/56 MINIMUM (CSGMIN)
00113 8* C CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
00113 8* C CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
00113 8* C CC20 CONTAINS THE NUMBER OF RELAXATION LOOP'S USED, INTEGER (LOOPCT)
00113 8* C CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEL)
00113 8* C CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
00113 8* C CC23 CONTAINS THE C/56 MAXIMUM (CSGMAX)
00113 8* C CC24 CONTAINS THE C/56 RANGE ALLOWED (CSGRAL)
00113 8* C CC25 CONTAINS THE C/56 RANGE CALCULATED (CSGRCL)
00113 8* C CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
00113 8* C CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DPLXCC)
00113 8* C CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
00113 8* C CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
00113 8* C CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (APLXCC)
00113 8* C CC31 IS INDICATOR, 0=THERMAL SPCS, 1=THERMAL LPCS, 2=GENERAL (LSPCS)
00113 8* C CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENGRAL)
00113 8* C CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)
00113 8* C CC34 CONTAINS THE MOCOPY SWITCH FOR MATRIX USERS (MOCOPY)
00113 8* C CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
00113 8* C CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
00113 8* C CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
00113 8* C CC38 CONTAINS RELATIVE NODE NUMBER OF ATMPCC
00113 8* C CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
00113 8* C CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
00113 8* C CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
00113 8* C CC50 IS NOT USED AT PRESENT
00114 8* C END
00115 9* IF(KON(5),LE.0) GO TO 999
00117 10* IF(CON(9),LE.0) CON(9) = 1.0
00121 11* IF(CON(10),LE.0) CON(10) = 1.0
00123 12* IF(NNA.GT.0.AND.CON(19),LE.0) GO TO 998
00125 13* IF(NND.GT.0.AND.CON(26),LE.0) GO TO 997
00127 14* IF(KOR(31),NE.1) GO TO 994.
00131 15* PASS = -1.0
00132 16* NN = NND+1
00133 17* NNC = NNA+NND
00134 18* IE1 = NTH
00135 19* IE2 = IE1+NNC
00136 20* NLA = NDIM
00137 21* JJ = 2*NNC
00140 22* NTH = NTH+JJ
00141 23* NDIM = NDIM-JJ
00142 24* IF(NDIM.LT.0) GO TO 996
00144 25* CON(1) = CON(13)
00145 26* CON(2) = 0.0
00146 27* CON(14) = CON(13)
00147 28* GO TO 10
00150 29* CON(1) = CON(13)+CON(18)
00151 30* IF(CON(1)-CON(3).GT.0.) CON(1) = CON(3)
00153 31* CON(14) = (CON(1)+CON(13))/2.0
00154 32* CON(2) = CON(1)-CON(13)
00154 33* COMPUTE STEADY STATE TEMPERATURES
00155 34* LAX = KON(5)
00156 35* JJ = 0
00157 36* DO 145 K1 = 1,LAX
00162 37* JJ = JJ+1
00163 38* KON(20) = K1
00163 39* ZERO OUT ALL SOURCE LOCATIONS
00164 40* DO 15 I = 1,NNC

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CINOSL.CINOSL

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00167 15 Q(I) = 0.0
00171 CALL VARBL1
00172 IF(PASS.GE.0.) GO TO 20
00174 CALL OUTCAL
00175 PASS = 1.0
00176 20 J1 = 0
00177 J2 = 1
00200 RLXD = 0.0
00201 RLXA = 0.0
00202 IF(NND.LE.0) GO TO 75
00204 DN = CON(10)
00205 DD = 1.0-DN
00206 DO 70 I = 1,NND
00211 GSUM = 0.0
00212 INCLUDE DUMC*LIST
00213 IF(FLD(1,1),NSQ1(J1+1),EQ.0) GO TO 2000
00215 NTYPE = FLD(10,5),NSG2(J2)
00216 GO TO (1999,1998,1998,1998,1999,1998,1998,1998,1998,1999), NTYPE
00220 1998 J2 = J2+1
00221 1999 J2 = J2+1
00222 2000 CONTINUE
00223 END
00224 INCLUDE VARG*LIST
00225 IF(FLD(4,1),NSQ1(J1+1),EQ.0) GO TO 5000
00226 NTYPE = FLD(10,5),NSQ2(J2)
00227 LA = FLD(5,17),NSQ2(J2)
00230 LK = FLD(22,14),NSQ2(J2)
00231 GO TO (4005,4010,4015,4020,4025,4030,4035,4040,4030), NTYPE
00232 4005 O(I) = XK(LK)+G(I)
00233 4010 O1 = 0.0
00234 4012 CALL DID1WM(T(I),A(LA),XK(LK),Q2)
00236 4015 O1 = 0.0
00237 4017 CALL DID1WM(CON(14),A(LA),XK(LK),Q2)
00240 4020 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00242 4022 J2 = J2+1
00243 LA = FLD(5,17),NSQ2(J2)
00244 LK = FLD(22,14),NSQ2(J2)
00245 GO TO 4017
00246 4025 O1 = XK(LK)*XK(LA)
00250 4030 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00251 J2 = J2+1
00252 LA = FLD(5,17),NSQ2(J2)
00253 LK = FLD(22,14),NSQ2(J2)
00254 Q2 = XK(LK)*XK(LA)
00255 GO TO 4998
00256 4035 CALL DID1WM(CON(14),A(LA),XK(LK),Q1)
00257 4037 J2 = J2+1
00260 LA = FLD(5,17),NSQ2(J2)
00261 LK = FLD(22,14),NSQ2(J2)
00262 GO TO 4012
00263 4040 O1 = XK(LK)*XK(LA)
00264 4998 O(I) = O1+Q2*O(I)
00265 4999 J2 = J2+1
00266 5000 CONTINUE
00267 57*
00270 57*

```

*NEW
**--1

*NEW
*NEW
*NEW
**--3

*NEW
*NEW
**--2

*NEW
*NEW
**--2

*NEW
*NEW
**--2

CINDSL,CINDSL

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00271 57*
00272 58*
00273 59*
00274 60*
00275 61*
00276 62*
00300 62*
00301 62*
00302 62*
00303 62*
00304 62*
00305 62*
00306 62*
00307 62*
00310 62*
00311 62*
00312 62*
00313 62*
00314 62*
00315 62*
00316 62*
00317 62*
00320 62*
00321 62*
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00323 62*
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00346 62*
00347 62*
00350 62*
00351 62*
00352 62*
00353 62*
00354 62*
00355 62*
00356 62*
00357 62*
00360 62*
00361 62*
00362 62*
00364 62*

END
25 J1 = J1+1
LG = FLD(5,16,NSQ1(J1))
LTA = FLD(22,14,NSQ1(J1))
INCLUDE VARG,LIST
CHECK FOR RADIATION CONDUCTOR
IF(FLD(2,1,NSQ1(J1)).EQ.0) GO TO 3000
NTYPE = FLD(0,5,NSQ2(J2))
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
GO TO(2005,2010,2015,2020,2025,2030,2035,2040,2045,2050,2055,
2060,2065), NTYPE
2005 TM = (T(I)+T(LTA))/2.0
2007 CALL DID1WM(TM,A(LA),XK(LK),G(LG))
GO TO 2999
2010 TM = T(I)
GO TO 2007
2015 CALL DID1WM(T(I),A(LA),XK(LK),G1)
2017 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
CALL DID1WM(T(LTA),A(LA),XK(LK),G2)
GO TO 2998
2020 G1 = XK(LK)*XK(LA)
GO TO 2017
2025 CALL DID1WM(T(I),A(LA),XK(LK),G1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
G2 = XK(LK)*XK(LA)
GO TO 2998
2030 TM = (T(I)+T(LTA))/2.0
2032 CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),G(LG))
GO TO 2999
2035 TM = T(I)
GO TO 2032
2040 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
2042 J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),G2)
GO TO 2998
2045 G1 = XK(LK)*XK(LA)
GO TO 2042
2050 CALL PLYAWM(A(LA),T(I),A(LA+1),XK(LK),G1)
J2 = J2+1
LA = FLD(5,17,NSQ2(J2))
LK = FLD(22,14,NSQ2(J2))
G2 = XK(LK)*XK(LA)
GO TO 2998
2055 TM = (T(I)+T(LTA))/2.0
CALL D2D1WM(TM,CON(14),A(LA),XK(LK),G(LG))
GO TO 2999
2060 TM = T(LTA)
GO TO 2007
2065 TM = T(LTA)
GO TO 2032
2998 G(LG) = 1./(1./G1+1./G2)
IF(FLD(3,1,NSQ1(J1)).EQ.1) G(LG) = G1*G2
2999 J2 = J2+1

```

*NEW
*NEW
***-3

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

*NEW
*NEW
***-2

CINDSL,CINDSL

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00365 3000 CONTINUE
00366 END
00367 IF(FLD(3,1,NSQ1(JJ1)),EQ,0) GO TO 30
00371 T1 = T(I)+460.0
00372 T2 = T(LTA)+460.0
00373 GV = G(LG)*(T1+T1+T2*T2)*(T1+T2)
00374 GO TO 35
00375 30 GV = G(LG)
00376 35 GSUM = GSUM+GV
00377 G(I) = G(I)+GV*(LTA)
00377 CHECK FOR LAST CONDUCTOR TO THIS NODE
00400 IF(NSQ1(JJ1),GT,0) GO TO 25
00402 T2 = DD*(T1)+DN*G(I)/GSUM
00403 OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
00403 T1 = ABS(T(I)-T2)
00403 STORE THE NEW TEMPERATURES AND EXTRAPOLATION FACTORS
00404 GO TO(65,60,55),JJ
00405 55 LE1 = IE1+I
00406 LE2 = IE2+I
00407 R1 = T2-T(I)
00410 X(LE1) = T(I)
00411 X(LE2) = R1/(R1-X(LE2))
00412 GO TO 65
00413 60 LE2 = IE2+I
00414 X(LE2) = T2-T(I)
00415 65 T(I) = T2
00415 SAVE THE MAXIMUM DIFFUSION RELAXATION CHANGE
00416 IF(RLXD,GT,T1) GO TO 70
00420 RLXD = T1
00421 N1 = I
00422 70 CONTINUE
00424 CON(27) = RLXD
00425 IF(NRA,LE,0) GO TO 130
00427 DN = CON(9)
00430 DD = 1.0-DN
00431 JJ1 = J1
00432 JJ2 = J2
00432 DO SUCCESSIVE POINT ITERATION ON ARITHMETIC NODES
00433 DO 125 I = NN,NNC
00436 GSUM = 0.0
00437 L = I
00440 INCLUDE VRQ2,L1ST
00441 IF(FLD(4,1,NSQ1(JJ1+1)),EQ,0) GO TO 6000
00443 NTYPE = FLD(0,5,NSQ2(JJ2))
00444 LA = FLD(5,17,NSQ2(JJ2))
00445 LK = FLD(22,14,NSQ2(JJ2))
00446 GO TO (5005,5010,5015,5020,5025,5030,5035,5040,5030), NTYPE
00447 5005 G(L) = XK(LK)+G(L)
00450 GO TO 5999
00451 5010 O1 = 0.0
00452 5012 CALL DID1WM(T(L),A(LA),XK(LK),O2)
00453 GO TO 5998
00454 5015 O1 = 0.0
00455 5017 CALL DID1WM(CON(14),A(LA),XK(LK),O2)
00456 GO TO 5998
00457 5020 CALL DID1WM(CON(14),A(LA),XK(LK),O1)
00460 JJ2 = JJ2+1
00461 LA = FLD(5,17,NSQ2(JJ2))
00462 LK = FLD(22,14,NSQ2(JJ2))
00463 GO TO 5017

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*NEW
*NEW
*NEW
**3

*NEW
*NEW
**2

CINDSL,CINOSL

00464	102*	5025	01 = XK(LK)*XK(LA)	*NEW
00465	102*	GO TO 5022		*NEW
00466	102*	5030	CALL DID1WM(CON(14),A(LA),XK(LK),01)	**--2
00467	102*	JJ2 = JJ2+1		
00470	102*	LA = FLD(5,17,NSQ2(JJ2))		
00471	102*	LK = FLD(22,14,NSQ2(JJ2))		
00472	102*	02 = XK(LK)*XK(LA)		
00473	102*	GO TO 5998		
00474	102*	5035	CALL DID1WM(CON(14),A(LA),XK(LK),01)	*NEW
00475	102*	5037	JJ2 = JJ2+1	*NEW
00476	102*	LA = FLD(5,17,NSQ2(JJ2))		**--2
00477	102*	LK = FLD(22,14,NSQ2(JJ2))		
00500	102*	GO TO 5012		
00501	102*	5040	01 = XK(LK)*XK(LA)	
00502	102*	GO TO 5037		
00503	102*	5998	Q(L) = Q1+Q2+Q(L)	
00504	102*	5999	JJ2 = JJ2+1	
00505	102*	6000	CONTINUE	
00506	102*	END		
00507	103*	80	JJ1 = JJ1+1	
00510	104*	LG = FLD(5,16,HSQ1(JJ1))		
00511	105*	LTA = FLD(22,14,NSQ1(JJ1))		
00512	106*	INCLUDE VRG2,LIST		
00512	107*	CHECK FOR RADIATION CONDUCTOR		
00513	107*	IF(FLD(2,1,NSQ1(JJ1)).EQ.0) GO TO 4000		
00516	107*	NTYPE = FLD(10,5,NSQ2(JJ2))		*NEW
00517	107*	LA = FLD(5,17,NSQ2(JJ2))		*NEW
00517	107*	LK = FLD(22,14,NSQ2(JJ2))		**--3
00520	107*	GO TO(3005,3010,3015,3020,3025,3030,3035,3040,3045,3050,3055,		
00521	107*	3060,3065), NTYPE		
00522	107*	TM = (T(L)+T(LTA))/2.0		
00523	107*	3005	CALL DID1WM(TM,A(LA),XK(LK),6(LG))	
00524	107*	GO TO 3999		
00525	107*	3010	TM = T(L)	
00526	107*	GO TO 3007		
00527	107*	3015	CALL DID1WM(T(L),A(LA),XK(LK),61)	
00530	107*	3017	JJ2 = JJ2+1	
00531	107*	LA = FLD(5,17,NSQ2(JJ2))		
00532	107*	LK = FLD(22,14,NSQ2(JJ2))		
00533	107*	CALL DID1WM(T(LTA),A(LA),XK(LK),62)		
00534	107*	GO TO 3998		
00535	107*	3020	01 = XK(LK)*XK(LA)	
00536	107*	GO TO 3017		
00537	107*	3025	CALL DID1WM(T(L),A(LA),XK(LK),61)	
00540	107*	JJ2 = JJ2+1		
00541	107*	LA = FLD(5,17,NSQ2(JJ2))		
00542	107*	LK = FLD(22,14,NSQ2(JJ2))		
00543	107*	62 = XK(LK)*XK(LA)		
00544	107*	GO TO 3998		
00545	107*	3030	TM = (T(L)+T(LTA))/2.0	
00546	107*	3032	CALL PLYAWM(A(LA),TM,A(LA+1),XK(LK),6(LG))	
00547	107*	GO TO 3999		
00550	107*	3035	TM = T(L)	
00551	107*	GO TO 3032		
00552	107*	3040	CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),61)	
00553	107*	3042	JJ2 = JJ2+1	
00554	107*	LA = FLD(5,17,NSQ2(JJ2))		
00555	107*	LK = FLD(22,14,NSQ2(JJ2))		
00556	107*	CALL PLYAWM(A(LA),T(LTA),A(LA+1),XK(LK),62)		
		GO TO 3998		

*NEW
*NEW
**=2

CINDSL,CINDSL

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00557 107* 3045 G1 = XK(LK)*XK(LA)
00560 107* 60 TO 3042
00561 107* 3050 CALL PLYAWM(A(LA),T(L),A(LA+1),XK(LK),G1)
00562 107* JJ2 = JJ2+1
00563 107* LA = FLD(15,17,NSQ2(JJ2))
00564 107* LK = FLD(22,14,NSQ2(JJ2))
00565 107* G2 = XK(LK)*XK(LA)
00566 107* 60 TO 3998
00567 107* 3055 TM = (T(L)+T(LTA))/2.0
00570 107* CALL D2D1*(M(TM,CON(14),A(LA),XK(LK),G(LG))
00571 107* 60 TO 3999
00572 107* 3060 TM = T(LTA)
00573 107* 60 TO 3007
00574 107* 3065 TM = T(LTA)
00575 107* 60 TO 3032
00576 107* 3998 G(LG) = 1./(1./G1+1./G2)
00577 107* 3999 JJ2 = JJ2+1
00602 107* 4000 CONTINUE
00603 107* END
00604 108* IF(FLD(3,1,NSQ1(JJ1)),EG.0) 60 TO 85
00606 109* T1 = T(I)+460.0
00607 110* T2 = T(LTA)+460.0
00610 111* GV = G(LG)*(T1+T2)*T1+T2)
00611 112* 60 TO 90
00612 113* 85 GV = G(LG)
00613 114* 90 Q(I) = Q(I)+GV*T(LTA)
00614 115* GSUM = GSUM+GV
00615 117* C CHECK FOR LAST CONDUCTOR TO THIS NODE
00617 118* IF(NSQ1(JJ1),GT.0) 60 TO 80
00620 120* T2 = DO*(T(I)+DM*Q(I))/GSUM
00621 121* T1 = ABS(T(I)-T2)
00622 122* C STORE THE NEW TEMPERATURES AND EXTRAPOLATION FACTORS
00623 123* 60 TO(120,115,110),JJ
00624 124* 110 LE1 = IE1+I
00625 125* LE2 = IE2+I
00626 126* R1 = T2-T(I)
00627 127* X(LE1) = T(I)
00630 128* X(LE2) = R1/(R1-X(LE2))
00631 129* 60 TO 120
00632 130* 115 LE2 = IE2+I
00633 131* 120 T(I) = T2
00635 132* IF(RLXA,GT.T1) 60 TO 125
00636 133* RLXA = T1
00637 134* N2 = I
00641 135* 125 CONTINUE
00642 137* CON(30) = RLXA
00644 138* C SEE IF THE RELAXATION CRITERIA ARE MET
00646 139* JJ = 0
00647 140* DO 135 I = 1,NMC
00652 141* LE2 = IE2+I
00653 143* C SEE IF THE EXTRAPOLATION CRITERIA ARE MET
00654 144* IF(X(LE2),GE.0.) 60 TO 135
00657 145* IF(X(LF2),LT.-R.) X(LE2) = -8.
00660 146* LE1 = IE1+I
00661 147* T(I) = X(LE2)*X(LE1)+(1.0-X(LE2))*T(I)
00661 147* 135 CONTINUE
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CINDSL,CINDSL

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00663 148*
00665 149*
00667 150* WRITE(6,882)
00671 151* WRITE(6,885) LAX
00674 152* KON(28) = KON(2R)+2
00675 153* KON(37) = N2
00676 154* IF(RLXA.GT.RLXD) GO TO 155
00700 155* CON(30) = RLXD
00701 156* KON(37) = N1
00701 157* CHECK THE ENERGY BALANCE OF THE SYSTEM
00702 158* CALL NORLIN
00703 159* GOUT = 0.0
00704 160* QIN = 0.0
00705 161* J1 = 0
00706 162* DO 195 I = 1,NNC
00711 163* GIN = QIN+Q(I)
00712 164* J1 = J1+1
00713 165* LTA = FLD(22,14,NSQI(J1))
00714 166* IF(LTA.LE.NNC) GO TO 175
00716 167* LG = FLD(15,16,NSQI(J1))
00717 168* IF(FLD(3,1,NSQI(J1)).EQ.0) GO TO 170
00721 169* T1 = T(I)+460.0
00722 170* T2 = T(LTA)+460.0
00723 171* GOUT = GOUT+G(LG)*(T1**4-T2**4)
00724 172* GO TO 175
00725 173* GO TO 175
00725 174* GOUT = GOUT+G(LG)*(T(I)-T(LTA))
00726 175* CHECK FOR LAST CONDUCTOR TO THIS NODE
00730 176* 175 IF(NSQI(J1).GT.0) GO TO 165
00732 177* 195 CONTINUE
00733 178* CON(32) = ABS(QIN-GOUT)
00734 179* CALL VARBL2
00735 180* CON(13) = CON(1)
00736 181* CALL OUTCAL
00740 182* WRITE(6,882)
00744 183* WRITE(6,883) KON(20),CON(32)
00745 184* KON(28) = KON(28)+2
00747 185* IF(CON(3).GT.CON(1)*1.000001) GO TO 5
00750 186* NTH = IEI
00751 187* NDIM = NLA
00752 188* RETURN
00754 189* 994 WRITE(6,884)
00755 190* GO TO 1000
00760 191* 996 WRITE(6,886) NDIM
00761 192* GO TO 1000
00763 193* 997 WRITE(6,887)
00764 194* GO TO 1000
00766 195* 998 WRITE(6,888)
00771 197* 999 WRITE(6,889)
00772 198* 1000 CALL OUTCAL
00773 199* CALL EXIT
00774 200* 882 FORMAT(1H )
00775 201* 883 FORMAT(10H LOOPCT = 16,10H ENGBAL = E12.5)
00776 202* 884 FORMAT(45H CINDSL REQUIRES LONG PSEUDO-COMPUTE SEQUENCE)
00777 203* 885 FORMAT(35H ITERATION COUNT EXCEEDED, LOOPCT =, 110)
01000 204* 886 FORMAT(10H NO (RLXCA)
01001 205* 887 FORMAT(10H NO (RLXCA)
01002 206* 888 FORMAT(10H NO (RLXCA)
01003 207* 889 FORMAT(14H NO LOOP COUNT)
          END

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JI FOR** CINDSM UNIVAC 1108 FORTRAN V ATHENA VERSION 131K-10D CREATED ON 20 AUG 70 THIS COMPILATION WAS DONE ON 09 JUN 70 AT 22:46:55

SUBROUTINE CINDSM ENTRY POINT 001175

STORAGE USED (BLOCK, NAME, LENGTH)

0001	*CODE	001211
0000	*CONST+TEMP	000067
0002	*SIMPLE VAR	000044
0004	*ARRAYS	000600
0005	*BLANK	000000
0006	*TITLE	000001
0007	TEMP	000001
0010	CAP	000001
0011	SOURCE	000001
0012	COND	000001
0013	PC1	000001
0014	PC2	000001
0015	KONST	000001
0016	ARRAY	000001
0017	FIXCON	000001
0020	XSPACE	000003
0021	DIMENS	000010

EXTERNAL REFERENCES (BLOCK, NAME)

0022	NONLIN
0023	OUTCAL
0024	VARBL2
0025	EXIT
0026	NERR2\$
0027	NWDUS
0030	NI02\$
0031	NER10\$

STORAGE ASSIGNMENT FOR VARIABLES (BLOCK, TYPE, RELATIVE LOCATION, NAME)

0016	R	000000	A	0010	R	000000	C	0017	R	000000	CON
0012	R	000000	G	0002	R	000031	GSUM	0006	R	000000	H
0002	I	000005	IE2	0002	I	000006	IE3	0002	I	000007	IE4
0015	I	000000	K	0002	I	000037	KK	0017	I	000000	KON
0002	I	000036	LE1	0002	I	000023	LE2	0002	I	000023	LE3
0002	I	000000	LPASS	0021	I	000006	LS01	0021	I	000007	LS02
0021	I	000004	NCT	0020	I	000000	NDIM	0021	I	000003	NGT
0021	I	000001	NNA	0002	I	000001	NHC	0021	I	000000	NND
0014	I	000000	NS02	0020	I	000001	NTH	0020	I	000002	NX
0011	R	000000	G	0002	R	000020	GIN	0002	R	000043	GOUT
0002	R	000030	T2	0002	R	000034	RLXD	0002	R	000040	R1
0002	R	000030	TL	0020	R	000002	X	0015	I	000000	KK
0001	I	000116	10L	0001	I	001152	1000L	0001	I	000571	105L
0001	I	000716	130L	0001	I	000727	135L	0001	I	001905	140L
0001	I	001041	155L	0001	I	001053	160L	0001	I	000177	2046
0001	I	000157	25L	0001	I	000332	2516	0001	I	000162	30L
0002	R	000013	DELXXX	0002	R	000016	DAMP	0002	R	000012	XXX
0002	I	000004	IE1	0002	I	000022	I	0021	I	000002	NNT
0002	I	000021	J1	0002	I	000002	JJ	0002	I	000041	N1
0002	I	000010	LAXFAC	0002	I	000032	K1	0002	R	000035	GSUM
0002	I	000024	LG	0002	I	000025	LE4	0007	R	000000	T
0002	I	000005	NAT	0002	I	000026	LTA	0001	I	000644	120L
0002	I	000011	PLAX	0002	I	000003	NLA	0001	I	000712	128L
0013	I	000000	NS01	0021	I	000002	NNT	0001	I	000135	15L
0002	I	000042	N2	0002	I	000002	N1	0001	I	000320	2126
0002	R	000015	RELAX	0002	R	000035	GSUM	0001	I	000507	3206
0002	R	000027	T1	0002	R	000012	XXX	0002	R	000014	XXXDUM
0002	R	000014	XXXDUM	0001	I	000712	128L	0001	I	000135	15L
0001	I	000712	128L	0001	I	000164	145L	0001	I	000320	2126
0001	I	000135	15L	0001	I	000300	2316	0001	I	000507	3206
0001	I	000320	2126	0001	I	000177	2046				
0001	I	000507	3206								

0001	000211 35L	0001	000623 3516	0001	000663 3709	0001	000261 40L	0001	000727 410G
0001	000265 45L	0001	000074 5L	0001	000336 60L	0001	000410 65L	0001	000837 70L
0001	000446 75L	0001	000455 80L	0001	000470 85L	0000	000000 884F	0000	000006 885F
0000	000013 886F	0000	000016 887F	0000	000021 888F	0000	000026 889F	0001	000535 90L
0001	000557 95L	0001	001114 995L	0001	001123 996L	0001	001131 997L	0001	001137 998L
0001	001145 999L								

CINDSM

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SUBROUTINE CINDSM
STEADY STATE EXECUTION SUBROUTINE FOR SINDA, FORTRAN V
THE LONG PSEUDO-COMPUTE SEQUENCE IS REQUIRED
DIFFUSION NODES RECEIVE A SUCCESSIVE POINT ITERATION
ARITHMETIC NODES RECEIVE A SUCCESSIVE POINT ITERATION
A QUASI-LINEAR BLOCK ITERATION IS PERFORMED ON RADIATION
INCLUDE COMM,LIST
COMMON /TITLE/H(1) /TEMP/T(1) /CAP/C(1) /SOURCE/G(1) /COND/G(1)
COMMON /PCI/NSQ(11) /PC2/NSQ2(1) /KONST/K(1) /ARRAY/A(1)
COMMON /FIXCON/KON(1) /XSPACE/NDIM,NTH,X(1)
COMMON /DIMENS/ NND,NNA,NNT,NGT,NCT,NAT,LS01,LS02
DIMENSION CON(1),XK(1),NX(1)
EQUIVALENCE (KON(1),CON(1)),(K(1),XK(1)),(X(1),NX(1))
END
INCLUDE DEFF,LIST
CC49 CONTAINS THE NO. OF ITER. BEFORE RAD,COND. LINEAR. (LAXFAC)
C***** CONTROL CONSTANT DEFINITIONS AND NAMES *****
C CONTROL CONSTANT 1 CONTAINS THE NEW PROBLEM TIME (TIMEIN)
C CONTROL CONSTANT 2 CONTAINS THE TIME STEP USED (TIMEU)
C CONTROL CONSTANT 3 CONTAINS THE PROBLEM STOP TIME (TIMEO)
C CONTROL CONSTANT 4 CONTAINS THE TIME STEP FACTOR,EXPLICIT (CSGFAC)
CC5 IS THE INPUT NUMBER OF ITERATION DO LOOPS, INTEGER (NLOOP)
CC6 CONTAINS THE DIFFUSION TEMPERATURE CHANGE ALLOWED (DTMPCA)
CC7 CONTAINS THE OUTPUT EACH ITERATION SWITCH (OPEITR)
CC8 CONTAINS THE MAXIMUM ALLOWED TIME STEP (DTIMEH)
CC9 CONTAINS THE NEW ARITHMETIC TEMP. DAMPING FACTOR (DAMP)
CC10 CONTAINS THE NEW DIFFUSION TEMP. DAMPING FACTOR (DAMPD)
CC11 CONTAINS THE MAXIMUM ALLOWED ARITHMETIC TEMP. CHANGE (ATMPCA)
CC12 CONTAINS THE BACKUP SWITCH CHECKED AFTER VARIABLES (BACKUP)
CC13 CONTAINS THE PRESENT TIME OR PROBLEM START TIME (TIMEO)
CC14 CONTAINS THE MEAN TIME BETWEEN AN ITERATION (TIMEM)
CC15 CONTAINS THE DIFFUSION TEMPERATURE CHANGE CALCULATED (DTMPC)
CC16 CONTAINS ARITHMETIC TEMPERATURE CHANGE CALCULATED (ATMPC)
CONTROL CONSTANT 17 IS RESERVED FOR THE C/SG MINIMUM (CSGMIN)
CONTROL CONSTANT 18 CONTAINS THE OUTPUT INTERVAL (OUTPUT)
CC19 CONTAINS THE ARITHMETIC RELAXATION CRITERIA ALLOWED (ARLXCA)
CC20 CONTAINS THE NUMBER OF RELAXATION LOOPS USED,INTEGER (LOOPCT)
CC21 CONTAINS THE MINIMUM ALLOWED TIME STEP (DTIMEI)
CC22 IS FOR THE INPUT TIME STEP IMPLICIT (DTIMEI)
CC23 CONTAINS THE C/SG MAXIMUM (CSGMAX)
CC24 CONTAINS THE C/SG RANGE ALLOWED (CSGRAL)
CC25 CONTAINS THE C/SG RANGE CALCULATED (CSGRCL)
CC26 CONTAINS THE DIFFUSION RELAXATION CRITERIA ALLOWED (DRLXCA)
CC27 CONTAINS THE DIFFUSION RELAXATION CHANGE CALCULATED (DRLXCC)
CC28 CONTAINS THE LINE COUNTER, INTEGER (LINECT)
CC29 CONTAINS THE PAGE COUNTER, INTEGER (PAGECT)
CC30 CONTAINS ARITHMETIC RELAXATION CHANGE CALCULATED (ARLXCC)
CC31 IS INDICATOR, 0=THERMAL SPCS,1=THERMAL LPCS,2=GENERAL (LSPCS)
CC32 CONTAINS THE ENERGY BALANCE OF THE SYSTEM, IN - OUT (ENBRAL)
CC33 CONTAINS THE DESIRED ENERGY BALANCE, USER INPUT (BALENG)

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CINDSM

(NOCOPY)

CC34 CONTAINS THE NOCOPY SWITCH FOR MATRIX USERS
 CC35 CONTAINS RELATIVE NODE NUMBER OF CSGMIN
 CC36 CONTAINS RELATIVE NODE NUMBER OF DTMPC
 CC37 CONTAINS RELATIVE NODE NUMBER OF ARLXCC
 CC38 CONTAINS RELATIVE NODE NUMBER OF ATWPC
 CC39-40-41-42-43 CONTAIN DUMMY INTEGER CONSTANTS (I-J-K-L-MTEST)
 CC44-45-46-47-48 CONTAIN DUMMY FLOATING CONSTANTS (R-S-T-U-VTEST)
 CC49 IS THE QUASI-LINEARIZATION INTERVAL FOR CINDSM (LAXFAC)
 CC50 IS NOT USED AT PRESENT

9* 00113
 9* 00114
 9* 00115
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 9* 00119
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 9* 00205
 9* 00206
 9* 00207

IF(KON(5).LE.0) GO TO 999
 IF(CON(9).LE.0.0) CON(9) = 1.0
 IF(CON(10).LE.0.0) CON(10) = 1.0
 IF(KON(11).NE.1) GO TO 998
 IF(CON(13).LE.0.0) GO TO 997
 IF(KON(49).LE.0) GO TO 996
 LPASS = -1
 NNC = NNA+RND
 JJ = 3*NMC+NGT
 NLA = NDIM
 NDM = NDIM-JJ
 IF(NDIM.LT.0) GO TO 995
 IE1 = NTH
 NTH = NTH+JJ
 IE2 = IE1+MNC
 IE3 = IE2+NMC
 IE4 = IE3+NMC
 CON(1) = CON(13)
 CON(2) = 0.0
 CON(14) = CON(13)
 GO TO 10
 5 CON(1) = CON(13)+CON(18)
 IF(CON(1)-CON(3).GT.0.0) CON(1) = CON(3)
 CON(14) = (CON(1)+CON(13))/2.0
 CON(2) = CON(1)-CON(13)
 LPASS = 0
 10 CONTINUE
 KON(20) = 0
 LAXFAC = KON(49)
 NLAX = KON(5)/LAXFAC
 XXX = 0.05
 DELXXX = 0.05/FLOAT(NLAX)
 XXXDUM = 0.001
 EVALUATE NONLINEARITIES
 15 CALL NONLIN
 RELAX = XXX
 DAMP = CON(10)
 RLX = 1.0-DAMP
 IF(LPASS.GT.-1) GO TO 25
 CALL OUTCAL
 LPASS = 1
 GO TO 30
 25 LPASS = LPASS+1
 30 CONTINUE
 C SUM ALL SOURCES, SAVE ALL TEMPERATURES AND LINEARIZE RADIATION
 QIN = 0.0
 J1 = 0
 DO 50 I = 1,NMC
 QIN = QIN+Q(I)
 LE3 = IE3+I

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CINDSM

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60* 00210 X(LE3) = T(I)
61* 00211 J1 = J1+1
62* 00212 LG = FLD(5,16,NSO1(J1))
63* 00213 LE4 = IE4+LG
64* 00214 CHECK FOR RADIATION CONDUCTOR
65* 00215 IF(FLD(3,1,NSO1(J1)).EQ.0) GO TO 40
66* 00216 LTA = FLD(22,14,NSO1(J1))
67* 00217 T1 = T(I)+460.0
68* 00220 T2 = T(LTA)+460.0
69* 00221 X(LE4) = G(LG)*(T1+T2)*T1+T2
70* 00222 GO TO 45
71* 00223 40 X(LE4) = G(LG)
72* 00224 CHECK FOR LAST CONDUCTOR INTO THIS NODE
73* 00225 45 IF(NSO1(J1).GT.0) GO TO 35
74* 00226 50 CONTINUE
75* 00227 SWITCH CONDUCTOR VALUES WITH FOURTH EXTRA ARRAY IN X
76* 00230 DO 55 I = 1,NGT
77* 00233 LE4 = IE4+I
78* 00234 GSUM = X(LE4)
79* 00235 X(LE4) = G(I)
80* 00236 G(I) = GSUM
81* 00237 55 PERFORM LAXFAC ITERATIONS USING THE LINEARIZED CONDUCTORS
82* 00240 JJ = 0
83* 00241 DO 100 K1 = 1,LAXFAC
84* 00244 KK = K1
85* 00245 JJ = JJ+1
86* 00246 J1 = 0
87* 00247 RLXD = 0.0
88* 00248 DO SUCCESSIVE POINT ITERATION ON ALL CALCULATED NODES
89* 00250 DO 85 I = 1,NNC
90* 00253 GSUM = 0.0
91* 00254 GSUM = G(I)
92* 00255 J1 = J1+1
93* 00256 LG = FLD(5,16,NSO1(J1))
94* 00257 LTA = FLD(22,14,NSO1(J1))
95* 00260 GSUM = GSUM+G(LG)
96* 00261 GSUM = GSUM+G(LG)*T(LTA)
97* 00262 CHECK FOR LAST CONDUCTOR INTO THIS NODE
98* 00263 IF(NSO1(J1).GT.0) GO TO 60
99* 00264 T2 = GSUM/GSUM
100* 00264 60 OBTAIN THE CALCULATED TEMPERATURE DIFFERENCE
101* 00265 T1 = ABS(T1)-T2
102* 00265 STORE THE NEW TEMPERATURES AND EXTRAPOLATION FACTORS
103* 00266 GO TO (80,75,65), JJ
104* 00267 LE1 = IE1+I
105* 00270 LE2 = IE2+I
106* 00271 R1 = T2-T(I)
107* 00272 X(LE1) = T(I)
108* 00273 IF(R1-X(LE2)).NE.0.0) GO TO 70
109* 00275 X(LE2) = 1.0
110* 00276 GO TO 80
111* 00277 70 X(LE2) = R1/(R1-X(LE2))
112* 00300 GO TO 80
113* 00301 75 LE2 = IE2+I
114* 00302 X(LE2) = T2-T(I)
115* 00303 T1 = T2
116* 00304 80 SAVE THE MAXIMUM TEMPERATURE CHANGE AND LOCATION
117* 00304 IF(RLXD.GT.T1) GO TO 85
118* 00306 RLXD = T1
119* 00307 N1 = I

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00310 120*      85 CONTINUE
00311 121*      SEE IF THE RELAXATION CRITERIA WAS MET
00312 122*      IF (RLXD.LE.RELAX) GO TO 105
00313 123*      IF (JJ.LE.2) GO TO 95
00314 124*      JJ = 0
00315 125*      DO 90 I = 1,NNC
00316 126*      LE1 = IE1+I
00317 127*      LE2 = IE2+I
00318 128*      SEE IF THE EXTRAPOLATION CRITERIA ARE MET
00319 129*      IF (X(LE2).GE.0.0.OR.ABS(X(LE2)*T(I)-X(LE1)).GE.RLXD) GO TO 90
00320 130*      T(I) = X(LE2)*X(LE1)+(1.0-X(LE2))*T(I)
00321 131*      90 CONTINUE
00322 132*      CON(30) = RLXD
00323 133*      KON(37) = N1
00324 134*      IF (KON(7).NE.0) CALL OUTCAL
00325 135*      100 CONTINUE
00326 136*      105 KON(20) = KON(20)+KK
00327 137*      STORE CONDUCTANCE VALUES BACK IN THE G ARRAY
00328 138*      DO 110 I = 1,NGT
00329 139*      LE4 = IE4+I
00330 140*      G(I) = X(LE4)
00331 141*      CHECK IF THE INITIAL NLAX ITERATIONS HAVE BEEN PERFORMED
00332 142*      IF (LPASS.GE.NLAX) GO TO 120
00333 143*      THE NLAX INITIAL ITERATIONS HAVE NOT BEEN PERFORMED,
00334 144*      APPLY DAMPING FACTOR AND REDUCE RELAX BY ONLY 0.005/NLAX
00335 145*      DO 115 I = 1,NNC
00336 146*      LE3 = IE3+I
00337 147*      T(I) = DAMP*T(I)+RLX*X(LE3)
00338 148*      XXX = XXX-DELXXX
00339 149*      GO TO 15
00340 150*      AFTER THE NLAX INITIAL ITERATIONS, REDUCE RELAX TO 0.001
00341 151*      120 CONTINUE
00342 152*      N2 = 0
00343 153*      CHECK TO SEE IF MAXIMUM NUMBER OF ITERATIONS HAS BEEN EXCEEDED
00344 154*      IF (KON(120).GE.KON(5)) GO TO 130
00345 155*      HAS NOT BEEN EXCEEDED, REDUCE RELAX TO 0.001
00346 156*      N2 = 1
00347 157*      XXX = XXXDUM
00348 158*      GSUM = 0.0
00349 159*      OBTAIN THE MAXIMUM TEMPERATURE CHANGE
00350 160*      DO 125 I = 1,NNC
00351 161*      LE3 = IE3+I
00352 162*      GOUT = ABS(T(I)-X(LE3))
00353 163*      IF (GOUT.GT.GSUM) GSUM = GOUT
00354 164*      125 CONTINUE
00355 165*      IF THE MAXIMUM TEMPERATURE CHANGE OVER THE QUASI-INTERVAL EXCEEDS
00356 166*      RELAX, REDUCE CRITERIA AND PERFORM MORE ITERATIONS
00357 167*      IF (GSUM.LE.RELAX) GO TO 130
00358 168*      REDUCE LAXFAC TO THE DIFFERENCE BETWEEN NLOOP AND LOOPCT, SO
00359 169*      THAT NLOOP REMAINS THE MAXIMUM NUMBER OF ITERATIONS POSSIBLE
00360 170*      LAXFAC = KON(5)-KON(120)
00361 171*      GO TO 15
00362 172*      CHECK THE ENERGY BALANCE OF THE SYSTEM
00363 173*      NLOOP HAS BEEN EXCEEDED
00364 174*      N2 = N2+1
00365 175*      GOUT = 0.0
00366 176*      J1 = 0
00367 177*      DETERMINE NET HEAT FLOW TO BOUNDARY NODES
00368 178*      DO 150 I = 1,NHC
00369 179*      135 J1 = J1+1

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00413 160* LTA = FLD(22,14,NSO1(J1))
00414 161* CHECK FOR BOUNDARY NODE
00415 162* IF(LTA.LE.NNC) GO TO 145
00416 163* LG = FLD(5,16,NSO1(J1))
00417 164* CHECK FOR RADIATION CONDUCTOR
00418 165* IF(FLD(3,1,NSO1(J1)).EQ.0) GO TO 140
00419 166* T1 = T(I)+460.0
00420 167* T2 = T(LTA)+460.0
00421 168* GOUT = GOUT+(LG)*(T1**4-T2**4)
00422 169* GO TO 145
00423 170* GOUT = GOUT+(LG)*(T(I)-T(LTA))
00424 171* CHECK FOR LAST CONDUCTOR INTO THIS NODE
00425 172* IF(NSO1(J1).GT.0) GO TO 135
00426 173* 150 CONTINUE
00427 174* CON(32) = ABS(QIN-GOUT)
00428 175* GO TO (160,155), N2
00429 176* 155 IF(CON(32).LE.CON(33)) GO TO 160
00430 177* NLOOP HAS NOT BEEN EXCEEDED, MAXIMUM TEMPERATURE CHANGE IS LESS
00431 178* THAN OR EQUAL TO RELAXATION CRITERIA BUT ENGBAL IS GREATER THAN
00432 179* BALENG. MORE ITERATIONS WILL BE PERFORMED WITH A TIGHTER CRITERIA
00433 180* XXXDUM = XXXDUM/5.0
00434 181* XXX = XXXDUM
00435 182* GO TO 128
00436 183* EITHER NLOOP HAS BEEN EXCEEDED OR ENGBAL IS LESS THAN OR
00437 184* EQUAL TO BALENG. IN EITHER CASE PRINT LOOPCT AND ENGBAL,
00438 185* INCREMENT TIME AND PROCEED WITH THE PROBLEM.
00439 186* 160 WRITE(6,884) KON(20),CON(32)
00440 187* KON(28) = KON(28)+2
00441 188* CON(30) = RLXD
00442 189* KON(37) = N1
00443 190* CALL VARBL2
00444 191* CON(13) = CON(1)
00445 192* CALL OUTCAL
00446 193* TEST TO SEE IF TIMEND HAS BEEN REACHED
00447 194* IF(CON(3).GT.CON(1)*1.000001) GO TO 5
00448 195* NTH = IEI
00449 196* NDIM = NLA
00450 197* RETURN
00451 198* 995 WRITE(6,885) NDIM
00452 199* GO TO 1000
00453 200* 996 WRITE(6,886)
00454 201* GO TO 1000
00455 202* 997 WRITE(6,887)
00456 203* GO TO 1000
00457 204* 998 WRITE(6,888)
00458 205* GO TO 1000
00459 206* 999 WRITE(6,889)
00460 207* GO TO 1000
00461 208* 1000 CALL OUTCAL
00462 209* CALL EXIT
00463 210* 884 FORMAT(10H0LOOPCT = I6,10H ENGBAL = E12.5)
00464 211* 885 FORMAT(18,20H LOCATIONS AVAILABLE)
00465 212* 886 FORMAT(10H NO LXFAC)
00466 213* 887 FORMAT(10H NO BALENG)
00467 214* 888 FORMAT(21H CINDSM REQUIRES LPCS)
00468 215* 889 FORMAT(10H NO NLOOP )
00469 216*
00470 217*
00471 218*
00472 219*
00473 220*
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