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AEC RESEARCH AND DEVELOPMENT REPORT

USER'S MANUAL FOR SRL - HERESY I

HEterogeneous REactor SYstems Code for the IBM 704

D. R. Finch

SRL
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Savannah River Laboratory

Aiken, South Carolina

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Mathematics and Computers
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USER'S MANUAL FOR SRL - HERESY I
HEterogeneous REactor SYstems Code for the IBM 704

by

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March 1966

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ABSTRACT

A detailed description of the SRL - HERESY I code (a production version of the HERESY I code developed by C. N. Klahr) is presented in a manner primarily intended for the user of the code. The heterogeneous reactor theory on which the code is based is developed briefly and followed by detailed coding information. Methods are given for solving finite or semi-infinite lattices containing up to 6500 lattice assemblies of up to 50 different types with unique geometrical or flux symmetry properties. Special techniques used with reflected reactor lattices are also described. Several examples of computations are presented as aids in testing and using the code. Detailed descriptions of programming and operating procedures are presented as appendices.

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USER'S MANUAL FOR SRL - HERESY I

HEterogeneous REactor SYstems Code for the IBM 704

INTRODUCTION

In most heterogeneous reactors, essentially all of the neutron births occur in the lattice fuel elements. Most of the neutron deaths take place in the fuel elements, the target elements, or outside of the reactor. Heterogeneous reactor theory treats these facts explicitly by considering the individual lattice elements as neutron sinks and sources of appropriate strength and then summing the interactions between all the individual elements to obtain the over-all reactor equations. This approach is particularly attractive to the reactor operator because the output of the theory is directly in terms of events at the individual lattice elements, as are the monitoring information and most of the power limitations in the actual reactor. Further, the heterogeneous theory is uniquely well equipped to deal with reactors composed of mixtures of lattice elements or with the insertion of a foreign element into a uniform lattice, since each of the lattice assemblies may be treated individually without recourse to homogenization procedures.

Several computer codes have been developed to take advantage of these attractive features of heterogeneous theory in practical reactor calculations. The HERESY I code^(1,2), written by C. N. Klahr, was one of the first of the heterogeneous codes. It is a rather direct formulation of the equations developed by Feinberg⁽³⁾ and Galanin⁽⁴⁾. Although the original code gave good results, it required a very extensive input, necessitated a very careful verification of all input data, and presented only limited output information. A recoding of the original code into the SRL - HERESY I code was undertaken to eliminate these inconveniences and to permit the use of stored data to reduce the code running time in repetitive problems.

SUMMARY

The SRL - HERESY I code performs reactor lattice calculations based on the heterogeneous reactor theory formulated by Feinberg⁽³⁾. The code will perform calculations on finite lattices, with or without reflector, or semi-infinite lattices. The lattice is limited to a maximum of 6500 total lattice assemblies and to 50 types of lattice locations possessing unique geometrical or flux symmetry properties. The allowed number of physically differing sorts of lattice assemblies is 48.

The code solves for the thermal neutron absorptions in each lattice assembly normalized to the most highly absorbing assembly in the lattice. Resonance neutron absorptions are computed on the basis of a single resonance energy. The edit of the solved problem produces lattice-averaged values of the resonance escape probability, p ; the thermal utilization, f ; the average, η ; and the leakage-corrected static coefficient of reactivity, k .

Three criticality searches built into the code permit variation of a selected parameter in the calculations until critical requirements are reached. The variations may be made in the thermal absorption or thermal η parameter of a specified group of physically identical lattice assemblies or in axial buckling of the lattice with a cosine axial flux shape.

DISCUSSION

I. SUMMARY OF HERESY I EQUATIONS

This section will present a summary of the equations used in SRL - HERESY I. It is not meant to be a complete treatment of the theory underlying the code, nor of heterogeneous reactor systems in general. Insofar as is possible, reference is made to original papers in which the theory is developed.

A. Equations Solved^(2,3,4)

The general approach of heterogeneous reactor theory is to consider the individual reactor lattice elements (referred to as rods in most of the text) as line sources and sinks of neutrons and to sum the interactions between these elements to obtain the over-all reactor equation.

The fundamental equation solved by the SRL - HERESY I Code is

$$\gamma_n i_n = \sum_{m=1}^N \left(\frac{\eta_m}{k} F_{nm}^* - f_{nm} \right) i_m \quad m = 1, 2 \dots N \quad (\text{I.A.1})$$

where

i_n = number of thermal neutron absorptions per unit length per second in n^{th} rod in the lattice

γ_n = ratio of asymptotic neutron flux at the surface of the rod (flux that would be calculated at the rod surface using infinite-medium line-source kernels) to the thermal neutron absorptions per unit length per second in the rod

η_m = number of fission neutrons produced per thermal neutron absorption in the m^{th} rod

F_{nm}^* = fission-to-thermal source kernel (with resonance absorption included). This factor gives the thermal neutron flux at the n^{th} rod due to fissions in m^{th} rod. It is evaluated for an infinitely long line source of fast neutrons in an infinite moderator at a source intensity of 1 fission neutron per second per unit length of the m^{th} rod. This factor includes absorptions in the moderator.

f_{nm} = thermal source kernel. This factor gives the thermal neutron flux at the n^{th} rod due to thermal neutrons emitted at the m^{th} rod. It is evaluated for an infinitely long line source of thermal neutrons in an infinite moderator at a source intensity of 1 thermal neutron per second per unit length of the m^{th} rod.

k = static reactivity coefficient. It corresponds to the effective multiplication constant and is used as a criticality condition.

N = number of rods in the lattice. (Lattice symmetry will enable this number to be reduced later, Section II-C.)

The F_{nm}^* fission-to-thermal source kernel accounts for resonance absorption by subtracting a neutron sink at the resonance absorption energy from the fission-to-thermal source kernel without resonance absorption. If F_{nm} is the fission-to-thermal source kernel with no resonance absorption, then

$$F_{nm}^* = F_{nm} - \sum_{t=1}^N F_{nt}^r A_t \varepsilon_{tm} \quad (\text{I.A.2})$$

where

A_t = ratio of resonance absorptions per unit rod length at an equivalent single resonance energy to the asymptotic slowing down density averaged over rod t at this single equivalent energy

ε_{tm} = fission to resonance slowing down kernel giving number of neutrons per cubic centimeter per second slowing down to the equivalent resonance energy at rod t due to an infinite line source at rod m having a source strength of 1 fission neutron per second per unit length

F_{nt}^r = thermal neutron flux at rod n due to an infinite line source at rod t with source strength of one neutron per second per unit length emitted at the single equivalent resonance energy

Section II-D gives the details of the age diffusion theory kernels used.

B. Self-Consistent Calculation for γ Values

It is not immediately clear how to compute a value of γ for any rod. C. N. Klahr has made detailed studies of this problem for a variety of lattices⁽²⁾, and concluded that the best values are provided by the so-called Self-Consistent Method.

The value of γ , in the Self-Consistent Method, is calculated by considering an infinite regular lattice of rods of the same physical kind. Equation I.A.1 then becomes

$$\gamma_n = \frac{\eta_n}{k} \sum_{m=1}^{\infty} F_{nm}^* - \sum_{m=1}^{\infty} f_{nm} \quad (\text{I.B.1})$$

where η_n is the eta value for the n^{th} rod kind, and use has been made of the facts that $i_m = i_n$ and $\eta_m = \eta_n$. From a relation shown by Klahr⁽²⁾ that

$$\sum_{m=1}^{\infty} F_{nm}^* = p \sum_{m=1}^{\infty} F_{nm} \quad (\text{I.B.2})$$

where p is the resonance escape probability, and the fact discussed in Section I-C that η_n includes ϵ and fuel rod captures, equation I.B.1 becomes, after setting $k = \eta pf$ and using I.B.2.

$$\gamma_n = \frac{1}{f_n} \sum_{m=1}^{\infty} F_{nm} - \sum_{m=1}^{\infty} f_{nm} \quad (\text{I.B.3})$$

where f_n is obtained by a cell calculation or by experiment as

$$f_n = \frac{\text{abs. in all cell components except mod. and coolant}}{\text{abs. in all cell components including mod. and coolant}}$$

Feinberg shows⁽³⁾ that the first sum is approximately given by

$$\sum_{m=1}^{\infty} F_{nm} \approx \frac{1}{V_n \Sigma_a} \quad (\text{I.B.4})$$

where V_n is the cell volume, and Σ_a is the moderator macroscopic absorption cross section. Except for the diagonal elements, the second term of I.B.3 can be summed by use of the Poisson Sum Formula for an infinite regular lattice.

In general the diagonal terms of the f_{nm} sum are very sensitive to the geometrical radius used for evaluation, since in the diffusion theory approximation they are infinite if evaluated at $r = 0$. To circumvent this difficulty, the physical radius of the rod may be used to yield the following expression (valid in the diffusion theory approximation):

$$f_{nn} = \frac{1}{2\pi D_n} K_0\left(\frac{r_n}{L'}\right) \quad (\text{I.B.5})$$

where L' is the moderator diffusion length corrected for axial leakage, and $K_0\left(\frac{r_n}{L'}\right)$ is the modified Bessel function of the second kind.

Using then the results stated in equations I.B.3, I.B.4, and I.B.5 reduces all of equation I.B.3 (N identical equations) to

$$\gamma_n = \frac{1}{V_n \Sigma_a f_n} - \sum_{m=1}^{\infty'} f_{nm} - \frac{1}{2\pi D_n} K_0\left(\frac{r_n}{L'}\right)$$

where the prime on the sum indicates the omission of the n^{th} term. After detailed work on summing over the f_{nm} terms, this takes the following form:

$$\gamma_n = \frac{1}{V_n \Sigma_a} \frac{(1-f)}{f} - \frac{1}{2\pi D_n} K_0\left(\frac{r_n}{L'}\right) + \frac{1}{4\pi D_n} \left[\ln\left(\frac{L'^2}{V_n}\right) + 2.8741 \right] \quad (\text{I.B.6})$$

The constant 2.8741 is geometrical in origin and corresponds to the proper value for a hexagonal lattice; however, this constant does not change appreciably for most lattices, so will give reasonably accurate values for all lattice types.

Axial leakage is taken into account in the computation of γ_n by assuming a cosine axial flux shape. The axial buckling then is defined by

$$B_z^2 = \left(\frac{\pi}{H + \lambda} \right)^2 \quad (\text{I.B.7})$$

where H is the height of the reactor, and λ is the sum of extrapolation distances at the top and bottom of the reactor. A correction may be applied to the moderator absorption cross section and thermal diffusion area to account for the leakage in the form

$$L'^2 = \frac{L^2}{1 + L^2 B_z^2} \quad (\text{I.B.8})$$

$$\Sigma_a' = \Sigma_a (1 + L^2 B_z^2)$$

where Σ_a and L^2 are the uncorrected or axially infinite values of Σ_a' and L'^2 .

Fast axial leakage is accounted for by decreasing the reactivity k by a factor $\exp(-\tau B_z^2)$:

$$k_{\text{corr}} = k_{\text{heresy}} e^{-\tau B_z^2} \quad (\text{I.B.9})$$

C. Fast Fission Correction

A first order correction may be made to η to take account of fast fission. Basically this amounts to redefining η used in HERESY to be

$$\eta_{\text{heresy}} = \frac{1}{f_{\text{heresy}}} (\eta f)_{\text{cell}} \epsilon$$

where ϵ and $(\eta f)_{\text{cell}}$ are obtained separately from cell calculations, and f_{heresy} is the heterogeneous f defined in Section B.

D. Resonance Absorption Parameter

Several different methods have been devised to give reasonable values of A . Klahr⁽²⁾ suggests one of the following:

1. An approximate method deriving A from resonance escape probability as

$$p = 1 - \frac{A}{V_n} \quad (\text{I.D.1})$$

where V_n is the lattice area per rod. This is applicable (in D_2O lattices) only for lattice pitches ≤ 8 inches, for which the resonance flux is approximately flat.

2. A more exact form of equation I.D.1 derived from the self-consistent form is

$$p = 1 - A_n \sum_{m=1}^{\infty} \varepsilon_{nm} \quad (\text{I.D.2})$$

3. If an effective resonance integral is available for the rod in question, A is given by:

$$A = \frac{N_f V_f}{(\xi \Sigma_s)_{\text{cell}}} (R I_{\text{eff}}) \quad (\text{I.D.3})$$

where all terms have their conventional meaning.

E. Method of Solving Equations

Equation I.A.1 is first reduced to a standard eigenvalue problem as follows:

1. The sum on the right side of I.A.1 is broken into two parts, the $\sum_{m=1}^N f_{nm} i_m$ being brought to the left side of the equation.

2. The equation is now rewritten in the form:

$$\sum_{m=1}^N (f_{nm} + \gamma_m \delta_{nm}) i_m = \frac{1}{k} \sum_{m=1}^N \eta_m F_{nm}^* i_m \quad (\text{I.E.1})$$

or, in matrix form:

$$A I = \frac{1}{k} B I \quad (\text{I.E.2})$$

3. Inverting matrix A and multiplying I.E.2 through by A^{-1} from the left yield

$$I = \frac{1}{k} A^{-1} B I \quad \text{or} \quad C I = k I \quad (\text{I.E.3})$$

where

$$C = A^{-1} B$$

4. A standard power iteration is then used to solve for k as the largest eigenvalue of equation I.E.3. The eigenvectors i_n normalized to the largest value are also determined from this same power iteration.

5. Criticality searches enter at this point, if desired. The three types of searches are:

a. Searches on η merely change the B matrix in equation I.E.2, thus reforming matrix C of equation I.E.3, and re-solving the eigenvalue problem is all that is required to re-solve the equation. Convergence is tested externally.

b. Searches on gamma demand that matrix A of equation I.E.2 be re-formed and inverted to form equation I.E.3. Hence matrix inversion and solving the eigenvalue problem are required for each iteration of a criticality search on gamma.

c. Searches over axial buckling alter the kernel matrices of equation I.E.1, hence require regeneration of these matrices from new kernels formed with the current buckling value. Each iteration must therefore, re-solve the entire problem.

II. DESCRIPTION OF CODE

A. General Description - Machine Requirements

SRL - HERESY I was written for an IBM 704 computer having a 32K core storage, an on-line printer, and a minimum of five tape units.

The program itself is in two parts or links, these being connected by means of COMMON STORAGE using the CHAIN subroutine (see Appendix D). Link 1 of the code has 13 primary subprograms, a package of library function subroutines, and a package of FORTRAN II control subroutines. The purpose of this link is to (a) read data; (b) check input data for gross errors; (c) generate appropriate kernels and geometries; (d) compute input parameters by the Self-Consistent Method of Section I-B; and (e) output a summary of all input data and generated quantities.

Link 2 of the code takes the data generated in Link 1 and proceeds to solve the problem by the method discussed in Section I-E. This link contains 15 primary subprograms, a package of library subroutines, and a package of FORTRAN II control subroutines.

Links 1 and 2 have 15,725 words of COMMON STORAGE shared between them. This enables information to be permanently stored in the machine such that both Links 1 and 2 can use it. In general, Link 1 reads and stores, computes, or modifies the information contained in the shared storage portion of COMMON STORAGE, and Link 2 draws from this modified information the necessary data to solve any particular problem. After a particular problem is completed, the numbers that are stored in the shared storage are those that correspond to the solved problem. For example, the initial guess to the eigenvectors for the next problem to be run will be the solution to the problem just completed unless this quantity is re-specified in the input to the next problem. Again, if a criticality search has been called for and completed, the η (or γ or B_z^2) value in the shared storage will be that value corresponding to the converged criticality search just completed. One might utilize this latter feature in a sequence of problems in which the first problem normalizes a particular parameter to correspond to $k_{eff} = 1$, and later problems then use this value of the parameter to determine changes that arise from changing other lattice or rod parameters.

Several operations have been streamlined in the SRL code to enable faster running time than for the original HERESY code. An example occurs in the first stages of computation in Link 2. The matrices corresponding to the symbols $\{F_{nm}\}$, $\{F_{nm}^F\}$, $\{g_{nm}\}$, and $\{f_{nm}\}$

are formed in a subprogram, MATELM, and depend only upon rod coordinates and the kernel used. These matrices are placed in the shared portion of storage and re-used in sequences of problems using the same kernel and geometry. This is of particular significance in problems containing a large number of rod types, since the time required to form these matrices for any problem increases approximately as the square of the dimensionality of the matrices.

B. Geometry Tapes

The re-use of matrix elements in large sets of problems leads to a convenience feature with rod coordinates. A total of 13,000 locations in shared storage are allocated to store the X and Y coordinates of all rods (i.e., 6,500 rods maximum). To enable the matrices mentioned in Section II-A to be placed in the shared storage (a total of 10,000 locations are required), the X and Y coordinates of each rod are written out on binary tape for either temporary or permanent storage as the user specifies. The coordinates are always available for use, when properly specified, if a new kernel requires the matrices to be re-formed. If new coordinates are generated in Link 1, a new binary tape, either temporary or permanent, will be written out as specified by the user.

Temporary geometry tapes are written on, or read from, tape unit 2, while permanent geometry tapes are written on or read from, tape unit 3.

A temporary tape will contain one binary record of 13,000 words of storage, or the entire contents of the X and Y coordinate arrays. If a new geometry is generated by Link 1 the subroutine XYLIB will overwrite the previous record with the new coordinates. No END of FILE mark is placed at the end of this tape at any time. The input variable NFILE must be zero to use this method of geometry storage.

A permanent tape will contain several sets of coordinates, each set occupying two binary records on the tape. The first of these records is called the File Number and is specified by the number NFILE in the input data. The second record is the 13,000 binary words contained in the X and Y coordinate arrays plus a checksum word as an add-and-carry logical check on tape operation. Three indicators are necessary to specify the use of the permanent tape: (1) the number NFILE, which is the file number of the desired geometry to be used in the problem; (2) the indicator NWRITE, which tells if a new geometry is to be written out or not; and (3) KØØR if the initial geometry is to be read from a permanent tape. NFILE may be any number positive or negative (but not zero) having up to five fixed point digits, but

not exceeding 2^{15} . NWRITE is either zero (do not write new coordinates on tape) or 1 (write new coordinates on tape). Any time that tape 3 is read all coordinates taken from the tape will be scaled by the current value of an arbitrary scale factor XCOR.

The following operations may be performed with the permanent tapes:

1. If NFILE is specified, and no new coordinates occupy the X and Y coordinate arrays (NWRITE = 0), the coordinates in the second record corresponding to NFILE in the first record may be read, but writing out the coordinates is redundant, hence is skipped.

2. If NFILE is specified and new coordinates occupy the X and Y coordinate arrays (NWRITE = 1), two cases arise:

- a. If NFILE is a file number already on the tape, the new coordinates will be written in place of the old coordinates on the tape.

- b. If NFILE is a file number not on the tape, the two record file number and coordinate geometry with checksum will be added at the end of the tape. An END of FILE mark will be placed following the last record on the tape.

Use of the permanent tapes allows often used geometries to be utilized repeatedly for most rapid running of the code. (See Appendix C for operating instructions on generating permanent tapes.)

C. Rod Specifications

The rods that make up any particular lattice may be made of several different materials, and may be placed in the reactor with different degrees of flux symmetry. The following scheme is used in SRL - HERESY I for specifying material properties and symmetry properties.

The term "rod" is used to designate a lattice assembly of any variety, i.e., a rod may be a fuel element, a control rod, a control rod cluster, an instrument thimble, or any other assembly as specified by the user.

The term "rod kind" is used to designate a rod with specific material properties independent of any geometrical or flux symmetry properties in the problem.

The term "rod type" is used to designate rods of the same "rod kind" that have the same geometrical and flux symmetry properties in the problem. Usually several rods in the problem will have equivalent geometrical position and flux symmetry properties and, therefore, be of the same "rod type" by definition. The absorptions in these rods are identical; it is only necessary to solve for the absorption in one rod of each given "type". This property enables the dimensionality of the matrices considered in solving the problem (see Section I-E) to be reduced to the number of different "rod types" that occur in the problem instead of the total number of rods.

In forming any problem, the rods are classified primarily according to their respective "types". Coordinates are produced and classified according to "type", matrices are formed having indices based on the number of "types", and running times are dependent mainly on the number of "rod types". The maximum number of "types" is 50.

Each "rod type" (numbered from 1 to NTYPE) has associated with it several other parameters in the code. The mnemonics NRPRT(I) give the number of rods in the lattice of each "rod type", the mnemonics KIND(I) give the "rod kind" that is associated with each "rod type", and the mnemonics NICPRT(I) supply information to the geometry generators to indicate the number of initial equivalent coordinate sets that are input for a given "rod type".

All rod material parameters are entered by "kind" rather than by "type", and through the mnemonics KIND(I) are internally associated with the correct rod type. The properties for each rod "kind" are specified by three parameters:

η_m = number of fast neutrons per thermal absorption
= fission parameter

γ_m = ϕ asymp./ i_m
= thermal absorption parameter

A_m = resonance absorptions per slowing down density
= resonance absorption parameter

Calculations of γ_m by the Self-Consistent Method are performed for a rod "kind" by setting γ for the "kind" to be calculated equal to zero and supplying a value of f (mnemonic THERU(I)) and V_{cell} (mnemonic VCELL(I)) for this rod "kind".

The maximum number of rod kinds is 48 for any problem.

D. Kernel Generators

The SRL - HERESY I code utilizes two kernel generators: one in Link 1 and a shorter version in Link 2. The larger generator in Link 1 examines the input information and calculated information for errors prior to attempting to use the information in solving the problem. The small generator in Link 2 does not contain any error checking routines (data are assumed to be sufficiently checked in Link 1), and is used only for the Buckling Criticality Search.

The kernels formed are age diffusion kernels having slowing down distributions $q(r)$ of the form

$$q(r) = \sum_{i=1}^3 \frac{B_i}{4\pi\tau_i} e^{-\frac{r^2}{4\tau_i}}$$

where r is the radial distance from the source line; τ_i corresponds to either fission-to-thermal age, or to fission-to-effective resonance energy age; and B_i is a weighting fraction corresponding to the specified τ_i value. This is a triple gaussian sum with a single effective neutron age, τ .

Using the above definition of the slowing down source in the diffusion theory approximation leads to the following kernels⁽¹⁾ (see Chapter I for definitions of symbols).

$$g(r) = \sum_{i=1}^3 \frac{B_i}{4\pi\tau_i^R} e^{-\frac{r^2}{4\tau_i^R}}$$

$$F(r) = \sum_{i=1}^3 \frac{B_i}{2\pi D} e^{-\frac{\tau_i}{L^2}} \left[K_0\left(\frac{r}{L}\right) + \frac{\tau_i}{2L^2} e^{-\frac{r^2}{4\tau_i}} - \frac{1}{2} \left(1 + \frac{r^2}{4L^2} \right) E_1\left(\frac{r^2}{4\tau_i}\right) \right]$$

$$F^R(r) = \sum_{i=1}^3 \frac{B_i}{2\pi D} e^{-\frac{(\tau_i - \tau_i^R)}{L^2}} \left[K_0\left(\frac{r}{L}\right) + \frac{(\tau_i - \tau_i^R)}{2L^2} e^{-\frac{r^2}{4(\tau_i - \tau_i^R)}} - \frac{1}{2} \left(1 + \frac{r^2}{4L^2} \right) E_1\left(\frac{r^2}{4(\tau_i - \tau_i^R)}\right) \right] \quad (\text{II.D.1})$$

In the thermal region, a standard diffusion theory kernel is used of the form

$$f(r) = \frac{1}{2\pi D} K_0\left(\frac{r}{L^1}\right) \quad (\text{II.D.2})$$

Self-contributions (i.e., diagonal elements of f_{nm}) of any rod 1 to itself in the thermal region have a special form since

$$r \xrightarrow{L^1} 0 K_0\left(\frac{r}{L^1}\right) \rightarrow \infty$$

This problem is eliminated in the standard heterogeneous fashion of evaluating these kernel elements at the rod radius. This leads to a group of numbers (one for each rod type) designated the rod SMF(I) and defined by

$$\text{SMF}(I) = f(R_1) = \frac{1}{2\pi D} K_0\left(\frac{R_1}{L^1}\right) \quad (\text{II.D.3})$$

where R_1 is the radius of the 1th rod.

Reference to equation I.B.6 shows that this quantity SMF(I) appears as the second term of the right side of that equation. However, the quantities γ_n and SMF(I) are never used separately, only as the sum of the two. Because of this fact it is more convenient usually to define a new quantity γ_n^1 as

$$\gamma_n^1 = \gamma_n + \text{SMF}(I)$$

$$\gamma_n^1 = \frac{1}{V_n \Sigma_a} \frac{(1-f)}{f} + \frac{1}{4\pi D_n} \left(\ln \frac{L^1{}^2}{V_n} + 2.8741 \right) \quad (\text{II.D.4})$$

This equation now has no dependence on the physical radius of the lattice assembly and is far more convenient to use.

The use of the quantities γ_n^1 in the code is facilitated in the following manner.

1. If rod radii are provided in the input data, γ_n and SMF(I) are calculated and have their respectively defined meanings.

2. If rod radii are not provided (i.e., are zero) in the input data a signal value of 0.10101 is assigned to each $SMF(I)$ and the γ_n corresponding to this value of $SMF(I)$ is calculated. In this case, γ_n and $SMF(I)$ do not individually have physical meaning, but their sum γ_n^1 is correct.

Kernels are stored as tabular arrays of maximum size 250 points. The finite height of a reactor is taken into account in the kernel generators through a correction on L^2 and Σ_a . This correction is of the form

$$L_{corr}^2 = \frac{L^2}{1 + L^2 B_z^2} \text{ and } \Sigma_a^{corr} = (1 + L^2 B_z^2) \Sigma_a$$

and is the same correction as used in calculating certain terms in the expression for γ in equation I.B.6.

Two restrictions are applied to the kernels to assure that proper Bessel's interpolation may be carried out for points not directly on a kernel point. Such a condition occurs in the subprograms MATELM and USKER where the kernel matrices are formed. The restrictions are:

- The mesh spacing must be less than or equal to one-half of the smallest center-to-center rod spacing in the lattice

$$XMESH \leq \text{smallest value } \frac{|r_1 - r_j|}{2}$$

- The kernel shall overlap the edge of the lattice on all sides, by at least two points

$$247 * XMESH \geq \text{largest value } |r_1 - r_j|$$

The first restriction arises due to the possibility that in some problems the kernel points used for interpolation might fall as much or more than a full lattice pitch away from the position desired. This does not affect the equations formulated in Chapter I; however, it is physically unsatisfying to know that the interpolation is using points this widely spaced. The value of 1/2 as the smallest center-to-center rod spacing simply assures that the points used for interpolation are at most exactly one lattice pitch away from the rod in question.

The second restriction arises in order to prevent the use of the 251st and 252nd points of a 250-point kernel from being used. Obviously these points would pick up values from other quantities stored at these locations in the memory (see Chapter V, problem 3). Since one point is assigned at position zero, the outer edge of the kernel is at $249 \cdot \text{XMESH}$ cm from the origin point. Now, because Bessel's interpolation uses points $m-1$, m , $m+1$, and $m+2$ for interpolation, m can have a value of 247 at most. Therefore, the smallest kernel mesh that may be used is equal to

$$\text{XMESH} \geq \frac{\text{largest center-to-center spacing in lattice}}{247}$$

For infinite lattices (Section G), this requirement is modified to the extent that the largest center-to-center spacing for a rod type is considered to be the distance between the initial coordinates of the "rod type" as input to the infinite lattice generator and the rod that is the greatest distance from it.

In addition to the radially infinite problems for which the above kernels are defined, the code can also treat, in an approximate manner, the problem of a radially finite reactor. This is accomplished by forming kernel matrices that extrapolate to zero at some radius from the center of the lattice.

The technique used employs the method of images. Each rod in the reactor is treated as having an image rod which provides a weighted negative contribution to each kernel element balancing the positive contribution from the real rod. The weighting of the image rod is provided by a factor $\sqrt{\rho_1/\rho_0}$ where ρ_1 is the distance of the image rod from the center of the reactor, and ρ_0 is the distance of the real rod from the center of the reactor. The image rod is positioned a distance outside the reactor equal to the radial distance; the real rod is situated inside the specified reflector radius (mnemonic RADRFL) which may be loosely defined as the reactor radius.

In practice it has been found that a ratio of the quantity RADRFL to the actual extrapolation radius (R_{extrap}) for a uniform fuel lattice varies over an approximate range

$$0.9 \leq \frac{\text{RADRFL}}{R_{\text{extrap}}} \leq 1.0$$

depending on the lattice. Because of this fact it is recommended that a test problem using a uniform fuel lattice be run to determine the real kernel extrapolation distance before using a particular finite kernel.

Another radius that may be input is the quantity ROEMIN which causes the code to omit calculating image contributions for rods within this radius, since their images are very far away from the lattice.

E. Criticality Searches

The SRL - HERESY I code has three built-in criticality searches that may be selected as options. These enable any specified eigenvalue (corrected for axial buckling) to be searched for as a function of γ , η , or B_z^2 .

Searches on γ and η require a "rod kind" to be specified for the search. Searches on B_z^2 require that the SMF(I) all be equal to 0.10101 (Section II-D).

Two edit options are available with all searches. The user may obtain edits for each iteration of the search, or edit only the final iteration following convergence. A short summary of each iteration will be printed out in this latter case.

The γ searches recompute the thermal utilization (f_{heresy}) based on the γ value at convergence. This value is inserted back into the proper storage position for the "rod kind" specified in the search.

F. Finite Lattice Geometry Generators

Three geometry generators incorporated into the code accept a bare minimum of data to produce uniform lattices of rods on hexagonal or square pitches, or patterns of rods on circles about the axis of a cylindricized reactor.

1. Hexagonal Lattices

Hexagonal lattices require for input the X and Y coordinates of each rod with angles in the range $0^\circ \leq \theta < 60^\circ$ inclusive. The code then generates the remaining five rotationally symmetric elements around the reactor. Provision is made for more than six rods at a given radius to be of the same rod type. This is accomplished by providing more than one set of initial coordinates for a particular rod type to a limit of three sets of coordinates.

For hexagonal lattices, this latter provision is often used in the outer hexes of the lattice. A second reason for its inclusion is that in studying large lattices it is often convenient to approximate the outer edge of a reactor by making several distinctly different rods equivalent to a single rod type.

2. Square Lattices

Square lattices require for input X and Y coordinates of each rod in quadrant 1 with angles in the range $0^\circ \leq \theta \leq 45^\circ$ inclusive. The code makes use of the fact that if a rod in this sector has polar coordinates R and θ , then the rods of this type will have polar coordinates of radius R and angles θ , $\pi/2 - \theta$, $\pi/2 + \theta$, $\pi - \theta$, $\pi + \theta$, $3\pi/2 - \theta$, $3\pi/2 + \theta$, and $2\pi - \theta$. Rods at 0° and 45° are not included twice. Provision is made for two rods at different positions (same radius) to be of the same rod type by allowing more than one set of initial coordinates to be input for a particular rod type.

3. Circular Patterns of Rods

This generator allows a pattern of rods all centered on circles about the axis of the reactor to be produced. Three initial coordinates are required in this case, these being the X and Y coordinates of the first rod at a particular radius as one proceeds counterclockwise from the positive X axis, and the angular separation (in degrees) between any two rods at the particular radius. Rods are generated from the initial position (at angle θ) to an angle equal to the initial angle plus 360 degrees (or once around the circle). Single rods may be generated by making the angular separation between rods greater than 360 degrees.

It will be noted that a hexagonal lattice is a special case of this lattice; however, hexagonal lattices are used far more than these circular patterns, hence are included as a special generator.

With the use of these three basic types of lattice patterns, most lattices may be generated quite easily. If, in addition, the permanent library tape option is used the generation of new lattices becomes more and more infrequent.

All three of the generators internally compute the number of rods for each "rod type" and store them. It is therefore unnecessary to enter these data when geometry is being generated.

G. Infinite Lattice Geometry Generator

The infinite lattice geometry generator considers only overlapping patterns of rods on rectangular pitches. However, if a block of rods in a hexagonal or other type of lattice can be found that displays infinite repeatability in rectilinear translation, the infinite lattice generator may be used. The user would then specify each rod in the block as a distinct "rod type" and the X and Y pitch of the rectangular lattice of this type as being the X and Y translational distance to the next rod of that type in the nearest symmetry block.

The information necessary to generate lattices of this type is
(a) the X and Y pitches for each rod type (as defined above), (b) the X and Y coordinates of the nearest rod of each type to the origin, and
(c) the maximum radius to which rods are to be generated.

As in the finite lattice, the number of rods for each rod type are determined by the generator and stored.

H. Print, Punch, and Plot Options

Along with the geometry generators go these three auxiliary sub-programs which will respectively list the rod coordinates for each rod type, punch a deck of cards suitable for direct input back into the code (this is restricted to finite lattices only), and plot to a specified scale (XSCALE and YSCALE values) the lattice geometry. These routines may be used to visually check each coordinate geometry for a particular lattice, and as such are recommended during the first use of any new lattice geometry that may be generated. The punch subroutine allows a seldom used lattice to be stored away on reloadable cards for future use, but is hardly the most convenient method of storage. In addition it requires a large amount of time to punch a deck for a big lattice.

III. INPUT INFORMATION

A. Data Preparation

All data are broken down into three groups and input in the following order:

1. Alphanumeric data (1 card only)
2. Fixed point data
3. Floating point data

For each problem run there must be at least one card of each type of data for input. The three types of data are described separately in the following paragraphs.

1. Alphanumeric Data

These data consist of one card of alphanumeric symbols contained in columns 1-72. If a 1 is placed in column 1, the output from the code will begin on a new page; any other standard carriage control symbol may be placed in column 1 (see IBM 704 FORTRAN Reference Manual, C28-6106, page 45).

2. Fixed Point Data

Fixed point data are input in a fixed card format that enables any number of individual datum up to five to be read from a single card. The card format is schematically drawn below:

<u>Card Columns</u>							
<u>1</u>	<u>5</u>	<u>6-12</u>	<u>13-24</u>	<u>25-36</u>	<u>37-48</u>	<u>49-60</u>	<u>61-72</u>
Last card indicator	Number of datum on this card (if blank or zero is assumed to be 5)	Relative location of datum 1 of this card (see following list)	Datum 1	Datum 2	Datum 3	Datum 4	Datum 5

All fixed point input data are stored in a large block of core designed as the LL(225) array. Each datum is assigned a relative location in this array, and a list of these locations follows. The relative location of datum 1 of the card being read is punched in columns 6-12 of the card.

Data 2, 3, 4, and 5 are then stored in sequential relative locations in the LL(225) array. If it is not desirable to input a full card of data (5 datum), it is only necessary to place the number of datum actually appearing on the card to be read in column 5 (a blank or zero will cause 5 datum to be read). For instance, if it is necessary to change one property of one rod type in a list of 50 rod types, the relative location of the datum may be determined from the following list and placed in columns 6-12, a number 1 placed in column 5 of the card, and the new datum placed in the datum 1 field of the card as defined above. All fixed point data are entered to the extreme right of the field defined, blank columns being read as zeros.

The last card of fixed point data must have a 1 punched in column 1 to so indicate it is the last card.

Relative locations for fixed point data are contained in the following list:

Problem Specification Information

- | | | |
|----|-------------------------|------------------------------------|
| 1. | LL(1) = INF | <u>Finite - Infinite Indicator</u> |
| | = 0 | Finite lattice. |
| | = 1 | Infinite lattice. |
| 2. | LL(2) = KERNEL | <u>Kernel Generator Indicator</u> |
| | = 0 | Use old kernels. |
| | = 1 | Generate new kernels. |
| | | (Must be 1 for first problem). |
| 3. | LL(3) = KOOR | <u>Coordinate Indicator</u> |
| | = 0 | Use previous coordinates. |
| | = 1 | Generate new coordinates |
| | | (Finite or infinite). |
| | = 2 | New coordinates are contained |
| | | in the floating point input |
| | | data. |
| | = 3 | Read geometry from permanent |
| | | tape on unit 3. |

Iteration Criteria

- | | |
|---------------|---|
| 4. LL(4) = N1 | Maximum number of iterations for matrix inversion correction (set equal to 20 if left blank). |
| 5. LL(5) = N2 | Maximum number of iterations on eigenvalue iteration (set equal to 200 if left blank). |
| 6. LL(6) = N3 | Maximum number of iterations on criticality searches (set equal to 25 if left blank). |

Kernel Information

(Must be included if KERNEL = 1)

- | | |
|-------------------|---|
| 7. LL(7) = NOGRID | Number of grid points in kernel (≤ 250). |
|-------------------|---|

Geometry Generator Information

(Must be included if KGOR = 1)

- | | |
|--------------------------|---|
| 8. LL(9) = LAT | <u>Lattice Type Indicator</u>
(Finite lattice only) |
| = 0 | Uniform hexagonal lattice. |
| = 1 | Uniform square lattice. |
| = 2 | Pattern of rods centered on circles about axis of reactor. |
| 9. LL(10)... = NICPRT(J) | Number of sets of initial equivalent coordinates input for each rod type. NTYPE NUMBERS. NICPRT(J) ≤ 3 . Include for finite lattices only. |
| 10. LL(60) = KPRNT | <u>Print Option Indicator</u> |
| = 0 | Do not list geometry. |
| = 1 | List geometry. |
| 11. LL(61) = KPLOT | <u>Plot Option Indicator</u> |
| = 0 | Do not plot generated geometry. |
| = 1 | Plot generated geometry. |

- | | | |
|-----|----------------|---|
| 12. | LL(62) = KSPIP | <u>Punch Option Indicator</u>
(Finite lattices only) |
| | = 0 | Do not punch geometry deck. |
| | = 1 | Punch geometry deck. |

Rod Information

- | | | |
|-----|---|---|
| 13. | LL(63) = NTYPE | Number of rod types (≤ 50) |
| 14. | LL(64) = NKIND | Number of rod kinds (≤ 48) |
| 15. | LL(65)... = NRPRT(J) | Number of individual rods for each rod type. NTYPE NUMBERS (may be omitted if KØØR = 1; the geometry generators then supply these numbers). |
| 16. | LL(115)... = KIND(J) | Rod kind associated with each rod type. NTYPE NUMBERS. |
| 17. | LL(165) = LCRIT
= 0
= 1
= 2
= 3 | <u>Criticality Search Indicator</u>
No criticality search.
Search on ETA.
Search on GAMMA.
Search on axial buckling. |
| 18. | LL(166) = KSERCH | Rod <u>KIND</u> to be searched over. |
| 19. | LL(167) = NSEPR
= 0
= 1 | <u>Edit Option on Criticality Search</u>
Edit only converged solution.
Edit each iteration of criticality search. |

Miscellaneous Information

- | | | |
|-----|-------------------------------|---|
| 20. | LL(168) = KDAT
= 0
= 1 | <u>Input Data Print Option</u>
Data will be printed.
Omit data print. |
| 21. | LL(169) = KTEST
= 0
= 1 | <u>Last Problem Indicator</u>
More problems follow.
Last problem. |

- | | | |
|-----|------------------|--|
| 22. | LL(224) = NWRITE | <u>New File for Permanent Tape Indicator</u> |
| | = 0 | Do not write new geometry file. |
| | = 1 | Write new geometry file. |
| 23. | LL(225) = NFILE | File number of geometry on designated permanent tape (on unit 3)(if zero uses scratch tape on unit 2). |

3. Floating Point Data

Floating point data are also input in a fixed card format that enables any number of individual datum up to five to be read from a single card. The card format is identical to that discussed for fixed point data, the only change made for floating point data is for the data to be read in FORTRAN FIXED DECIMAL FLOATING POINT notation (standard FORTRAN F type conversion). The number 10^{-6} would thus appear as 0.000001 on the input card. A floating point datum may be placed anywhere within the defined fields on the card so long as the decimal point is assigned the correct position in the number. If no decimal point is written, it is automatically assigned a position following the character appearing in the last column of the datum field. Blank columns are read as zeros.

All floating point input data are stored in a large block of core designated as the AA(15,500) array. Each datum is assigned a relative location in this array, and a list of these relative locations appears below. The relative location of datum 1 on the card must be punched into columns 6-12 of the card, and provision for reading less than five datum is identical to that for fixed point data.

The last card of floating point information must have a 1 punched in column 1 to so indicate it is the last.

Relative locations for fixed point data are contained in the following list:

Convergence Criteria

1. AA(1) = CR1 Convergence criteria on matrix inversion (if left blank is set equal to 0.000001).
2. AA(2) = CR2 Convergence criteria on eigenvalue iteration (if left blank is set equal to 0.00001).
3. AA(3) = CR3 Convergence criterion on criticality search (if left blank is set equal to 0.00001).

Scale Factors

4. AA(4) = XCØR Arbitrary scale factor (cm) applied to all coordinates (finite or infinite lattices).
5. AA(5) = XSCALE Plotting increment along X axis in geometry plotter (if left blank is set equal to XCØR).
6. AA(6) = YSCALE Plotting increment along Y axis in geometry plotter (if left blank is set equal to XCØR).

Reflector Information

7. AA(7) = RADRFL Reflector outer radius (cm) (if zero problem is treated as having an infinite reflector).
8. AA(8) = RØEMIN Radius (cm) within which the reflector may be treated as infinite.

Criticality Search Information

9. AA(9) = EIGEN Eigenvalue to be searched for (if left blank will be set equal to 1.0).

Kernel Information

(Must be included if KERNEL = 1)

10. AA(10) = XMESH Mesh spacing between kernel points (cm) must be such that

$$\frac{\text{largest } |r_1 - r_j|}{247} \ll \text{XMESH} \ll \frac{\text{smallest } |r_1 - r_j|}{2}$$

- 11. AA(11) = BSQ Axial buckling ($1/\text{cm}^2$)
- 12. AA(12) = XLSQ L^2 of moderator uncorrected for leakage (cm^2)
- 13. AA(13) = SIGA Σ_a moderator ($1/\text{cm}$)
- 14. AA(14)... = B(J) Weighting factors on TAU's and TAUR's for each term of triple Gaussian. (if left blank uses values 0.56, 0.44, 0.0)
- 15. AA(17)... = TAU(J) Fission to thermal age for each term of triple Gaussian (cm^2)(if left blank uses 154.3, 68.3, 0.0).
- 16. AA(20)... = TAUR(J) Fission to resonance energy age for each term of triple Gaussian (cm^2)(if left blank uses 126.5, 40.5, 0.0).

Rod Information

- 17. AA(25)... = XLSQ2(J) L^2 for each rod kind. NKIND VALUES (cm^2), if left blank are set equal to XLSQ.
- 18. AA(75)... = SIGC(J) Σ_a used to compute γ for each kind. NKIND VALUES (if left blank are set equal to Σ_a moderator).
- 19. AA(125)... = VCELL(J) Cell area used to compute γ for each rod kind. NKIND VALUES (cm^2).
- 20. AA(175)... = GAMMA(J) Thermal absorption parameter. NKIND VALUES. (if left blank must have THERU(J) value).
- 21. AA(225)... = THERU(J) Thermal utilization. NKIND VALUES (If GAMMA(J) is zero, computes GAMMA(J) using this value).
- 22. AA(275)... = ETA(J) Number of fast neutrons produced/thermal absorption. NKIND VALUES.
- 23. AA(325)... = AR(J) Resonance absorption parameter. NKIND VALUES.
- 24. AA(375)... = RADIUS(J) Rod radius (cm). NKIND VALUES (If zero assigns SMF(J) = 0.10101).

Finite Geometry Generator Information

(included if $K\emptyset\emptyset R = 1$ and $INF = 0$)

25. AA(425) ... = $X\emptyset(I,1)$ Initial X coordinates
AA(475) ... = $X\emptyset(I,2)$ $\{1 \leq I \leq NTYPE\}$
AA(525) ... = $X\emptyset(I,3)$ (see Notes 2 and 3 and Section-II-E)
26. AA(575) ... = $Y\emptyset(I,1)$ Initial Y coordinates
AA(625) ... = $Y\emptyset(I,2)$ $\{1 \leq I \leq NTYPE\}$
AA(675) ... = $Y\emptyset(I,3)$ (see Notes 2 and 3 and Section-II-F)
27. AA(725) ... = $THETA(I,1)$ Angular separation in degrees between rods
AA(775) ... = $THETA(I,2)$ on same radius about axis of reactor.
AA(825) ... = $THETA(I,3)$ $\{1 \leq I \leq NTYPE\}$ (see Note 2) (Not required for square or hexagonal lattices).

Infinite Geometry Generator Information

(Included if $K\emptyset\emptyset R = 1$ and $INF = 1$)

28. AA(875) ... = $WDX(J)$ X pitch for lattice J. NTYPE VALUES (see Note 3).
29. AA(925) ... = $WDY(J)$ Y pitch for lattice type J. NTYPE VALUES (see Note 3.)
30. AA(975) ... = $WX1(J)$ X coordinate of rod nearest origin for lattice J. NTYPE VALUES (see Note 3.)
31. AA(1025) ... = $WY1(J)$ Y coordinate of rod nearest origin for lattice J. NTYPE VALUES (see Note 3).
32. AA(1075) ... = $RADII$ Radius beyond which no more rods are considered (cm).

Miscellaneous Information

33. AA(1080) ... = $V\emptyset(J)$ Initial guess to eigenvectors. NTYPE VALUES (if left blank are set equal to 1.0)
34. AA(2500) ... = $X(J)$ X coordinates for all rods. (see Notes 1 and 3).

NTYPE

$$\sum_{I=1} NRPRT(I) \text{ VALUES}$$

35. AA(9000) ... = Y(J) Y coordinates for all rods. (see Notes 1 and 3).

$$\sum_{I=1}^{NTYPE} NRPRT(I) \text{ VALUES}$$

Note 1

Coordinates are input in sequence starting with rod type 1 and running over all rods of that type. These are followed by rod type 2, 3, 4..., where, for each type, the coordinates of all rods of that type are successively entered.

Note 2

Entering the initial coordinates $X\emptyset(I,J)$, $Y\emptyset(I,J)$, and $THETA(I,J)$ for the finite lattice geometry generator must be done in a special order. The matrices for these quantities are 50 by 3 or 50 rows and 3 columns. Data are entered columnwise which implies that for each value of J the index I runs through its full range. Hence using X0 as an example the values would be entered as follows:

$X\emptyset(1,1) \dots, X\emptyset(I,1) \dots, X\emptyset(50,1), X\emptyset(1,2) \dots, X\emptyset(I,2) \dots$
 $X\emptyset(50,2), X\emptyset(1,3) \dots, X\emptyset(I,3) \dots, X\emptyset(50,3)$

Some relative locations associated with these are:

F(425) = $X\emptyset(1,1)$	F(575) = $Y\emptyset(1,1)$	F(725) = $THETA(1,1)$
F(475) = $X\emptyset(1,2)$	F(625) = $Y\emptyset(1,2)$	F(775) = $THETA(1,2)$
F(525) = $X\emptyset(1,3)$	F(675) = $Y\emptyset(1,3)$	F(825) = $THETA(1,3)$

Note 3

All coordinates and pitches contained in items 25, 26, 28, 29, 30, 31, 34, and 35 will be multiplied by the number XCØR contained in 4.

B. Problem Sequencing

The data that are input into the code according to the format described in Section III-A are stored permanently in the memory of the computer and may be re-used at any time. In survey type calculations or where a sequence of different values of one or more parameters are the only changes from one problem to the next, it is only necessary to enter the changes to the immediately previous problem. However, the restriction that each problem has at least one card of each of the three types of data (see Section III-A) is maintained and must be satisfied.

This type of data input, while convenient, has some pitfalls that require care to avoid. The more important of these pitfalls are the following:

1. If geometry is generated in the 1st problem and is re-used in the 2nd problem, the data for the second problem must "turn off" the geometry generator or else the geometry will be regenerated. A similar comment applies to the kernel generator, geometry print, geometry plot, and geometry punch options.
2. If the self-consistent method is to be used to calculate a γ in any problem after the first, the γ value must be set equal to zero in the input data as a signal to perform this calculation.
3. In buckling searches an additional requirement is made that the SMF(I) values be those specified by the code (see Section II-E).
4. At the end of any criticality search the values stored in the memory will be those corresponding to the converged search values. If the original input data are desired, it must be restored in the input for the next problem.
5. When a new kernel is to be generated with different moderator properties from a previous problem, care must be taken to see that Σ_a and L^2 values for each rod kind used in the Self-Consistent calculation are correctly reassigned for the new moderator.

IV. ERROR MESSAGES

Error messages are produced at all levels in the SRL - HERESY I Code. A special subroutine checks the input information for each problem to see that no gross errors have been made that will lead to an error stop at a later time. In each subroutine as it is called, both in Link 1 and in Link 2, certain conditions can arise which will lead to error messages being placed on tape 10 and/or on the on-line printer.

Another class of messages that are printed out give information concerning certain conditions in the program that are not harmful to the successful solution of the problem.

All error messages, information messages, or error conditions displayed in the computer address lights that are built into this code are shown in the following list, subroutine by subroutine, as well as the meaning of the message. The notation in parentheses indicates where the message will be output (i.e. on tape 10 or on-line).

Link 1

A. Main Program 1 (Link 1 executive program)

AN ERROR HAS BEEN DETECTED IN THE INPUT DATA FOR THIS PROBLEM. AN ATTEMPT WILL BE MADE TO CONTINUE WITH SUCCESSIVE PROBLEMS (on-line) - This message appears on the on-line printer following the alphanumeric card for the problem. It means that an error has been detected in the input information and a diagnostic placed on tape 10. Problems that follow will be attempted, but no guarantee is made they will run even if the data are correctly input. Any error detected in main Program 1 will produce this on-line message.

B. Subroutine DATCHK (data checking routine)

1. THE NUMBER OF ROD KINDS IS GREATER THAN THE NUMBER OF ROD TYPES (tape 10) - self-explanatory.
2. THE NUMBER OF ROD TYPES EXCEEDS 50 (tape 10) - self-explanatory.
3. A ROD TYPE HAS ZERO RODS (tape 10) - This can arise from incorrect input data, or errors in inputting data to the geometry generators.
4. A ROD TYPE HAS NO KIND NUMBER ASSIGNED (tape 10) - self-explanatory.

5. THE LATTICE TYPE TO BE GENERATED IS NOT OF A RECOGNIZED TYPE (tape 10) - The mnemonic LAT >2.
6. A ROD KIND IS NOT SPECIFIED FOR THE CRITICALITY SEARCH (tape 10) - Can only occur for criticality searches on γ and η .
7. THE INFORMATION GIVEN FOR THE KERNEL GENERATOR IS INSUFFICIENT (tape 10) - One or more of the three quantities XMESH, XLSQ, or SIGA is ≤ 0 .
8. THE INFORMATION GIVEN FOR ROD KIND (I) IS INSUFFICIENT TO COMPUTE A GAMMA (tape 10) - γ_I is zero; hence calculation of a gamma value is called for using the thermal utilization f_I . One of the two quantities THERU(I) or VCELL(I) is ≤ 0 ; hence calculation of the γ_I value cannot be done.

Following each of the above errors, the following message will be placed on tape 10:

BECAUSE OF THE ERROR(S) DETECTED ABOVE THIS PROBLEM
CANNOT BE CONTINUED. SUCCESSIVE PROBLEMS WILL BE
ATTEMPTED, ALTHOUGH THEY MAY BE AFFECTED BY THE
ERRORS ABOVE (tape 10).

C. Subroutine KERGEN (kernel generator)

1. THE ARGUMENT OF THE E_1 INTEGRAL IN THE KERNEL GENERATOR EXCEEDS 87.3/THE VALUE OF THE INTEGRAL WILL BE SET EQUAL TO $1.0E-30$ AND COMPUTATIONS WILL CONTINUE (tape 10) - An information statement arising due to the fact that the exponential subroutine has a maximum argument of 87.3.
2. THE ARGUMENT OF THE E_1 INTEGRAL IN THE KERNEL GENERATOR IS WITHIN $1.0 E-05$ OF ZERO/THE VALUE OF THE INTEGRAL WILL BE SET EQUAL TO $1.0E+30$ AND COMPUTATIONS WILL CONTINUE (tape 10) - An information statement concerning a condition that could create an accumulator overflow and lead to a HPR 77(EFM) program stop.
3. A TAU(R) VALUE IS LARGER THAN OR EQUAL TO ITS CORRESPONDING TAU VALUE. A KERNEL CANNOT BE GENERATED UNDER THESE CONDITIONS (tape 10) - An error condition arising from placing the effective resonance energy lower than thermal energy.
4. A MACHINE ERROR HAS OCCURRED IN THE KERNEL GENERATOR - RESTART PROBLEM (tape 10 and on-line) - Arises from a branching error by the machine.

D. Subroutine GEOM (geometry executive routine)

1. THE NUMBER OF RODS EXCEEDS 6500 (tape 10) - Arises due to the fact that 6500 rods is the maximum number available for solving problems.

E. Subroutine INGEOM (infinite geometry generator)

1. THE NUMBER OF RODS EXCEEDS 6500 (tape 10) - Same as above for subroutine GEOM.
2. A MACHINE ERROR HAS OCCURRED IN GENERATING AN INFINITE LATTICE - RESTART PROBLEM (tape 10 and on-line) - Arises from a branching error that occurred.

F. Subroutine HEX and SQUARE (hexagonal and square lattice geometry generators)

1. ERROR IN GEOMETRY GENERATOR - INITIAL COORDINATES ARE IN WRONG SECTOR (tape 10) - Arises when the initial X and Y coordinates for a rod type are not in the sector $0^\circ \leq \theta \leq 45^\circ$ for the square lattice generator, or not in the sector $0^\circ \leq \theta \leq 60^\circ$ for the hexagonal lattice generator.

Link 2

G. Subroutine USKER (Bessel's interpolation of kernel)

1. ERROR - KERNEL SIZE IS TOO SMALL FOR LATTICE BEING USED (tape 10) - Arises because the required condition that the distance corresponding to the 250th kernel point be at least two mesh intervals larger than the largest center-to-center rod spacing in the lattice is not met.
2. ERROR - KERNEL MESH IS TOO LARGE (tape 10) - Arises because the required condition that the mesh spacing $XMESH \leq 1/2$ times smallest center-to-center rod spacing is not met.

Following either of the above two error conditions, the following will be placed on tape 10 and on the on-line printer respectively.

- THIS ERROR OCCURS IN FORMING THE MATRICES FROM THE KERNEL AND GEOMETRY/AN ATTEMPT WILL BE MADE TO RUN SUCCESSIVE PROBLEMS, BUT IT IS UNLIKELY THEY WILL RUN (tape 10).
- PROBLEM DID NOT RUN TO END (on-line).

H. Subroutine INVERT (matrix inversion)

1. THE SPECIFIED NUMBER OF ITERATIONS ON MATRIX INVERSION HAS BEEN EXCEEDED/NO. ITERATIONS COMPLETED = $(N1 + 1)$ CONVERGENCE CRITERION = $(CR1)/MATRIX$ TO BE INVERTED FOLLOWS (tape 10) - Arises from the number of iterations through the inversion process requiring more than $N1$ passes. Following this message a listing of each matrix element will appear in the order $M(1,1) \dots M(1,I) \dots M(1,NTYPE)$, $M(2,1) \dots M(2,I) \dots M(2,NTYPE)$, $M(3,1) \dots M(3,I) \dots$

Double spaces will appear in the listing at each point at which the first index changes. Following this listing $N1$ is increased by 10 and the iterations will continue. If the number of iterations exceeds $N1 + 10$ the problem is aborted and conditions are reset to start another problem. If this occurs a line will be placed on the on-line printer which reads PROBLEM DID NOT RUN TO END (on-line).

I. Subroutine EIGEN1 (power iteration to determine maximum eigenvalue)

1. THIS PROBLEM HAS EXCEEDED THE MAXIMUM NUMBER OF ITERATIONS SPECIFIED FOR THE EIGENVALUE ITERATION ($N2$)/CURRENT ITERATION NUMBER = $(N2 + 10M, M = 1, 2 \dots)$, CURRENT EIGENVALUE = $(FK2)/$

(followed by heading)	EIGENVECTOR	V2/A2
	"	"
	"	"
	"	" (tape 10)

Arising from the stated conditions, the above heading is shown first at iteration number $N2 + 11$ and then for every tenth iteration until the iteration number reaches $2*N2$. At this time the problem is aborted and conditions are reset for a new problem. If this aborting occurs, a line is printed on the on-line printer which reads PROBLEM DID NOT RUN TO END (on-line). The values listed under the headings of EIGENVECTOR and V2/A2 are respectively the absorptions computed during the current iteration (un-normalized) and the ratio of absorptions to maximum absorption (i.e. normalized eigenvectors).

J. Subroutine CONVER (convergence test for criticality searches)

1. THE MAXIMUM NUMBER OF ITERATIONS HAVE BEEN EXCEEDED IN THE SEARCH WITHOUT CONVERGING. THIS PROBLEM WILL BE SKIPPED OVER AND SUCCESSIVE PROBLEMS RUN (tape 10) - Arises because the iteration counter for

criticality search has exceeded N3. Due to the time involved in some searches no extra allowance is made for this condition. An edit of the current iteration will follow this statement so that the speed of convergence can be estimated and proper alteration of the problem made. If this occurs a line will be put out by the on-line printer which reads PROBLEM DID NOT RUN TO END (on-line); conditions are then reset for more problems to be run.

K. Subroutine KERGEN (kernel generator for Buckling Search)

1. A MACHINE ERROR HAS OCCURRED IN THE SUBROUTINE KERGEN-RESTART PROBLEM FROM THE BEGINNING (on-line) - Arises from a machine branching error that occurred.

L. Errors Leading to Program Stops

The standard BSS or LUCY loader errors are all in force during the running of both links of the CHAIN PROGRAM. In addition, the following program stops may occur in the course of running a problem which may be interrogated by the address lights.

<u>Storage Register Address Lights</u>	<u>Subroutine in Which Stop Occurs</u>	<u>Reason for Stop</u>
HPR 77	LFM - EFM	Accumulator overflow.
HTR 3,4	CHAIN	Program link called for is not on the designated tape unit. (Push start to transfer back to calling program).
HTR 5,4	XYLIB	File number called for is not on the tape mounted on tape 3. (Push start to transfer back to calling program).
HPR 464	XYLIB	Checksum error in reading tape 2. (Push start to ignore this).
HPR 547	XYLIB	Checksum error in reading tape 3. (Push start to ignore this).
HPR 610	XYLIB	Misplaced EOF or misread of tape 3. (Push start to return to calling program).

V. EXAMPLE PROBLEMS

The objective of this chapter is to give several examples of calculations made with the SRL - HERESY I code so that verification of the code's accuracy may be initially or periodically tested, and to give examples of the input and the output to and from the code in a manner designed to clarify questions that may arise.

A. Finite Lattice Problems

Problem 1. Control Rod at Center of Small Fuel Lattice

Because of its rapid running time, this problem is the basic debugging problem for SRL - HERESY I. The lattice is shown in Figure 1. It consists of a control rod situated at the center of a small array of fuel assemblies. The lattice is a square type with a pitch of 10.0 inches (25.4 cm). It has 8-fold symmetry, and hence can be generated by the square lattice generator. The problem was treated without axial leakage ($B_z^2 = 0$), and the values of $B(I)$, $TAU(I)$, and $TAUR(I)$ within the code were used. An infinite moderator was used. GAMMA values were computed by the self-consistent method from thermal utilizations. SMF's were specified internally by setting rod radii to zero.

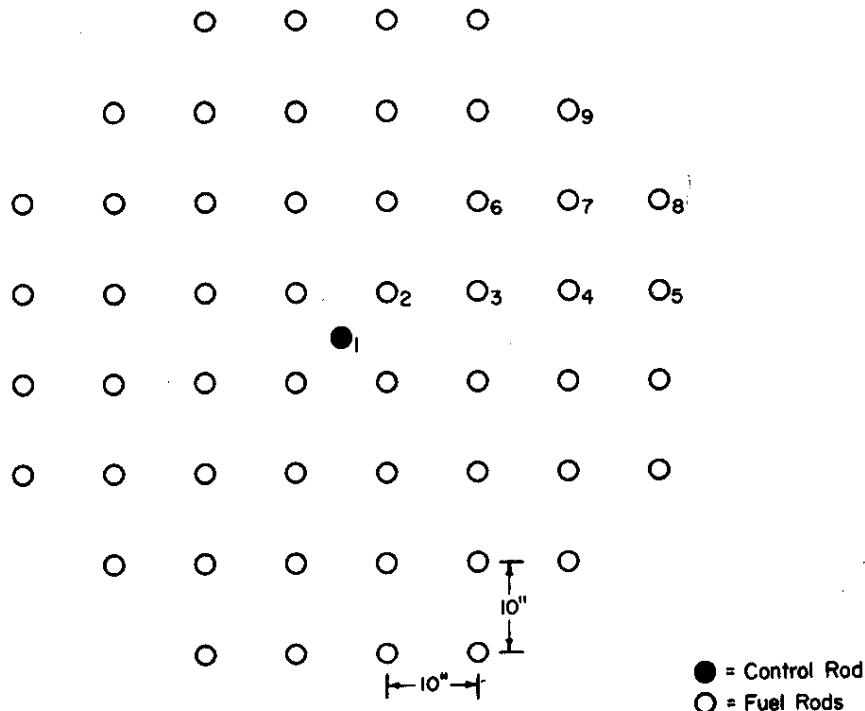


FIG. 1 LATTICE FOR SMALL FINITE TEST PROBLEM

The basic problem was run with no criticality searches, then each of the criticality searches was applied in turn to the basic problem. A listing and plot of geometry were obtained, but are not shown since they were used only to verify the correctness of the geometry as generated. Figure 2 shows a summary, as produced by the code, of the input data for the basic problem. Figure 3 shows the solution to the problem.

FIGURE 2

1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**BASIC PROBLEM

SUMMARY OF HERESY INPUT DATA

FINITE/INFINITE OPTION = 0 KERNEL OPTION = 1 INPLT OPTIONS COORDINATE OPTION = 1 PURE THERMAL KERNEL OPTION = 0
LATTICE TYPE = 1 GEOMETRY PRINT = 0 GEOMETRY PLOT = 0 GEOMETRY PUNCH = 0

MATRIX INVERSION ITERATIONS = 20 EIGENVALUE ITERATIONS = 200 CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001 EIGENVALUE CONVERGENCE = .000010 CRITICALITY CONVERGENCE = .000010

CONVERGENCE CRITERIA

KERNEL DATA

NUMBER OF GRID POINTS = 250 MESH SIZE = 1.000 CM
B(1) = 0.56 B(2) = 0.44 B(3) = 0.
TAU(1) = 154.30 TAU(2) = 68.30 TAU(3) = 0.
TAUR(1) = 126.50 TAUR(2) = 40.50 TAUR(3) = 0.
B SQUARE = . L SQUARE = 12626.7000 MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = C. RHO(MIN) = 0.

DATA FOR EACH ROD KIND

ROD KIND	GAMMA	SMF	1./(GAM+SMF)	ETA	RES. CCEF.	LSQ	SIGMA(A)	F	CELL VOL.	RADIUS
1	1.14909	0.10101	0.799934	0.	0.	12626.7	.0000696	0.98544	322.48000	0.
2	0.74668	0.10101	1.179676	1.30530	58.8260	12626.7	.0000696	0.98590	645.16000	0.

DATA FOR EACH ROD TYPE

ROD TYPE	ROD KIND	NO. RODS	GAMMA	SMF	1./(GAMMA+SMF)
1	1	1	1.14909	0.10101	0.799934
2	2	4	0.74668	0.10101	1.179676
3	2	8	0.74668	0.10101	1.179676
4	2	4	0.74668	0.10101	1.179676
5	2	8	0.74668	0.10101	1.179676
6	2	8	0.74668	0.10101	1.179676
7	2	4	0.74668	0.10101	1.179676
8	2	8	0.74668	0.10101	1.179676
9	2	8	0.74668	0.10101	1.179676

FIGURE 3

NO EIGENVALUE ITERATIONS EIGENVALUE PREVIOUS EIGENVALUE
21 1.0828493 1.0828522

ABSORPTIONS FOR EACH ROD TYPE

ROD TYPE	NO RODS/TYPE	ROD SMF	GAMMA	ETA	A	THERMAL ABSRB	RESONANCE ABSRB
1	1	0.10101	1.149093	0.	0.	0.417634	0.
2	4	0.10101	0.746680	1.30530	58.8260	0.943900	0.121672
3	8	0.10101	0.746680	1.30530	58.8260	1.000000	0.125911
4	4	0.10101	0.746680	1.30530	58.8260	0.975386	0.122873
5	8	0.10101	0.746680	1.30530	58.8260	0.936836	0.118093
6	8	0.10101	0.746680	1.30530	58.8260	0.892689	0.111465
7	4	0.10101	0.746680	1.30530	58.8260	0.827345	0.091342
8	8	0.10101	0.746680	1.30530	58.8260	0.815398	0.094986
9	8	0.10101	0.746680	1.30530	58.8260	0.782690	0.085284

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.08284929 AVERAGE ROD ABSORPTION = 0.8834916

AVERAGE ETA = 1.2936520 THERMAL UTILIZATION = 0.98589589 RESONANCE ESCAPE PROBABILITY = 0.89936797

APPRX MODERATOR THERMAL ABSORPTIONS = 0.66987 TOTAL LEAKAGE = 2.8166 NON-LEAKAGE PROBABILITY = 0.9440176

AVERAGE ABSORPTIONS FOR EACH ROD KIND

ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.417634	0.
2	52	0.892450	0.108259

Using this basic problem as a starting point for each criticality search is next in order. Figure 4 shows the summary of input data produced by the code for a search over the η value of rod kind 2. It was specified that this problem edit each iteration of the search, and these edits are shown in Figures 5, 6, and 7. Figure 8 shows the input data summary for a gamma criticality search and the summary of the search that is provided when each iteration is not edited. The edited final solution is shown in Figure 9. A new thermal utilization is computed in this case; however, if γ is not originally obtained from a self-consistent calculation, the re-computation of the f corresponding to the new γ value is not possible.

FIGURE 4

```

1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**ETA SEARCH

SUMMARY OF HERESY INPUT DATA

FINITE/INFINITE OPTION = -0      KERNEL OPTION = -0      INPUT OPTIONS
                                COORDINATE OPTION = -0      PURE THERMAL KERNEL OPTION = 0

MATRIX INVERSION ITERATIONS = 20  EIGENVALUE ITERATIONS = 200  CONVERGENCE CRITERIA
INVERSION CONVERGENCE = .000001  EIGENVALUE CONVERGENCE = .000010  CRITICALITY SEARCH ITERATIONS = 25
                                CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 250  MESH SIZE = 1.000 CM
B(1) = 0.56  B(2) = 0.44  B(3) = 0.
TAU(1) = 154.30  TAU(2) = 68.30  TAU(3) = 0.
TAUR(1) = 126.50  TAUR(2) = 40.50  TAUR(3) = 0.
B SQUARE = .  L SQUARE = 12626.7000  MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = 0.  RHO(MIN) = 0.

CRITICALITY SEARCH ON ETA - EIGENVALUE TO BE SEARCHED FOR = 1.0000000
ROD KIND TO BE SEARCHED ON = 2

DATA FOR EACH ROD KIND
ROD KIND  GAMMA  SMF  1./(GAM+SMF)  ETA  RES. COEF.  LSQ  SIGMA(A)  F  CELL VOL.  RADIUS
1  1.14909  0.10101  0.799934  0.  0.  12626.7  .0000696  0.98544  322.48000  0.
2  0.74668  0.10101  1.179676  1.30530  58.8260  12626.7  .0000696  0.98590  645.16000  0.

DATA FOR EACH ROD TYPE
ROD TYPE  ROD KIND  NO. RODS  GAMMA  SMF  1./(GAMMA+SMF)
1  1  1  1.14909  0.10101  0.799934
2  2  4  0.74668  0.10101  1.179676
3  2  8  0.74668  0.10101  1.179676
4  2  4  0.74668  0.10101  1.179676
5  2  8  0.74668  0.10101  1.179676
6  2  8  0.74668  0.10101  1.179676
7  2  4  0.74668  0.10101  1.179676
8  2  8  0.74668  0.10101  1.179676
9  2  8  0.74668  0.10101  1.179676

```

FIGURE 5

CRITICALITY SEARCH ITERATION NO 0

NO EIGENVALUE ITERATIONS		EIGENVALUE	PREVIOUS EIGENVALUE		ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB
15		1.0828390	1.0828367		GAMMA	ETA	A		
ROD TYPE	NO RODS/TYPE	ROD SMF							
1	1	0.10101	1.149093	0.	0.	0.417642	0.		
2	4	0.10101	0.746680	1.30530	58.8260	0.943913	0.121674		
3	8	0.10101	0.746680	1.30530	58.8260	1.000000	0.125912		
4	4	0.10101	0.746680	1.30530	58.8260	0.975375	0.122872		
5	8	0.10101	0.746680	1.30530	58.8260	0.936817	0.118091		
6	8	0.10101	0.746680	1.30530	58.8260	0.892664	0.111463		
7	4	0.10101	0.746680	1.30530	58.8260	0.827315	0.091339		
8	8	0.10101	0.746680	1.30530	58.8260	0.815367	0.094984		
9	8	0.10101	0.746680	1.30530	58.8260	0.782659	0.085281		

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.08283895 AVERAGE ROD ABSORPTION = 0.8834738

AVERAGE ETA = 1.2936575 THERMAL UTILIZATION = 0.98589589 RESONANCE ESCAPE PROBABILITY = 0.89936788

APPROX MODERATOR THERMAL ABSORPTIONS = 0.66986 TOTAL LEAKAGE = 2.8170 NON-LEAKAGE PROBABILITY = 0.9440090

AVERAGE ABSORPTIONS FOR EACH ROD KIND			
ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.417642	0.
2	52	0.892432	0.108258

FIGURE 6

CRITICALITY SEARCH ITERATION NO 1

NO EIGENVALUE ITERATIONS		EIGENVALUE	PREVIOUS EIGENVALUE		ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB
1		0.9931400	0.9931391		GAMMA	ETA	A		
ROD TYPE	NO RODS/TYPE	ROD SMF							
1	1	0.10101	1.149093	0.	0.	0.417640	0.		
2	4	0.10101	0.746680	1.19717	58.8260	0.943910	0.121673		
3	8	0.10101	0.746680	1.19717	58.8260	1.000000	0.125912		
4	4	0.10101	0.746680	1.19717	58.8260	0.975378	0.122872		
5	8	0.10101	0.746680	1.19717	58.8260	0.936821	0.118092		
6	8	0.10101	0.746680	1.19717	58.8260	0.892670	0.111463		
7	4	0.10101	0.746680	1.19717	58.8260	0.827322	0.091340		
8	8	0.10101	0.746680	1.19717	58.8260	0.815374	0.094985		
9	8	0.10101	0.746680	1.19717	58.8260	0.782667	0.085282		

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 0.99314001 AVERAGE ROD ABSORPTION = 0.8834781

AVERAGE ETA = 1.1864924 THERMAL UTILIZATION = 0.98589590 RESONANCE ESCAPE PROBABILITY = 0.89936790

APPROX MODERATOR THERMAL ABSORPTIONS = 0.66986 TOTAL LEAKAGE = 2.8169 NON-LEAKAGE PROBABILITY = 0.9440110

AVERAGE ABSORPTIONS FOR EACH ROD KIND			
ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.417640	0.
2	52	0.892437	0.108258

FIGURE 7

CRITICALITY SEARCH ITERATION NO 2

NO EIGENVALUE ITERATIONS		EIGENVALUE	PREVIOUS EIGENVALUE		ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB
0		1.0000008	1.0000000		GAMMA	ETA	A		
ROD TYPE	NO RODS/TYPE	ROD SMF							
1	1	0.10101	1.149093	0.	0.	0.417640	0.		
2	4	0.10101	0.746680	1.20544	58.8260	0.943909	0.121673		
3	8	0.10101	0.746680	1.20544	58.8260	1.000000	0.125912		
4	4	0.10101	0.746680	1.20544	58.8260	0.975378	0.122872		
5	8	0.10101	0.746680	1.20544	58.8260	0.936823	0.118092		
6	8	0.10101	0.746680	1.20544	58.8260	0.892671	0.111463		
7	4	0.10101	0.746680	1.20544	58.8260	0.827324	0.091340		
8	8	0.10101	0.746680	1.20544	58.8260	0.815376	0.094985		
9	8	0.10101	0.746680	1.20544	58.8260	0.782669	0.085282		

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.00000083 AVERAGE ROD ABSORPTION = 0.8834792

AVERAGE ETA = 1.1946882 THERMAL UTILIZATION = 0.98589590 RESONANCE ESCAPE PROBABILITY = 0.89936791

APPROX MODERATOR THERMAL ABSORPTIONS = 0.66986 TOTAL LEAKAGE = 2.8168 NON-LEAKAGE PROBABILITY = 0.9440117

AVERAGE ABSORPTIONS FOR EACH ROD KIND			
ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.417640	0.
2	52	0.892438	0.108258

FIGURE 8

1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**GAMMA SEARCH

SUMMARY OF HERESY INPUT DATA

FINITE/INFINITE OPTION = -0 KERNEL OPTION = -0 INPUT OPTIONS
COORDINATE OPTION = -0 PURE THERMAL KERNEL OPTION = 0

CONVERGENCE CRITERIA
MATRIX INVERSION ITERATIONS = 20 EIGENVALUE ITERATIONS = 200 CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001 EIGENVALUE CONVERGENCE = .000010 CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 250 MESH SIZE = 1.000 CM
B(1) = 0.56 B(2) = 0.44 B(3) = 0.
TAU(1) = 154.30 TAU(2) = 68.30 TAU(3) = 0.
TAUR(1) = 126.50 TAUR(2) = 40.50 TAUR(3) = 0.
B SQUARE = . L SQUARE = 12626.7000 MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = C. RHO(MIN) = 0.

CRITICALITY SEARCH ON GAMMA - EIGENVALUE TO BE SEARCHED FOR = 1.0000000
ROD KIND TO BE SEARCHED ON = 2

DATA FOR EACH ROD KIND

ROD KIND	GAMMA	SMF	1./(GAM+SMF)	ETA	RES. COEF.	LSQ	SIGMA(A)	F	CELL VOL.	RADIUS
1	1.14909	0.10101	0.799934	0.	0.	12626.7	.0000696	0.98544	322.48000	0.
2	0.74668	0.10101	1.179676	1.30530	58.8260	12626.7	.0000696	0.98590	645.16000	0.

DATA FOR EACH ROD TYPE

ROD TYPE	ROD KIND	NO. RODS	GAMMA	SMF	1./(GAMMA+SMF)
1	1	1	1.14909	0.10101	0.799934
2	2	4	0.74668	0.10101	1.179676
3	2	8	0.74668	0.10101	1.179676
4	2	4	0.74668	0.10101	1.179676
5	2	8	0.74668	0.10101	1.179676
6	2	8	0.74668	0.10101	1.179676
7	2	4	0.74668	0.10101	1.179676
8	2	8	0.74668	0.10101	1.179676
9	2	8	0.74668	0.10101	1.179676

SUMMARY OF CRITICALITY SEARCH
EIGENVALUE BEING SEARCHED FOR = 1.0000000

ITERATION NO	CURRENT EIGENVALUE	PREVIOUS EIGENVALUE	CURRENT GAMMA VALUE	ROD KIND BEING SEARCHED ON
1	1.0947788	1.0828446	0.6818	2
2	0.9079699	1.0947788	1.8940	2
3	1.0335259	0.9079699	1.0307	2
4	1.0096160	1.0335259	1.1785	2
5	0.9985274	1.0096160	1.2495	2
6	1.0000483	0.9985274	1.2396	2
7	1.0000032	1.0000483	1.2399	2

FIGURE 9

NO EIGENVALUE ITERATIONS EIGENVALUE PREVIOUS EIGENVALUE
1 1.0000032 1.0000027

ABSORPTIONS FOR EACH ROD TYPE

ROD TYPE	NO RODS/TYP	ROD SMF	GAMMA	ETA	A	THERMAL ABSRB	RESONANCE ABSRB
1	1	0.10101	1.149093	C.	0.	0.936613	0.
2	4	0.10101	1.239935	1.30530	58.8260	0.932519	0.130471
3	8	0.10101	1.239935	1.30530	58.8260	1.000000	0.136275
4	4	0.10101	1.239935	1.30530	58.8260	0.978130	0.133267
5	8	0.10101	1.239935	1.30530	58.8260	0.940163	0.128065
6	8	0.10101	1.239935	1.30530	58.8260	0.893862	0.120481
7	4	0.10101	1.239935	1.30530	58.8260	0.807285	0.096972
8	8	0.10101	1.239935	1.30530	58.8260	0.804716	0.101615
9	8	0.10101	1.239935	1.30530	58.8260	0.761840	0.090280

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.00000319 AVERAGE ROD ABSORPTION = 0.8870377

AVERAGE ETA = 1.2792953 THERMAL UTILIZATION = 0.96522497 RESONANCE ESCAPE PROBABILITY = 0.89929765

APPROX MODERATOR THERMAL ABSORPTIONS = 1.69378 TOTAL LEAKAGE = 5.3800 NON-LEAKAGE PROBABILITY = 0.9005308

AVERAGE ABSORPTIONS FOR EACH ROD KIND

ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.936613	0.
2	52	0.886084	0.116473

THE VALUE OF THERMAL UTILIZATION CORRESPONDING TO A GAMMA VALUE OF 1.239935 FOR ROD TYPE 2 IS 0.96482265

The last problem of this set of four problems is that of a buckling search starting from the basic problem. This problem requires a longer running time than the previous three combined. Figure 10 summarizes the input data and search summary, and Figure 11 shows the converged solution.

FIGURE 10

1 SMALL TEST PROBLEM FOR FINITE LATTICE==SRL-HERESY I==BUCKLING SEARCH

SUMMARY OF HERESY INPUT DATA

FINITE/INFINITE OPTION = -0 KERNEL OPTION = -0 INPUT OPTIONS
COORDINATE OPTION = -0 PURE THERMAL KERNEL OPTION = 0

CONVERGENCE CRITERIA
MATRIX INVERSION ITERATIONS = 20 EIGENVALUE ITERATIONS = 200 CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001 EIGENVALUE CONVERGENCE = .000010 CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 250 MESH SIZE = 1.000 CM
B(1) = 0.56 B(2) = 0.44 B(3) = 0.
TAU(1) = 154.30 TAU(2) = 68.30 TAU(3) = 0.
TAUR(1) = 126.50 TAUR(2) = 40.50 TAUR(3) = 0.
B SQUARE = . L SQUARE = 12626.7000 MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = C. RHO(MIN) = 0.

CRITICALITY SEARCH ON BUCKLING - EIGENVALUE TO BE SEARCHED FOR = 1.0000000

DATA FOR EACH ROD KIND

ROD KIND	GAMMA	SMF	1./(GAM+SMF)	ETA	RES. COEF.	LSQ	SIGMA(A)	F	CELL VOL.	RADIUS
1	1.14909	0.10101	0.799934	0.	0.	12626.7	.0000696	0.98544	322.48000	0.
2	0.74668	0.10101	1.179676	1.30530	58.8260	12626.7	.0000696	0.98590	645.16000	0.

DATA FOR EACH ROD TYPE

ROD TYPE	ROD KIND	NO. RODS	GAMMA	SMF	1./(GAMMA+SMF)
1	1	1	1.14909	0.10101	0.799934
2	2	4	0.74668	0.10101	1.179676
3	2	8	0.74668	0.10101	1.179676
4	2	4	0.74668	0.10101	1.179676
5	2	8	0.74668	0.10101	1.179676
6	2	8	0.74668	0.10101	1.179676
7	2	4	0.74668	0.10101	1.179676
8	2	8	0.74668	0.10101	1.179676
9	2	8	0.74668	0.10101	1.179676

SUMMARY OF CRITICALITY SEARCH
EIGENVALUE BEING SEARCHED FOR = 1.0000000

ITERATION NO	CURRENT EIGENVALUE	PREVIOUS EIGENVALUE	CURRENT BUCKLING VALUE
1	0.9837358	1.0028429	0.00025000
2	0.9975945	0.9837358	0.00020897
3	1.0000624	0.9975945	0.00020185
4	1.0000096	1.0000624	0.00020203

FIGURE 11

NO EIGENVALUE ITERATIONS EIGENVALUE PREVIOUS EIGENVALUE
2 1.0238174 1.0238195

ABSORPTIONS FOR EACH ROD TYPE

ROD TYPE	NO RODS/TYPE	ROD SMF	GAMMA	ETA	A	THERMAL ABSRB	RESONANCE ABSRB
1	1	0.10101	1.034394	0.	0.	0.432026	0.
2	4	0.10101	0.631981	1.30530	58.8260	0.968079	0.134405
3	8	0.10101	0.631981	1.30530	58.8260	1.000000	0.135665
4	4	0.10101	0.631981	1.30530	58.8260	0.950584	0.129016
5	8	0.10101	0.631981	1.30530	58.8260	0.889412	0.120772
6	8	0.10101	0.631981	1.30530	58.8260	0.821611	0.110742
7	4	0.10101	0.631981	1.30530	58.8260	0.706140	0.085673
8	8	0.10101	0.631981	1.30530	58.8260	0.701128	0.089272
9	8	0.10101	0.631981	1.30530	58.8260	0.646464	0.077601

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.00000955 AVERAGE ROD ABSORPTION = 0.8188709

AVERAGE ETA = 1.2923064 THERMAL UTILIZATION = 0.98589543 RESONANCE ESCAPE PROBABILITY = 0.89892645

APPROX MODERATOR THERMAL ABSORPTIONS = 0.62090 TOTAL LEAKAGE = 6.3959 NON-LEAKAGE PROBABILITY = 0.8731395

AVERAGE ABSORPTIONS FOR EACH ROD KIND

ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	1	0.432026	0.
2	52	0.826310	0.109015

A listing of the input cards for these problems is shown in Figure 12.

FIGURE 12

```

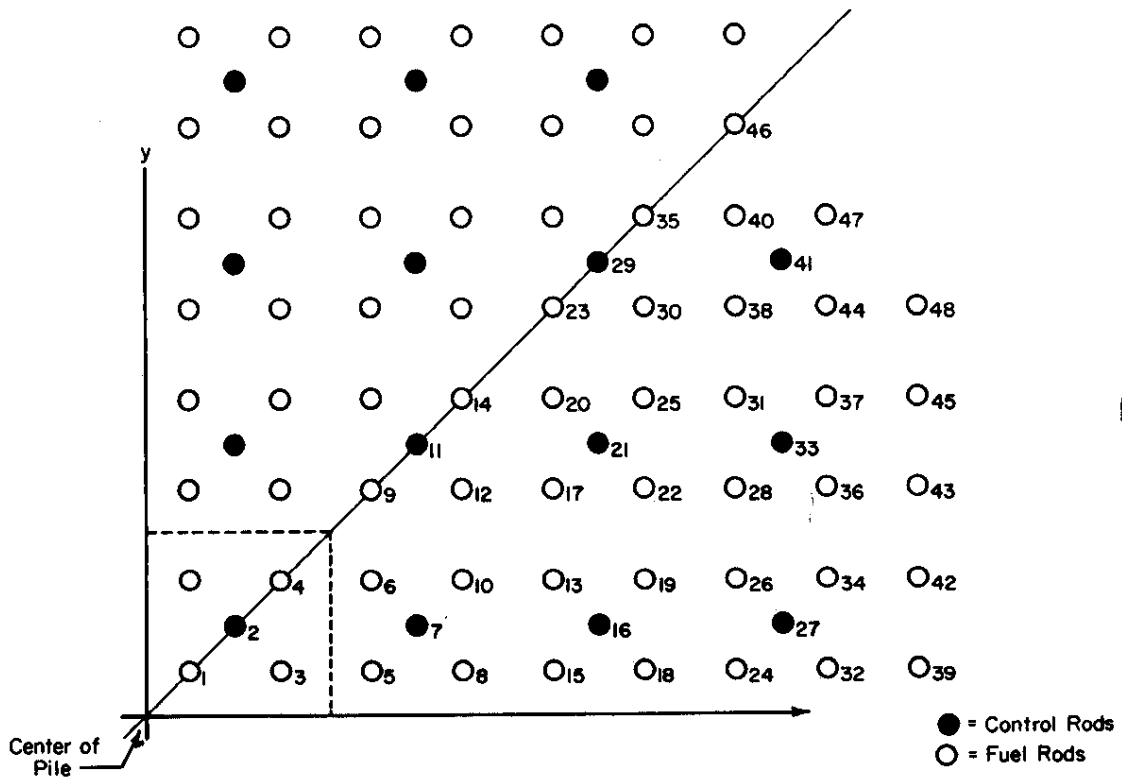
1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**BASIC PROBLEM
  3      1      0      1      1
  5      7      250    0      1      1      1
  5     12      1      1      1      1
  2     17      1      1
  2     63      9      2
  5    115      1      2      2      2
1  4    120      2      2      2      2
  1      4 12.7
  4     10 1.0      0.0      12626.7      .00006963
  2    125 322.48    645.16
  2    225 .98544    .98590
  2    275 0.0      1.3053
  2    325 0.0      58.826
  5    425 0.0      1.0      3.0      3.0      5.0
  4    430 5.0      5.0      7.0      7.0
  5    575 0.0      1.0      1.0      3.0      1.0
1  4    580 3.0      5.0      1.0      3.0
1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL HERESY I**XYLIB TEST
1  2    224      1      1
1  2    175
1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**GAMMA SEARCH
  5      1
  2    224      1
1  2    165      2      2
  1      4 1.0
1  1    276 1.3053
1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**ETA SEARCH
1  3    165      1      2      1
  2    175
  2    225 .98544    .9859
1 SMALL TEST PROBLEM FOR FINITE LATTICE**SRL-HERESY I**BUCKLING SEARCH
1  5    165      3
1  1    276 1.3053
C

```

Problem 2. Large Lattice of Fuel and Control Rods

This problem is an extension of the previous problem to that of determining the control rod effects in a large lattice of fuel assemblies. Total rod types in this case is 48, of which 8 are control rods, as shown in Figure 13. Once more 8-fold symmetry is exhibited by the square lattice taken at a 10-inch (25.4-cm) pitch; hence, the internal lattice generator was used. Parameters for input were prepared in exactly the same manner as for Problem 1, and it will be noted that the fuel is the same as for Problem 1.

FIGURE 13 Sector of Large Finite Lattice



EXAMPLE PROBLEM

The input data for Problem 1 is summarized in Figure 14.

FIGURE 14

```

1 LARGE TEST PROBLEM FOR FINITE LATTICE ** SRL-HERESY I

SUMMARY OF HERESY INPUT DATA

INPUT OPTIONS
FINITE/INFINITE OPTION = 0   KERNEL OPTION = 1   COORDINATE OPTION = 1   PURE THERMAL KERNEL OPTION = 0
LATTICE TYPE = 1   GEOMETRY PRINT = 1   GEOMETRY PLOT = 1   GEOMETRY PUNCH = 0

CONVERGENCE CRITERIA
MATRIX INVERSION ITERATIONS = 20   EIGENVALUE ITERATIONS = 200   CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001   EIGENVALUE CONVERGENCE = .000010   CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 250   MESH SIZE = 2.500 CM
B(1) = 0.56   B(2) = 0.44   B(3) = 0.
TAU(1) = 154.30   TAU(2) = 68.30   TAU(3) = 0.
TAUR(1) = 126.50   TAUR(2) = 40.50   TAUR(3) = 0.
B SQUARE = .   L SQUARE = 12626.7000   MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = -0.   RHO(MIN) = -0.

DATA FOR EACH ROD KIND
ROD KIND   GAMMA   SMF   1./(GAMMA+SMF)   ETA   RES. COEF.   LSQ   SIGMA(A)   F   CELL VOL.   RADIUS
1   0.74668   0.10101   1.179676   1.30530   58.8260   12626.7   .0000696   0.98590   645.16000   0.
2   4.17239   0.10101   0.234006   0.   0.   12626.7   .0000696   0.92365   322.48000   0.

DATA FOR EACH ROD TYPE
ROD TYPE   ROD KIND   NO. RODS   GAMMA   SMF   1./(GAMMA+SMF)
1   1   4   0.74668   0.10101   1.179676
2   2   4   4.17239   0.10101   0.234006
3   1   8   0.74668   0.10101   1.179676
4   1   4   0.74668   0.10101   1.179676
5   1   8   0.74668   0.10101   1.179676
6   1   8   0.74668   0.10101   1.179676
7   2   8   4.17239   0.10101   0.234006
8   1   8   0.74668   0.10101   1.179676
9   1   4   0.74668   0.10101   1.179676
10   1   8   0.74668   0.10101   1.179676
11   2   8   4.17239   0.10101   0.234006
12   1   8   0.74668   0.10101   1.179676
13   1   8   0.74668   0.10101   1.179676
14   1   8   0.74668   0.10101   1.179676
15   1   8   0.74668   0.10101   1.179676
16   2   8   4.17239   0.10101   0.234006
17   1   8   0.74668   0.10101   1.179676
18   1   8   0.74668   0.10101   1.179676
19   1   8   0.74668   0.10101   1.179676
20   1   8   0.74668   0.10101   1.179676
21   2   8   4.17239   0.10101   0.234006
22   1   8   0.74668   0.10101   1.179676
23   1   4   0.74668   0.10101   1.179676
24   1   8   0.74668   0.10101   1.179676
25   1   8   0.74668   0.10101   1.179676
26   1   8   0.74668   0.10101   1.179676
27   2   8   4.17239   0.10101   0.234006
28   1   8   0.74668   0.10101   1.179676
29   2   4   4.17239   0.10101   0.234006
30   1   8   0.74668   0.10101   1.179676
31   1   8   0.74668   0.10101   1.179676
32   1   8   0.74668   0.10101   1.179676
33   2   8   4.17239   0.10101   0.234006
34   1   8   0.74668   0.10101   1.179676
35   1   4   0.74668   0.10101   1.179676
36   1   8   0.74668   0.10101   1.179676
37   1   8   0.74668   0.10101   1.179676
38   1   8   0.74668   0.10101   1.179676
39   1   8   0.74668   0.10101   1.179676
40   1   8   0.74668   0.10101   1.179676
41   2   8   4.17239   0.10101   0.234006
42   1   8   0.74668   0.10101   1.179676
43   1   8   0.74668   0.10101   1.179676
44   1   8   0.74668   0.10101   1.179676
45   1   8   0.74668   0.10101   1.179676
46   1   4   0.74668   0.10101   1.179676
47   1   8   0.74668   0.10101   1.179676
48   1   8   0.74668   0.10101   1.179676

```

The solution to Problem 2 is shown in Figure 15. The number of iterations for the eigenvalue power iteration have markedly increased over the 9 rod types of Problem 1. This, however, is expected due simply to the increase of rod types. In fact, a simple ratio of rod types shows:

$$\frac{\text{No. types}_2}{\text{No. types}_1} = \frac{48}{9} = 5.3333 \approx \frac{\text{No. iterations}_2}{\text{No. iterations}_1} = \frac{101}{21} = 4.8095$$

from which the time increase is evidently fairly linear with number of rod types. Not shown by the output is the increase in time required to form the matrix elements. This should be proportional to the number of matrix elements formed times the total number of rods in the lattice:

$$\frac{(\text{No. elements } 2)(\text{No. rods } 2)}{(\text{No. elements } 1)(\text{No. rods } 1)} = \frac{(48)^2(352)}{(9)^2(53)} = 188.9$$

FIGURE 15

NO EIGENVALUE ITERATIONS 101		EIGENVALUE 1.1031582	PREVIOUS EIGENVALUE 1.1031590					
ROD TYPE	NO RODS/TYPE	ROD SMF	ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB	
			GAMMA	ETA	A			
1	4	0.10101	0.746680	1.30530	58.8260	1.000000	0.124763	
2	4	0.10101	4.172394	0.	0.	0.097868	0.	
3	8	0.10101	0.746680	1.30530	58.8260	0.980397	0.122319	
4	4	0.10101	0.746680	1.30530	58.8260	0.957996	0.119513	
5	8	0.10101	0.746680	1.30530	58.8260	0.946001	0.118022	
6	4	0.10101	0.746680	1.30530	58.8260	0.910903	0.116074	
7	8	0.10101	4.172394	0.	0.	0.090236	0.	
8	8	0.10101	0.746680	1.30530	58.8260	0.894804	0.113923	
9	4	0.10101	0.746680	1.30530	58.8260	0.860294	0.110252	
10	8	0.10101	0.746680	1.30530	58.8260	0.777796	0.123812	
11	8	0.10101	4.172394	0.	0.	0.081147	0.	
12	8	0.10101	0.746680	1.30530	58.8260	0.770296	0.122271	
13	8	0.10101	0.746680	1.30530	58.8260	0.764106	0.118380	
14	8	0.10101	0.746680	1.30530	58.8260	0.825550	0.111798	
15	8	0.10101	0.746680	1.30530	58.8260	0.756574	0.122955	
16	8	0.10101	4.172394	0.	0.	0.077800	0.	
17	8	0.10101	0.746680	1.30530	58.8260	0.867611	0.102111	
18	8	0.10101	0.746680	1.30530	58.8260	0.829320	0.105578	
19	8	0.10101	0.746680	1.30530	58.8260	0.840363	0.101647	
20	8	0.10101	0.746680	1.30530	58.8260	0.841596	0.101997	
21	8	0.10101	4.172394	0.	0.	0.082167	0.	
22	8	0.10101	0.746680	1.30530	58.8260	0.819485	0.101046	
23	4	0.10101	0.746680	1.30530	58.8260	0.803758	0.099230	
24	8	0.10101	0.746680	1.30530	58.8260	0.792883	0.098896	
25	8	0.10101	0.746680	1.30530	58.8260	0.791338	0.098810	
26	8	0.10101	0.746680	1.30530	58.8260	0.784054	0.097838	
27	8	0.10101	4.172394	0.	0.	0.075375	0.	
28	8	0.10101	0.746680	1.30530	58.8260	0.765694	0.095521	
29	4	0.10101	4.172394	0.	0.	0.075121	0.	
30	8	0.10101	0.746680	1.30530	58.8260	0.753404	0.094021	
31	8	0.10101	0.746680	1.30530	58.8260	0.734786	0.091678	
32	8	0.10101	0.746680	1.30530	58.8260	0.725331	0.090617	
33	8	0.10101	4.172394	0.	0.	0.071332	0.	
34	8	0.10101	0.746680	1.30530	58.8260	0.715163	0.089354	
35	4	0.10101	0.746680	1.30530	58.8260	0.700009	0.087429	
36	8	0.10101	0.746680	1.30530	58.8260	0.694434	0.086751	
37	8	0.10101	0.746680	1.30530	58.8260	0.660562	0.082562	
38	8	0.10101	0.746680	1.30530	58.8260	0.691595	0.086294	
39	8	0.10101	0.746680	1.30530	58.8260	0.659965	0.075347	
40	8	0.10101	0.746680	1.30530	58.8260	0.635764	0.078599	
41	8	0.10101	4.172394	0.	0.	0.062264	0.	
42	8	0.10101	0.746680	1.30530	58.8260	0.649481	0.074159	
43	8	0.10101	0.746680	1.30530	58.8260	0.627870	0.071709	
44	8	0.10101	0.746680	1.30530	58.8260	0.615239	0.076105	
45	8	0.10101	0.746680	1.30530	58.8260	0.593526	0.067842	
46	4	0.10101	0.746680	1.30530	58.8260	0.598279	0.064728	
47	8	0.10101	0.746680	1.30530	58.8260	0.578067	0.062829	
48	8	0.10101	0.746680	1.30530	58.8260	0.556709	0.059473	

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.10315818 AVERAGE ROD ABSORPTION = 0.6349238

AVERAGE ETA = 1.2760129 THERMAL UTILIZATION = 0.98441141 RESONANCE ESCAPE PROBABILITY = 0.89282393

APPROX MODERATOR THERMAL ABSORPTIONS = 3.53911 TOTAL LEAKAGE = 3.7739 NON-LEAKAGE PROBABILITY = 0.9836492

AVERAGE ABSORPTIONS FOR EACH ROD KIND			
ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	288	0.758607	0.096203
2	64	0.078352	0.

This problem is a reasonable example of the use of SRL - HERESY I in practical problems. The tendency to approach nearly 50 rod types in most problems has been observed in practice, and arises from attempting to approximate real lattices to as fine a detail as possible.

Data cards for this problem are shown in Figure 16.

FIGURE 16

```

1 LARGE TEST PROBLEM FOR FINITE LATTICE ** SRL-HERESY I
  1      0      1      1      0      0
  7      250     0      1      1      1
 12      1      1      1      1      1
 17      1      1      1      1      1
 22      1      1      1      1      1
 27      1      1      1      1      1
 32      1      1      1      1      1
 37      1      1      1      1      1
 42      1      1      1      1      1
 47      1      1      1      1      1
 52      1      1      1      1      1
 57      1      1      1      1      1
 63      48      2      1      1      1
115      1      2      1      1      1
120      1      2      1      1      1
125      2      1      1      1      1
130      2      1      1      1      1
135      2      1      1      1      1
140      1      2      1      2      1
145      1      1      2      1      1
150      1      1      1      1      1
155      2      1      1      1      1
160      1      1      1      1      1
  4 12.7
 10 2.5      12626.7      .00006963
125 645.16      322.48
225 .9859      .92365
275 1.3053      0.0
325 58.826      0.0
425 1.0      2.0      3.0      3.0      5.0
430 5.0      6.0      7.0      5.0      7.0
435 6.0      7.0      9.0      7.0      9.0
440 10.0      9.0      11.0      11.0      9.0
445 10.0      11.0      9.0      13.0      11.0
450 13.0      14.0      13.0      10.0      11.0
455 13.0      15.0      14.0      15.0      11.0
460 15.0      15.0      13.0      17.0      13.0
465 14.0      17.0      17.0      15.0      17.0
470 13.0      15.0      17.0
575 1.0      2.0      1.0      3.0      1.0
580 3.0      2.0      1.0      5.0      3.0
585 5.0      4.0      2.0      6.0      1.0
590 2.0      5.0      1.0      3.0      7.0
595 6.0      5.0      9.0      1.0      7.0
600 3.0      2.0      5.0      10.0      9.0
605 7.0      1.0      6.0      3.0      11.0
610 5.0      7.0      9.0      1.0      11.0
615 10.0      3.0      5.0      9.0      7.0
620 13.0      11.0      9.0

```

B. Infinite Lattice Problems

Problem 3. Fundamental Methods Associates, Inc. Test Problem II

As the first example of the infinite lattice problem, the test problem formulated by Fundamental Methods Associates, Inc. described in Reference 1, p 59, was run on the SRL - HERESY I code. Input data are summarized in Figure 17 by the code, and reference to the original work cited above shows its source. (Note the change of the maximum rod radius to 260.0 cm rather than 280.0 cm as given by Fundamental Methods Associates.)

Results for this problem are shown in Figure 18 and are not the same as those quoted by Fundamental Methods Associates, Inc. Using their original version of HERESY I produces the results shown in Figure 18 for a maximum outer radius of 260.0 cm, but SRL - HERESY I cannot produce Fundamental Methods Associates' HERESY I results for a maximum outer radius of 280.0 cm. This problem arises because the original HERESY I code uses the 61st kernel point (with a maximum kernel size of 60 points) in its Bessel's interpolation of the kernel. Reducing the maximum outer radius of rods considered to 260.0 cm avoids this problem and produces agreement between the two codes.

Input for this problem is shown in Figure 19.

FIGURE 17

```

1 FUNDAMENTAL METHODS ASSOCIATES, INC. TEST PROBLEM II ** SRL-HERESY I

SUMMARY OF HERESY INPUT DATA

INPUT OPTIONS
FINITE/INFINITE OPTION = 1   KERNEL OPTION = 1   COORDINATE OPTION = 1   PURE THERMAL KERNEL OPTION = 0
LATTICE TYPE = -0           GEOMETRY PRINT = 1   GEOMETRY PLOT = 1       GEOMETRY PUNCH = 0

CONVERGENCE CRITERIA
MATRIX INVERSION ITERATIONS = 20   EIGENVALUE ITERATIONS = 200   CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001   EIGENVALUE CONVERGENCE = .000010   CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 60   MESH SIZE = 5.000 CM
B(1) = 1.00   B(2) = 0.00   B(3) = 0.
TAU(1) = 350.00   TAU(2) = 68.30   TAU(3) = 0.
TAUR(1) = 262.50   TAUR(2) = 40.50   TAUR(3) = 0.
B SQUARE = .   L SQUARE = 2800.0000   MODERATOR SIGMA(A) = 0.00040000

REFLECTOR RADIUS = -0.   RHO(MIN) = -0.
RADIUS BEYOND WHICH NO MORE RODS ARE CONSIDERED = 260.000

DATA FOR EACH ROD KIND
ROD KIND   GAMMA   SMF   1./(GAM+SMF)   ETA   RES. COEF.   LSQ   SIGMA(A)   F   CELL VOL.   RADIUS
1          0.25500   0.48216   1.356558   1.34000   51.0000   2800.0   .00040000   0.   0.   2.0000
2          0.17800   0.48216   1.514784   1.60500   51.0000   2800.0   .00040000   0.   0.   2.0000

DATA FOR EACH ROD TYPE
ROD TYPE   ROD KIND   NO. RODS   GAMMA   SMF   1./(GAMMA+SMF)
1          1       127   0.25500   0.48216   1.356558
2          1       127   0.25500   0.48216   1.356558
3          1       127   0.25500   0.48216   1.356558
4          1       127   0.25500   0.48216   1.356558
5          2       127   0.17800   0.48216   1.514784

```

FIGURE 18

NO EIGENVALUE ITERATIONS		EIGENVALUE	PREVIOUS EIGENVALUE				
3		1.1125614	1.1125614				

ROD TYPE		NO RODS/TYPE	ROD SMF	ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB
			GAMMA	ETA	A			
1	127	0.48216	0.255000	1.34000	51.0000	0.962513	0.193800	
2	127	0.48216	0.255000	1.34000	51.0000	0.962513	0.193800	
3	127	0.48216	0.255000	1.34000	51.0000	0.962513	0.193800	
4	127	0.48216	0.255000	1.34000	51.0000	0.962512	0.193800	
5	127	0.48216	0.178000	1.60500	51.0000	1.000000	0.194092	

A THERMAL UTILIZATION VALUE FOR A ROD KIND IS ZERO. MODERATOR THERMAL ABSORPTIONS CANNOT BE SOLVED FOR.
THE MODERATOR ABSORPTIONS WILL BE LUMPED INTO THE TOTAL LEAKAGE

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.11256137 AVERAGE ROD ABSORPTION = 0.9700101

AVERAGE ETA = 1.3946386 THERMAL UTILIZATION = 1.00000000 RESONANCE ESCAPE PROBABILITY = 0.84056976

APPROX MODERATOR THERMAL ABSORPTIONS = 0. TOTAL LEAKAGE = 33.0686 NON-LEAKAGE PROBABILITY = 0.9490488

INFINITE LATTICE MATERIAL BUCKLING = 0.00030475

AVERAGE ABSORPTIONS FOR EACH ROD KIND			
ROD KIND	TOTAL RODS	AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	508	0.962513	0.193800
2	127	1.000000	0.194092

FIGURE 19

1 FUNDAMENTAL METHODS ASSOCIATES, INC. TEST PROBLEM II ** SRL-HERESY I

1	1	1	1	1		
7	60					
60	1	1	0	5	2	
115	1	1	1	1	2	
165					1	
4 20.0	10.0	10.0				
10 5.0	0.0	2800.0	.0004	1.0		
15.000000000001	0.0	350.0	0.0	0.0		
20 262.5	0.0	0.0				
175 .255	.178					
275 1.34	1.605					
325 51.0	51.0					
375 2.0	2.0					
875 2.0	2.0	2.0	2.0	2.0		
925 2.0	2.0	2.0	2.0	2.0		
975 0.0	1.0	0.0	1.0	0.5		
1025 0.0	0.0	1.0	1.0	0.5		
1075 260.0						

1
C

Problem 4. Large Infinite Lattice Test Problem

This problem is the extension of Problem 2 to as large a radius as possible without exceeding a maximum of 6500 rods as specified by the code. The problem contains a total of 6355 rods or 1271 rods of each of five rod types. Figure 13 shows the infinitely repeating pattern of five rods within the dashed lines near the center of the lattice. Figure 20 shows a summary of the input data, and Figure 21 shows the results of the calculations.

Comparing "rod type" 2 of the Finite Lattice Problem 2 and "rod type" 5 of the infinite lattice problem indicates about 0.1% difference between them. This is reasonable agreement for the two problems since the finite lattice considered is large enough to approach being a semi-infinite lattice near its center, and the effects of the infinite moderator are clearly small near the center.

The input cards for this problem are shown in Figure 22.

FIGURE 20

```

1 LARGE TEST PROBLEM FOR INFINITE LATTICE ** SRL-HERESY I

SUMMARY OF HERESY INPUT DATA

INPUT OPTIONS
FINITE/INFINITE OPTION = 1   KERNEL OPTION = 1   COORDINATE OPTION = 1   PURE THERMAL KERNEL OPTION = 0
LATTICE TYPE = -0           GEOMETRY PRINT = 1   GEOMETRY PLOT = 1   GEOMETRY PUNCH = 0

CONVERGENCE CRITERIA
MATRIX INVERSION ITERATIONS = 20   EIGENVALUE ITERATIONS = 200   CRITICALITY SEARCH ITERATIONS = 25
INVERSION CONVERGENCE = .000001   EIGENVALUE CONVERGENCE = .000010   CRITICALITY CONVERGENCE = .000010

KERNEL DATA
NUMBER OF GRID POINTS = 250   MESH SIZE = 4.500 CM
B(1) = 0.56   B(2) = 0.44   B(3) = 0.
TAU(1) = 154.30   TAU(2) = 68.30   TAU(3) = 0.
TAUR(1) = 126.50   TAUR(2) = 40.50   TAUR(3) = 0.
B SQUARE = .   L SQUARE = 12626.7000   MODERATOR SIGMA(A) = 0.00006963

REFLECTOR RADIUS = -0.   RHO(MIN) = -0.
RADIUS BEYOND WHICH NO MORE RODS ARE CONSIDERED = 1025.000

DATA FOR EACH ROD KIND
ROD KIND   GAMMA   SMF   1./(GAM+SMF)   ETA   RES. COEF.   LSQ   SIGMA(A)   F   CELL VOL.   RADIUS
1          0.74668  0.10101  1.179676   1.30530  58.8260  12626.7  .0000696  0.98590  645.16000  0.
2          4.17239  0.10101  0.234006   0.       0.       12626.7  .0000696  0.92365  322.48000  0.

DATA FOR EACH ROD TYPE
ROD TYPE   ROD KIND   NO. RODS   GAMMA   SMF   1./(GAMMA+SMF)
1          1        1271     0.74668  0.10101  1.179676
2          1        1271     0.74668  0.10101  1.179676
3          1        1271     0.74668  0.10101  1.179676
4          1        1271     0.74668  0.10101  1.179676
5          2        1271     4.17239  0.10101  0.234006

```

FIGURE 21

NO EIGENVALUE ITERATIONS 1 EIGENVALUE 1.1232000 PREVIOUS EIGENVALUE 1.1232004

ROD TYPE	NO RODS/TYPE	ROD SMF	ABSORPTIONS FOR EACH ROD TYPE			THERMAL ABSRB	RESONANCE ABSRB
			GAMMA	ETA	A		
1	1271	0.10101	0.746680	1.30530	58.8260	0.999999	0.122910
2	1271	0.10101	0.746680	1.30530	58.8260	1.000000	0.122910
3	1271	0.10101	0.746680	1.30530	58.8260	1.000000	0.122910
4	1271	0.10101	0.746680	1.30530	58.8260	0.999999	0.122910
5	1271	0.10101	4.172394	0.	0.	0.099292	0.

EIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = 1.12320000 AVERAGE ROD ABSORPTION = 0.8198580

AVERAGE ETA = 1.2736834 THERMAL UTILIZATION = 0.98429321 RESONANCE ESCAPE PROBABILITY = 0.89423694

APPROX MODERATOR THERMAL ABSORPTIONS = 83.14140 TOTAL LEAKAGE = -9.9669 NON-LEAKAGE PROBABILITY = 1.0018865

INFINITE LATTICE MATERIAL BUCKLING = 0.00037757

ROD KIND	TOTAL RODS	AVERAGE ABSORPTIONS FOR EACH ROD KIND	
		AVG. THERMAL ABS.	AVG. RESONANCE ABS.
1	5084	1.000000	0.122910
2	1271	0.099292	0.

FIGURE 22

1 LARGE TEST PROBLEM FOR INFINITE LATTICE ** SRL-HERESY I

1	1	1	1	1	1
7	250	1	0	5	2
60	1	1	1	1	2
115	1	1	1	1	2
4	12.7				
10	4.5				
125	645.16	322.48	12626.7	.00006963	
175					
225	.9859	.92365			
275	1.3053	0.0			
325	58.826	0.0			
875	4.0	4.0	4.0	4.0	4.0
925	4.0	4.0	4.0	4.0	4.0
975	1.0	3.0	1.0	3.0	2.0
1025	1.0	1.0	3.0	3.0	2.0
1075	1025.0				

VI. OUTPUT INFORMATION

The SRL HERESY - I code outputs a variety of information both as a data summary and as an edit of the solved problem. In some cases information is printed more than once for convenience. The output format changes according to the particular problem being run so that additional lines appear in a few cases.

A. Equations Used in Edit of Solved Problem

At the end of the sequence of mathematical operations described in Section I-E, the code has solved for the thermal absorption in each rod type (normalized to the largest absorption), and the largest eigenvalue, k , of the system of equations given in equation I.A.1. These quantities are then used, with the input parameters, to obtain a variety of lattice parameters.

The lattice parameters calculated by the edit routine satisfy the relation

$$\bar{\eta} p f \chi = k_{\text{corr}} \quad (\text{VI.A.1})$$

where k_{corr} is the eigenvalue solved for above corrected for fast axial leakage, χ is the lattice nonleakage probability, f is the lattice thermal utilization, p is the lattice resonance escape probability, and $\bar{\eta}$ is the lattice average η value. The equations used to obtain these parameters, as well as other quantities of interest, are given in the following paragraphs.

1. Eigenvalue Corrected for Fast Axial Leakage

The eigenvalue obtained in solving equation I.A.1 takes into account both fast and slow radial leakage, but only slow axial leakage. Fast axial leakage is accounted for by assuming a cosine axial flux shape, and leads to a corrected k given by

$$k_{\text{corr}} = k \exp(-\tau B_z^2) \quad (\text{VI.A.2})$$

where k is the eigenvalue solved for in I.A.1, τ is the thermal age, and B_z^2 is the axial geometrical buckling. The age τ is given by

$$\tau = \sum_{j=1}^3 B_j \tau_j \quad (\text{VI.A.3})$$

where the B_j are weighting fractions and the τ_j are thermal ages for each of the three components of the triple Gaussian slowing down distribution given in Section II-D.

2. Resonance Absorptions

Resonance absorptions occur in rods at the single equivalent resonance level, and are normalized to the highest thermal absorption. They are computed from the equation

$$i_n^R = \frac{A_n}{k_{\text{corr}}} \sum_{m=1}^N g_{nm} i_m \eta_m \quad (\text{VI.A.4})$$

where N is the number of rod types in the lattice, η_m is the thermal η parameter for the m^{th} rod type, i_m is the thermal absorption for the m^{th} rod type, g_{nm} is the number of neutrons born in all rods of type m that reach the resonance energy at a rod type n , A_n is the resonance absorption coefficient (see Section I-D) for rod type n , and k_{corr} is the corrected k from equation VI.A.2.

3. Approximate Moderator Thermal Absorption

The calculation of moderator absorption makes use of the definition of the heterogeneous f parameter given in Section I-B. From this definition, the relationship between moderator absorptions and rod absorptions in a cell with reflecting boundaries is

$$i_{\text{mod}} = i_{\text{rod}} \frac{(1 - f_H)}{f_H} \quad (\text{VI.A.5})$$

where i_{rod} is the total rod thermal absorptions, i_{mod} is the total moderator thermal absorptions, and f_H is the HERESY-defined parameter taken from an appropriate cell calculation.

This same expression may be extended in approximate form to the lattice calculation as

$$M = \sum_{j=1}^K N_j \bar{i}_j \frac{(1 - f_j)}{f_j} \quad (\text{VI.A.6})$$

where K is the number of rod kinds, N_K is the number of rods in the lattice of the k^{th} rod kind, and f_j is the HERESY-defined thermal f parameter for the k^{th} rod kind. M is the moderator thermal absorptions.

4. Total Leakage

The total leakage is defined as the difference between the total fission source and the number of neutrons absorbed (1) at resonance energy, (2) thermally absorbed in rods, and (3) thermally in moderator. This is expressed by

$$L = \frac{1}{k_{\text{corr}}} \sum_{m=1}^N i_m \eta_m (N_m - \sum_{t=1}^N N_t A_t g_{tm}) - \sum_{m=1}^N N_m i_m - M \quad (\text{VI.A.7})$$

where N_m is the number of rods of the m^{th} rod type in the lattice, and all other symbols are defined as previously.

5. Nonleakage Probability

This quantity is defined to be the total thermal absorptions divided by total thermal neutron source.

$$\chi = \frac{\sum_{m=1}^N N_m i_m + M}{\sum_{m=1}^N N_m i_m + M + L} \quad (\text{VI.A.8})$$

where all symbols have been defined previously.

6. Thermal Utilization

The definition of this quantity for the lattice is much the same as for the individual rods.

$$f_{\text{lattice}} = \frac{\text{Total thermal absorptions in rods}}{\text{Total thermal absorptions in rods plus moderator}}$$

or

$$f_{\text{lattice}} = \frac{\sum_{m=1}^N N_m i_m}{\sum_{m=1}^N N_m i_m + M} \quad (\text{VI.A.9})$$

where all symbols have been defined previously.

7. Resonance Escape Probability

A standard definition is used for computations as

$$p_{\text{lattice}} = \frac{\sum_{m=1}^N N_m i_m + M + L}{\sum_{m=1}^N N_m i_m + M + L + \sum_{m=1}^N N_m i_m^R} \quad (\text{VI.A.10})$$

where all symbols have been defined previously.

8. Average Rod Absorption

This is obtained by a simple arithmetic average over the rods in the lattice.

$$\bar{i} = \frac{\sum_{m=1}^N N_m i_m}{\sum_{m=1}^N N_m} \quad (\text{VI.A.11})$$

where all symbols have been previously defined.

9. Average Thermal Eta

A simple arithmetic average weighted by thermal absorptions is used as

$$\bar{\eta} = \frac{\sum_{m=1}^N N_m \eta_m i_m}{\sum_{m=1}^N N_m i_m} \quad (\text{VI.A.12})$$

where all symbols have been defined previously.

10. Average Absorptions for Each Rod Kind

A simple arithmetic average is made as

$$\bar{i}_p = \frac{\text{Absorptions in all rods of kind } p}{\text{Total number of rods of kind } p} \quad (\text{VI.A.13})$$

11. Infinite Lattice Material Buckling

If the lattice considered is radially infinite and $B_z^2 = 0$, then the eigenvalue found from solving equation I.A.1 is k_∞ . For this case, the material buckling of the lattice may be solved for under the assumption that

$$L_{\text{lattice}}^2 = (1 - f_{\text{lattice}}) L_{\text{moderator}}^2 \quad (\text{VI.A.14})$$

The material buckling is obtained by successive iterations on the age-diffusion critical equation in the form

$$k_\infty = [1 + L_{\text{mod}}^2 (1 - f) B_m^2] \exp(\tau B_m^2) \quad (\text{VI.A.15})$$

where all symbols have been defined previously.

B. Output Formats

Several examples of output formats for both finite and infinite problems were shown in Section V. The quantities are output in a particular order subject to modification by selection of the many options. The order of output for a routine problem is given, then possible modifications that can occur are itemized.

1. Routine Problem Format

A standard problem will produce the following format:

- A line summarizing the number of iterations necessary to converge to the final eigenvalue, and the last two eigenvalues used in the iteration. A comparison of the two eigenvalues assures that the problem did indeed converge.
- A table summarizing the rod parameters, thermal absorptions, and resonance absorptions for each rod type.
- A list of the edited lattice parameters. This includes k_{corr} , \bar{i} , $\bar{\eta}$, f , p , M , L , and χ .
- A table summarizing the average thermal and resonance absorptions for each rod kind.

2. Modifications to Routine Output Format

- If an edit of each iteration of a criticality search is requested, a line is printed prior to the line of the routine format (first item) giving the number of the current iteration (Figure 5).
- If the thermal utilization is zero for a rod kind, it is not possible to calculate the moderator absorptions as given in Section VI-A-3. Under these conditions, the value of M is set to zero and a line is printed under second item of the routine format to so indicate this. The effect of setting $M = 0$ is to force f to equal 1.0 and lump the moderator absorptions into the total leakage, however equation VI.A.1 is still satisfied.

- If the problem concerns an infinite lattice and B_z^2 is zero, the lattice material buckling is computed as described in Section VI-A-11 and inserted following third item of the routine format.
- If a criticality search on gamma of a rod kind has been performed, a line will be inserted following fourth item of the routine format giving the thermal utilization corresponding to this gamma value if it is possible to re-compute f. If insufficient information is available to re-compute f, a line will be inserted to so indicate.

VII. REFERENCES

A. Cited

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B. General

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- (2) Ibid. NYO-2675 (1959). (TRG-129-QTR-3)
- (3) Ibid. NYO-2676 (1959). (TRG-129-QTR-4)
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- (8) W. A. Horning. Small Source Model of a Thermal Pile. USAEC Report HW-24282, General Electric, Hanford Works, Richland, Wash. (1952)(declassified 1957).

APPENDIX A - Mnemonics Used in Code

The list below contains all of the primary symbols used in SRL - HERESY I, their dimensions, and their locations in the three large storage arrays. The code uses three arrays in COMMON STORAGE to store data, designated as the LL(225) or fixed point array, AA(15,500) or Link 1 floating point array, and the AAA(10,250) or Link 2 floating point array. Main Program 1 uses only the LL(225) and AA(15,500) arrays while Main Program 2 uses the LL(225), AA(15,500), and AAA(10,250) arrays. Shared storage is made up of arrays LL(225) and AA(15,500), the AAA(10,250) being the necessary amount of temporary storage to handle the 50 x 50 arrays used in Link 2.

The following list contains the majority of the mnemonics used. Those used internally, hence not appearing in the input, are indicated by an asterisk.

<u>Mnemonic</u>	<u>Location</u>	<u>Use</u>
* A(50)	AA(2342)...	Resonance absorption parameters for each rod type
* A1	AA(1056)	Maximum value of V1 matrix
* A2	AA(1077)	Maximum value of V2 matrix
AR(48)	AA(325)...	Resonance absorption parameters for each rod kind
* ABSRB(50)	AAA(10001)...	Normalized thermal absorptions for each rod type
B(3)	AA(14)...	Weighting fractions for each term of triple Gaussian slowing down distribution
* BGF(250)	AA(1142)...	Fission to thermal kernel values
* BGFR(250)	AA(1392)...	Resonance to thermal kernel values
BSQ	AA(11)	Axial buckling
CRI	AA(1)	Convergence criterion on matrix inversion

<u>Mnemonic</u>	<u>Location</u>	<u>Use</u>
CR2	AA(2)	Convergence criterion on power iteration
CR3	AA(3)	Convergence criterion on criticality search
* D(48)	AA(2192)...	Moderator diffusion coefficient for each rod kind
EIGEN	AA(9)	Eigenvalue to be searched for
* ETA(50)	AA(2292)	Eta values for each rod type
ETAI(48)	AA(275)	Eta values for each rod kind
* FK1	AA(1078)	Value of <u>uncorrected</u> eigenvalue for the current iteration minus one
* FK2	AA(1069)	Value of <u>uncorrected</u> eigenvalue for current iteration
* FKMAX(50)	AA(2442)...	No. rods/type in floating point
GAM(I)	AA(175)...	Gamma values for each rod kind
* GAMMA(I)	AA(2242)...	Gamma values for each rod type
* IMAX1	LL(174)	Number of largest element of V1 array
* IMAX2	LL(175)	Number of largest element of V2 array
INF	LL(1)	Finite/infinite lattice indicator
* IT1	LL(170)	Iteration counter for matrix inversion
* IT2	LL(171)	Iteration counter for power eigenvalue iteration
* IT3	LL(172)	Print counter in power eigenvalue iteration
* IT4	LL(173)	Iteration counter for criticality search

<u>Mnemonic</u>	<u>Location</u>	<u>Use</u>
KDAT	LL(168)	Omit data print indicator
KERNEL	LL(2)	New kernel indicator
KIND(50)	LL(115)...	Rod kind for each rod type
KOOR	LL(3)	New coordinate indicator
KPLOT	LL(61)	Geometry plot indicator
KPRNT	LL(60)	Geometry list indicator
KSEPR	LL(167)	Separate edit indicator (criticality searches)
KSERCH	LL(166)	Rod kind to be searched over
KSKIP	LL(62)	Geometry punch indicator
KTEST	LL(169)	Last problem indicator
LAT	LL(9)	Lattice type indicator
LCRIT	LL(165)	Criticality search indicator
N1	LL(94)	Max. No. iterations on matrix inversion
N2	LL(95)	Max. No. iterations on eigenvalue
N3	LL(96)	Max. No. iterations in criticality search
* NERR	AA(24)	Error indicator
* NF(50)	LL(9176)...	Indicator for f calculation for each rod kind
NICPRT(50)	LL(10)...	No. initial coordinates for each rod type
NKIND	LL(64)	Number of rod kinds
NOGRID	LL(7)	Number of kernel mesh points

<u>Mnemonic</u>	<u>Location</u>	<u>Use</u>
NTYPE	LL(63)	Number of rod types
NWRITE	LL(224)	Write indicator for library tapes
RADII	AA(1075)	Radius beyond which no more rods are considered in infinite lattices
RADIUS(48)	AA(375)...	Radius of each rod kind
RADRFL	AA(7)	Reflector radius
* RESABS(50)	AAA(10051)...	Resonance absorption for each rod type
ROEMIN	AA(8)	A radius within which the reflector may be considered infinite
SIGA	AA(13)	Moderator absorption cross section
SIGC(48)	AA(75)...	Absorption cross section for gamma calculations for each rod kind
SM(48)	AA(2142)...	Self-contribution to thermal diffusion kernel for each rod kind
* SMF(250)	AA(1892)...	Thermal diffusion kernel points
* SMFO(50)	AA(2392)...	Self-contribution to thermal diffusion kernel for each rod type
* SMGR(250)	AA(1642)...	Resonance to thermal kernel values
* SR(50,50)	AA(8500)...	Resonance to thermal slowing down matrix
* T(50,50)	AA(3500)...	Fission to thermal slowing down matrix
* T1(50,50)	AAA(1)...	Temporary matrix for computations
* T2(50,50)	AAA(2501)...	Temporary matrix for computations
* T3(50,50)	AAA(5001)...	Temporary matrix for computations
* T4(50,50)	AAA(7501)...	Temporary matrix for computations

Mnemonic	Location	Use
* T5(50)	AA(13,500)...	Temporary matrix for computations
* T6(50)	AA(13,550)...	Temporary matrix for computations
TAU(3)	AA(17)...	Fission to thermal age
TAUR(3)	AA(20)...	Fission to resonance age
* TAUTOL	AA(23)	$= \text{EXPF} \{ -\text{BSQ} * [\text{B}(1) * \text{TAU}(1) + \text{B}(2) * \text{TAU}(2) + \text{B}(3) * \text{TAU}(3)] \}$
THERU(50)	AA(225)...	Thermal utilization for each rod kind
THETA(50,3)	AA(725)...	Angular initial coordinates
TITLE(12)	AA(1130)...	Problem alphanumeric data
* TR(50,50)	AA(6000)...	Fission to resonance slowing down matrix
* V1(50)	AAA(10151)...	Previous set of unnormalized eigenvectors
* V2(50)	AAA(10201)...	Current set of unnormalized eigenvectors
VCELL(48)	AA(125)...	Cell volume for each rod kind
VO(50)	AA(1080)...	Initial approximation to eigenvectors
* VRATIO(50)	AAA(10101)...	Ratio of V2/A2 for each rod type
* W(8)	AA(2492)...	Miscellaneous transfer locations
WDX(50)	AA(875)...	X direction pitch in infinite lattice
WDY(50)	AA(925)...	Y direction pitch in infinite lattice
WX1(50)	AA(975)...	Initial X position in infinite lattice

<u>Mnemonic</u>	<u>Location</u>	<u>Use</u>
WY1(50)	AA(1025)...	Initial Y position in infinite lattice
X(6500)	AA(2500)	X coordinates of all rods
XCOR	AA(5)	Scale factor on all coordinates
XLSQ	AA(12)	L^2 for moderator
XLSQ2(50)	AA(25)...	L^2 of moderator for each rod kind - used in self-consistent calculation of Y
XMESH	AA(10)	Mesh size for kernel
X0(50,3)	AA(425)...	Initial X coordinates
XSCALE	AA(5)	Scale increment along X axis in plotter
Y(6500)	AA(9000)...	Y coordinates of all rods
Y0(50,3)	AA(575)...	Initial Y coordinates
YSCALE	AA(6)	Scale increment along Y axis in plotter
* Z(50,50)	AA(11,000)...	Thermal diffusion kernel matrix

The three Arrays LL, AA, and AAA are stored in order in common storage, i.e. LL(1) = location 77462₈ in storage. All mnemonics are then assigned to locations in these three arrays by EQUIVALENCE statements. Information concerning the three large arrays follow.

<u>Array</u>	<u>First Position in Storage</u>	<u>Last Position in Storage</u>
LL(225 ₁₀)	32,562 ₁₀ or 77462 ₈	32,338 ₁₀ or 77122 ₈
AA(15,500 ₁₀)	32,337 ₁₀ or 77121 ₈	16,838 ₁₀ or 40706 ₈
AAA(10,250 ₁₀)	16,837 ₁₀ or 40705 ₈	6,588 ₁₀ or 14,674 ₈

APPENDIX B - FORTRAN II Listings

The source decks for all FORTRAN II and IBM 704 SAP subroutines used in SRL - HERESY I are listed on the following pages. Two sections appear which correspond to the two links in the CHAIN code. The CHAIN subroutine and related codes (i.e. CPTG, EPTC, and SLCI routines) are special versions not available from another source. Appendix D gives information on the CHAIN codes, as well as listings of their source cards.

Link 1

Link 1 consists of 13 FORTRAN II subprograms, a package of library function subroutines, and a package of FORTRAN II control subroutines. COMMON STORAGE begins at location 407058 in this link.

FORTRAN II Subprograms

<u>Subprogram Name</u>	<u>Function</u>
MAIN PROG. 1	Main Prog 1 executive routine
DATIN	Data reading subroutine
DATCHK	Data checking subroutine
KERGEN	Kernel generator subroutine
GEOM	Finite geometry subroutine
INGEOM	Infinite geometry subroutine
PLOT	Geometry plotting subroutine
PRINT	Geometry listing subroutine
DATPRT	Data summary output subroutine
HEX	Hexagonal geometry subroutine
SQARE	Square geometry subroutine
CIRCLE	Circular pattern geometry subroutine
PUNCH	Geometry punch subroutine

FORTRAN II Library Function Subroutines

The following functions must appear in the program with the call names as listed:

<u>FORTRAN Name</u>	<u>Function</u>
BESSKK (arg)	K_0 (arg)
ATANF (arg)	\tan^{-1} (arg)
LOGF (arg)	\ln (arg)
EXPF (arg)	e (arg)
EXPIF (arg)	E_1 (arg) integral
SINF (arg)	\sin (arg)
COSF (arg)	\cos (arg)
SQRFT (arg)	$\sqrt{\text{arg}}$

FORTRAN II Control Subroutines

<u>FORTRAN Name</u>	<u>Operation</u>
(FIL)	Format interpretive loader
(BDC)	Binary to decimal conversion
(DBC)	Decimal to binary conversion
(IOH)O	Input - Output Hollerith, Output
(IOH)I	Input - Output Hollerith, Input
(SPH)	Storage to printer Hollerith
(STH)	Storage to tape Hollerith
(TSH)	Tape to storage Hollerith
(SCH)	Storage to card Hollerith
(RTN)	Return
(LEV)	Leave
(LFM)	Leave floating point trap
(EFM)	Enter floating point trap
CHAIN	Link search routine
CPTG	Program tape link generator

C
C
C
C

DIMENSION DECK FOR SRL HERESY 1 MAIN PROGRAM 1

DIMENSION DECK FOR MAIN PROGRAM 1 VERSION OF 10/2/64

```

DIMENSION LL(225),AA(15500)
DIMENSION NICPRT(50),NRPRT(50),KIND(50),NF(50)
DIMENSION B(3),TAU(3),TAUR(3),XLSQ2(50),SIGC(50),VCELL(50),GAM(5
1  ,THERU(50),ETAI(50),AL(50),RADIUS(50),XO(50,3),YO(50,3),
2  THETA(50,3),WDX(50),WDY(50),WX1(50),WY1(50),VO(50),W(8),
3  TITLE(12),BGF(250),BGFR(250),SMGR(250),SMF(250),SM(50),D(50),
4  GAMMA(50),ETA(50),A(50),SMFO(50),FKMAX(50),X(6500),Y(6500)
COMMON LL,AA

```

C
C
C

EQUIVALENCES FOR LL ARRAY

```

EQUIVALENCE (INF,LL(1)),(KERNEL,LL(2)),(KOOR,LL(3)),(N1,LL(4)),
1  (N2,LL(5)),(N3,LL(6)),(NOGRID,LL(7)),(NHER,LL(8)),(LAT,
2  LL(9)),(NICPRT,LL(10)),(KPRNT,LL(60)),(KPLT,LL(61)),(KSKIP,
3  LL(62)),(NTYPE,LL(63)),(NKIND,LL(64)),(NRPRT,LL(65)),(KIND,
4  LL(115)),(LCRIT,LL(165)),(KSERCH,LL(166)),(NSEPR,LL(167)),
5  XKDAT,LL(168)),(KTEST,LL(169)),(IT1,LL(170)),(IT2,LL(171)),
6  (IT3,LL(172)),(IT4,LL(173)),(IMAX1,LL(174)),(IMAX2,LL(175)),
7  (NF,LL(176))

```

C
C
C

EQUIVALENCES FOR AA ARRAY

```

EQUIVALENCE (CR1,AA(1)),(CR2,AA(2)),(CR3,AA(3)),(XCOR,AA(4)),
1  (XSCALE,AA(5)),(YSCALE,AA(6)),(RADRFL,AA(7)),(ROEMIN,AA(8)),
2  (EIGEN,AA(9)),(XMESH,AA(10)),(BSQ,AA(11)),(XLSQ,AA(12)),
3  (SIGA,AA(13)),(B,AA(14)),(TAU,AA(17)),(TAUR,AA(20)),(TAUTOL,
4  AA(23)),(NERR,AA(24)),(XLSQ2,AA(25)),(SIGC,AA(75)),(VCELL,
5  AA(125)),(GAM,AA(175)),(THERU,AA(225)),(ETAI,AA(275)),(AL,
6  AA(325)),(RADIUS,AA(375)),(XO,AA(425)),(YO,AA(575)),(THETA,
7  AA( 725)),(WDX,AA( 875)),(WDY,AA( 925)),(WX1,AA( 975)),(WY1,
8  AA(1025)),(RADII,AA(1075)),(A1,AA(1076)),(A2,AA(1077)),(FK1,
9  AA(1078)),(FK2,AA(1079)),(VO,AA(1080)),(W,AA(2492))
EQUIVALENCE (BGF,AA(1142)),(BGFR,AA(1392)),(SMGR,AA(1642)),
1  (SMF,AA(1892)),(SM,AA(2142)),(D,AA(2192)),(GAMMA,AA(2242)),
2  (ETA,AA(2292)),(A,AA(2342)),(SMFO,AA(2392)),(FKMAX,AA(2442))
3  (X,AA(2500)),(Y,AA(9000)),(TITLE,AA(1130))

```

G

SRL HERESY I MAIN PROGRAM 1

VERSION OF 10/12/64

HER-P

SUBROUTINE MAIN

HER-P
DIME

INSERT DIMENSION DECK FOR MAIN PROGRAM 1

```

CALL EFM
IN=8
NOUT=10
5 READ INPUT TAPE IN,10,(TITLE(I),I=1,12)
10 FORMAT(12A6)
NERR=0
PRINT 15,(TITLE(I),I=1,12)
15 FORMAT(1H0,12A6)
WRITE OUTPUT TAPE NOUT,10,(TITLE(I),I=1,12)
CALL DATIN
CALL DATCHK
IF(NERR)40,40,20
20 PRINT 30
30 FORMAT(1H0,61HAN ERROR HAS BEEN DETECTED IN THE INPUT DATA FOR THIER-P
1S PROBLEM/1H ,60HAN ATTEMPT WILL BE MADE TO CONTINUE WITH SUCCESSIHER-P
2VE PROBLEMS)
GO TO 5
40 IF(KERNEL)60,60,50
50 CALL KERGEIN
IF(NERR)60,60,20
60 IF(KOOR-1)100,70,100
70 IF(INF)90,80,90
80 CALL GEOM
IF(NERR)100,100,20
90 CALL INGEOM
IF(NERR)100,100,20
100 IF(KPLOT)105,105,101
101 CALL PLOT
105 IF(KPRNT)109,109,106
106 CALL PRINT
109 DO 110 I=1,NTYPE
110 FKMAX(I)=NRPRT(I)
TAUTOL=EXP(-BSQ*(B(1)*TAU(1)+B(2)*TAU(2)+B(3)*TAU(3)))
DO 160 I=1,NKIND
D(I)=XLSQ2(I)*SIGC(I)
XYLSQ=XLSQ2(I)/(1.+BSQ*XLSQ2(I))
XXLSQ=SQRT(XYLSQ)
IF(RADIUS(I))120,120,130
120 SM(I)=0.101010
GO TO 140
130 ARG1=RADIUS(I)/XXLSQ
ARG2=0.0
ARG3=0.0
ARG4=0.0
ARG5=0.0
ARG6=0.0
ARG7=0.0
CALL BESSI(ARG1,ARG2,ARG3,ARG4,ARG5)
CALL BESSKN(ARG1,ARG2,ARG3,ARG4,ARG5,ARG6,ARG7)
SM(I)=(1./(6.2831852*D(I)))*ARG4
140 IF(GAM(I))160,150,160
150 GAM(I)=XLSQ2(I)*(1.-THERU(I))/(D(I)*VCELL(I)*THERU(I))+
1(1./(12.56637 *D(I)))*LOGF( XYLSQ/VCELL(I))*228714/D(I)-SM(I)

```

NF(I)=1	HER-2
160 CONTINUE	HER-2
DO 170 I=1,NTYPE	HER-2
NX=KIND(I)	HER-2
ETA(I)=ETA1(NX)	HER-2
GAMMA(I)=GAM(NX)	HER-2
A(I)=AL(NX)	HER-2
170 SMFO(I)=SM(NX)	HER-2
IF(KOOR-2)210,180,210	
180 NX=0	HER-2
DO 190 I=1,NTYPE	HER-2
190 NX=NX+NRPRT(I)	HER-2
DO 200 I=1,NX	HER-2
X(I)=X(I)*XCOR	HER-2
200 Y(I)=Y(I)*XCOR	HER-2
210 IF(KDAT)220,220,230	HER-2
220 CALL DATPRT	HER-2
230 CALL CHAIN(2,1)	HER-2
END(2,1,0,0,0)	HER-2

C	SRL HERESY 1 SUBROUTINE DATIN	VERSION OF 10/2/64	DA
C			DA
	SUBROUTINE DATIN		DA
C	INSERT DIMENSION DECK FOR MAIN PROGRAM 1		
	CALL EFM		DA
	IN=8		DA
	10 READ INPUT TAPE IN,20,LAST,NUMD,LOCATE,(KK(I),I=1,5)		DA
	20 FORMAT(I1,3X11,2X15,5I12)		DA
	IF(NUMD)22,22,24		DA
	22 NUMD=5		DA
	24 DO 30 I=1,NUMD		DA
	M=LOCATE+I-1		DA
	30 LL(M)=KK(I)		DA
	IF(LAST)10,10,40		DA
	40 READ INPUT TAPE IN,50,LAST,NUMD,LOCATE,(AX(I),I=1,5)		DA
	50 FORMAT(I1,3X11,2X15,5F12.)		DA
	IF(NUMD)52,52,54		DA
	52 NUMD=5		DA
	54 DO 60 I=1,NUMD		DA
	M=LOCATE+I-1		DA
	60 AA(M)=AX(I)		DA
	IF(LAST)40,40,70		DA
	70 RETURN		DA
	END(2,1,0,0,0)		DA

C	SRL HERESY 1 SUBROUTINE DATCHK VERSION OF 10/2/64	DAT
C		DAT
	SUBROUTINE DATCHK	DAT
C		DIME
	INSERT DIMENSION DECK FOR MAIN PROGRAM 1	
C		DIME
	CALL EFM	DAT
	NOUT=10	DAT
	40 IF(NKIND-NTYPE)70,70,50	DAT
	50 WRITE OUTPUT TAPE NOUT,60	DAT
	60 FORMAT(1H0,63HTHE NUMBER OF ROD KINDS IS GREATER THAN THE NUMBER 0	DAT
	IF ROD TYPES)	DAT
	NERR=1	DAT
	70 IF(NTYPE-50)100,100,80	DAT
	80 NERR=1	DAT
	WRITE OUTPUT TAPE NOUT,90	DAT
	90 FORMAT(1H0,34HTHE NUMBER OF ROD TYPES EXCEEDS 50)	DAT
	100 IF(KOOR-1)105,128,105	DAT
	105 DO 125 I=1,NTYPE	DAT
	IF(NRPRT(I))110,110,125	DAT
	110 WRITE OUTPUT TAPE NOUT,120	DAT
	120 FORMAT(1H0,26H A ROD TYPE HAS ZERO RODS)	DAT
	NERR=1	DAT
	125 CONTINUE	DAT
	128 DO 160 I=1,NTYPE	DAT
	130 IF(KIND(I))140,140,160	DAT
	140 WRITE OUTPUT TAPE NOUT,150	DAT
	150 FORMAT(1H0,38HA ROD TYPE HAS NO KIND NUMBER ASSIGNED)	DAT
	NERR=1	DAT
	160 CONTINUE	DAT
	IF(N1)170,170,180	DAT
	170 N1=20	DAT
	180 IF(N2)190,190,200	DAT
	190 N2=200	DAT
	200 IF(N3)203,203,203	DAT
	203 N3=25	DAT
	208 IF(KOOR-1)240,210,240	DAT
	210 IF(LAT-2)240,240,220	DAT
	220 WRITE OUTPUT TAPE NOUT,230	DAT
	NERR=1	DAT
	230 FORMAT(1H0,60HTHE LATTICE TYPE TO BE GENERATED IS NOT OF A RECOGNI	DAT
	IZED TYPE)	DAT
	240 IF(LCRIT)230,280,250	DAT
	250 IF(LCRIT-2)255,255,280	DAT
	255 IF(KSERCH)260,260,280	DAT
	260 NERR=1	DAT
	WRITE OUTPUT TAPE NOUT,270	DAT
	270 FORMAT(1H0,54HA ROD KIND IS NOT SPECIFIED FOR THE CRITICALITY SEAR	DAT
	1CH)	DAT
	280 IF(CR1)290,290,300	DAT
	290 CR1=.000001	DAT
	300 IF(CR2)310,310,320	DAT
	310 CR2=.00001	DAT
	320 IF(CR3)330,330,340	DAT
	330 CR3=.00001	DAT
	340 IF(B(1))350,350,360	DAT
	350 B(1)=0.56	DAT
	360 IF(B(2))370,370,380	DAT
	370 B(2)=0.44	DAT

380 IF(B(3))390,390,400	DAT
390 B(3)=0.0	DAT
400 IF(TAU(1))410,410,420	DAT
410 TAU(1)=154.3	DAT
420 IF(TAU(2))430,430,440	DAT
430 TAU(2)=68.3	DAT
440 IF(TAU(3))450,450,460	DAT
450 TAU(3)=0.0	DAT
460 IF(TAUR(1))470,470,480	DAT
470 TAUR(1)=126.5	DAT
480 IF(TAUR(2))490,490,500	DAT
490 TAUR(2)=40.5	DAT
500 IF(TAUR(3))510,510,520	DAT
510 TAUR(3)=0.0	DAT
520 DO 560 I=1,NKIND	DAT
IF(XLSQ2(I))530,530,540	DAT
530 XLSQ2(I)=XLSQ	DAT
540 IF(SIGC(I))550,550,560	DAT
550 SIGC(I)=SIGA	DAT
560 CONTINUE	DAT
IF(XCOR)558,555,558	DAT
555 XCOR=1.0	DAT
558 IF(KOOR-1)610,570,610	DAT
570 IF(XSCALE)580,580,590	DAT
580 XSCALE=XCOR	DAT
590 IF(YSCALE)600,600,610	DAT
600 YSCALE=XCOR	DAT
610 IF(LCRIT)630,630,620	DAT
620 IF(EIGEN)625,625,630	DAT
625 EIGEN=1.0	DAT
630 DO 650 I=1,NTYPE	DAT
IF(VO(I))640,640,650	DAT
640 VO(I)=1.0	DAT
650 CONTINUE	DAT
IF(KERNEL-1)720,675,675	DAT
675 IF(XMESH)700,700,680	DAT
680 IF(XLSQ)700,700,690	DAT
690 IF(SIGA)700,700,720	DAT
700 NERR=1	DAT
WRITE OUTPUT TAPE NOUT,710	DAT
710 FORMAT(1H0,62HTHE INFORMATION GIVEN FOR THE KERNEL GENERATOR IS IN	DAT
1SUFFICIENT)	DAT
720 DO 780 I=1,NKIND	DAT
IF(GAM(I))730,730,780	DAT
730 IF(ITERU(I))760,760,740	DAT
740 IF(VCELL(I))760,760,750	DAT
750 IF(SIGC(I))760,760,780	DAT
760 NERR=1	DAT
WRITE OUTPUT TAPE NOUT,770,1	DAT
770 FORMAT(1H0,35HTHE INFORMATION GIVEN FOR ROD KIND 12,35H IS INSUFFI	DAT
1CIENT TO COMPUTE A GAMMA)	DAT
780 CONTINUE	DAT
IF(NERR)810,810,790	DAT
790 WRITE OUTPUT TAPE NOUT,800	DAT
800 FORMAT(1H0,71HBECAUSE OF THE ERROR(S) DETECTED ABOVE THIS PROBLEM	DAT
1CANNOT BE CONTINUED/1H-88HSUCCESSIVE PROBLEMS WILL BE ATTEMPTED,	DAT
2ALTHOUGH THEY MAY BE AFFECTED BY THE ERRORS ABOVE)	DAT
810 RETURN	DAT
END(2,1,0,0,0)	DAT

```

C          SRL HERESY 1  SUBROUTINE KERGEN 1          VERSION OF 10/2/64
C
C          SUBROUTINE KERGEN
C
C          INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EPM
      NOUT=10
      I1=0
      I2=0
      DENL=1.+BSQ*XLSQ
      XXLSQ=XLSQ/DENL
      SIGMA=SIGA*DENL
      CC=1./(6.2831852*XXLSQ*SIGMA)
      RLSQ=SQRTF(XXLSQ)
      IF(ABSF(B(2)))340,20,10
10  IF(ABSF(B(3)))340,30,40
20  NN=1
      GO TO 50
30  NN=2
      GO TO 50
40  NN=3
50  DO 380 I=1,NOGRID
      FI=I-1
      DX=FI*XMESH
      ARG1=DX/RLSQ
      IF(ARG1)340,60,70
60  IFPATH=3
      GO TO 80
70  IFPATH=1
      ARG2=0.0
      ARG3=0.0
      ARG4=0.0
      ARG5=0.0
      ARG6=0.0
      ARG7=0.0
      CALL BESSI(ARG1,ARG2,ARG3,ARG4,ARG5)
      CALL BESSK(ARG1,ARG2,ARG3,ARG4,ARG5,ARG6,ARG7)
80  SMGR(I)=0.0
      D2BY4=DX**2/4.
      DO 90 J=1,NN
90  SMGR(I)=SMGR(I)+(B(J)/(TAUR(J)*12.5663705))*EXP(-D2BY4/TAUR(J))
      DO 100 J=1,NN
100 WTAU(J)=TAU(J)
105 WBG=0.0
      GO TO (210,210,110,110),IFPATH
110 DO 180 J=1,NN
      IF(ABSF(WTAU(J)/XXLSQ)-87.3)140,140,120
120 IF(I1)125,125,135
125 WRITE OUTPUT TAPE NOUT,130
130 FORMAT(1H0,68HTHE ARGUMENT OF THE E1 INTEGRAL IN THE KERNEL GENERA
      1TOR EXCEEDS 87.3/1H ,85HTHE VALUE OF THE INTEGRAL WILL BE SET EQUA
      2L TO 1.0E-30 AND COMPUTATIONS WILL CONTINUE)
      I1=1
135 CON(J)=1.0E-30
      GO TO 180
140 IF(ABSF(WTAU(J)/XXLSQ)-.00001)150,150,170
150 IF(I2)155,155,165
155 WRITE OUTPUT TAPE NOUT,160

```

160 FORMAT(1H0,81H THE ARGUMENT OF THE E1 INTEGRAL IN THE KERNEL GENERA
1TOR IS WITHIN 1.0E-05 OF ZERO/1H ,85H THE VALUE OF THE INTEGRAL WILL
2L BE SET EQUAL TO 1.0E+30 AND COMPUTATIONS WILL CONTINUE)

```

      I2=1
165 CON(J)=1.0E+30
      GO TO 180
170 CON(J)=EXP(-WTAU(J)/XXLSQ)
180 CONTINUE
      DO 190 J=1,NN
190 WBG=WBG-(B(J)/2.)*CC*EXP(WTAU(J)/XXLSQ)*CON(J)
      GO TO (340,340,200,360),IFPATH
200 BGF(I)=WBG
      IFPATH=4
      GO TO 290
210 DO 260 J=1,NN
      IF(ABS(D2BY4/WTAU(J))-87.3)230,230,220
220 IF(I1)225,225,228
225 WRITE OUTPUT TAPE NOUT,130
      I1=1
228 CON(J)=1.0E-30
      GO TO 260
230 IF(ABS(D2BY4/WTAU(J))-0.00001)240,240,250
240 IF(I2)245,245,248
245 WRITE OUTPUT TAPE NOUT,160
      I2=1
248 CON(J)=1.0E+30
      GO TO 260
250 CON(J)=EXP(-D2BY4/WTAU(J))
260 CONTINUE
      DO 270 J=1,NN
270 WBG=WBG+CC*EXP(WTAU(J)/XXLSQ)*B(J)*(ARG4+(WTAU(J)/2./XXLSQ)*
1EXP(-D2BY4/WTAU(J))+(.5+D2BY4/(2.*XXLSQ))*CON(J))
      GO TO (280,360,340,340),IFPATH
280 BGF(I)=WBG
      IFPATH=2
290 DO 300 J=1,NN
300 WTAU(J)=TAU(J)-TAUR(J)
      DO 330 J=1,NN
      IF(WTAU(J))310,310,330
310 WRITE OUTPUT TAPE NOUT,320
320 FORMAT(1H0,69H A TAU(R) VALUE IS LARGER THAN OR EQUAL TO ITS CORRES  

1PONDING TAU VALUE/1H ,51H A KERNEL CANNOT BE GENERATED UNDER THESE  

2CONDITIONS)
      NERR=1
      RETURN
330 CONTINUE
      GO TO (340,105,340,105),IFPATH
340 WRITE OUTPUT TAPE NOUT,350
350 FORMAT(1H0,70H A MACHINE ERROR HAS OCCURRED IN THE KERNEL GENERATOR  

1 - RESTART PROBLEM)
      PRINT 350
      NERR=1
      RETURN
360 BGR(I)=WBG
      SMF(I)=CC*ARG4
      IF(NTHER)380,380,370
370 BGE(I)=SMF(I)
      BGR(I)=SMF(I)
      SMGR(I)=1.
380 CONTINUE
430 RETURN
      END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE  GEOM      VERSION OF 10/14/64
C
C      SUBROUTINE GEOM
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL BPM
      NOUT=10
      IF(LAT)10,10,20
10  CALL HEX
      GO TO 50
20  IF(LAT-1)30,30,40
30  CALL SQUARE
      GO TO 50
40  CALL CIRCLE
50  IF(KSKIP)90,90,80
80  CALL PUNCH
90  NX=0
      DO 120 I=1,NTYPE
120  NX=NX+NRPT(I)
      DO 130 I=1,NX
          X(I)=X(I)*XCOR
130  Y(I)=Y(I)*XCOR
          IF(NX-6500)160,160,140
140  WRITE OUTPUT TAPE NOUT,150
150  FORMAT(1H0,37HTHE TOTAL NUMBER OF RODS EXCEEDS 6500)
          NERR=1
160  RETURN
          END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE HEX      VERSION OF 10/2/64
C
C      SUBROUTINE HEX
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NOUT=10
      K=0
      DO 160 I=1, NTYPE
      L=1
      J=0
      5 IF(YO(I,L))10,40,10
      10 IF(XO(I,L))60,20,60
      20 NERR=1
      WRITE OUTPUT TAPE NOUT,30
      30 FORMAT(1H0,69HERROR IN GEOMETRY GENERATOR - INITIAL COORDINATES AR
      IE IN WRONG SECTOR)
      RETURN
      40 ANGLE=0.0
      IF(XO(L,L))70,50,70
      50 NN=1
      M=K+NN
      X(M)=0.0
      Y(M)=0.0
      GO TO 150
      60 TAN=YO(I,L)/XO(I,L)
      ANGLE=ATANF(TAN)
      70 RADIX=SQRTF(XO(I,L)*XO(I,L)+YO(I,L)*YO(I,L))
      DO 120 N=1,6
      XN=FLOATF(N)
      NN=J+N
      M=J+N+K
      X(M)=RADIX*COSF(ANGLE+(XN-1.)*1.0471975)
      80 Y(M)=RADIX*SINF(ANGLE+(XN-1.)*1.0471975)
      IF(ABSF(X(M))-0.001)90,90,100
      90 X(M)=0.0
      100 IF(ABSF(Y(M))-0.001)110,110,120
      110 Y(M)=0.0
      120 CONTINUE
      130 IF(NICPRT(I)-L)150,150,140
      140 L=L+1
      J=J+6
      GO TO 5
      150 NRPRT(I)=NN
      K=K+NN
      160 CONTINUE
      RETURN
      END(2,1,0,0,0)

```



```

C      SRL HERESY 1 SUBROUTINE SQUARE      VERSION OF 10/2/64
C
C      SUBROUTINE SQUARE
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NOUT=10
      K=0
      DO 260 I=1,NTYPE
      L=1
      J=0
      N=1
10  IF(YO(I,L))20,40,20
20  IF(XO(I,L))60,30,60
30  NERR=1
      WRITE OUTPUT TAPE NOUT,35
35  FORMAT(1H0,69HERROR IN GEOMETRY GENERATOR - INITIAL COORDINATES ARE
      IE IN WRONG SECTOR)
      RETURN
40  ANGLE=0.0
      IF(XO(I,L))70,50,70
50  M=K+N
      X(M)=0.0
      Y(M)=0.0
      GO TO 250
60  TAN=YO(I,L)/XO(I,L)
      ANGLE=ATANF(TAN)
70  RADIX=SQRTF(XO(I,L)*XO(I,L)+YO(I,L)*YO(I,L))
      ANGLE1=ANGLE
80  M=K+N
      X(M)=RADIX*COSEF(ANGLE1)
      Y(M)=RADIX*SINF(ANGLE1)
      IF(ABSF(X(M))-0.001)90,90,100
90  X(M)=0.0
100 IF(ABSF(Y(M))-0.001)110,110,120
110 Y(M)=0.0
120 J=J+1
      GO TO (150,160,170,180,190,200,210,230),J
130 IF(ABSF(ANGLE2-ANGLE1)-0.001)120,120,140
140 N=N+1
      ANGLE1=ANGLE2
      GO TO 80
150 ANGLE2=1.5707963-ANGLE
      GO TO 130
160 ANGLE2=1.5707963+ANGLE
      GO TO 130
170 ANGLE2=3.1415926-ANGLE
      GO TO 130
180 ANGLE2=3.1415926+ANGLE
      GO TO 130
190 ANGLE2=4.7123889-ANGLE
      GO TO 130
200 ANGLE2=4.7123889+ANGLE
      GO TO 130
210 ANGLE2=6.2831852-ANGLE
      IF(ANGLE2-6.2831852)220,120,120
220 IF(ABSF(ANGLE2-6.2831852)-0.001)120,120,130
230 IF(NICPRT(I)-L)250,250,240
240 L=L+1
      N=N+1
      J=0
      GO TO 10
250 NRPT(I)=N
      K=K+N
260 CONTINUE
      RETURN
      END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE CIRCLE      VERSION OF 10/2/64
C
C      SUBROUTINE CIRCLE
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NOUT=10
      K=0
      DO 250 I=1, NTYPE
      L=1
      J=1
20  IF(X0(I,L))70,30,70
30  IF(Y0(I,L))40,60,50
40  ANGLE=4.7123889
      GO TO 140
50  ANGLE=1.5707963
      GO TO 140
50  M=K+J
      X(M)=0.0
      Y(M)=0.0
      GO TO 240
70  IF(X0(I,L))80,80,110
80  IF(Y0(I,L))90,100,100
90  Y0(I,L)=ABSF(Y0(I,L))
      X0(I,L)=ABSF(X0(I,L))
      ANGLE=3.1415926+ATANF(Y0(I,L)/X0(I,L))
      GO TO 140
100 Y0(I,L)=ABSF(Y0(I,L))
      X0(I,L)=ABSF(X0(I,L))
      ANGLE=3.1415926-ATANF(Y0(I,L)/X0(I,L))
      GO TO 140
110 IF(Y0(I,L))120,120,130
120 Y0(I,L)=ABSF(Y0(I,L))
      X0(I,L)=ABSF(X0(I,L))
      ANGLE=6.2831852-ATANF(Y0(I,L)/X0(I,L))
      GO TO 140
130 Y0(I,L)=ABSF(Y0(I,L))
      X0(I,L)=ABSF(X0(I,L))
      ANGLE=ATANF(Y0(I,L)/X0(I,L))
140 RADIX=SQRTF(X0(I,L)*X0(I,L)+Y0(I,L)*Y0(I,L))
      ANGLE1=ANGLE
150 M=J+K
      X(M)=RADIX*COSF(ANGLE1)
      Y(M)=RADIX*SINF(ANGLE1)
      IF(ABSF(X(M))- .001)160,160,170
160 X(M)=0.0
170 IF(ABSF(Y(M))- .001)180,180,190
180 Y(M)=0.0
190 ANGLE2=ANGLE1+THETA(I,L)/57.29578
      IF(ANGLE2-ANGLE-6.2831822)200,200,220
200 IF(ABSF(ANGLE2-ANGLE-6.2831852)-.001)220,220,210
210 ANGLE1=ANGLE2
      J=J+1
      GO TO 150
220 IF(NICPRT(I)-L)240,240,230
230 L=L+1
      J=J+1
      GO TO 20

240 NRPRT(I)=J
      K=K+J
250 CONTINUE
      RETURN
      END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE PRINT          VERSION OF 10/2/64
C
C      SUBROUTINE PRINT
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NOUT=10
      K=0
      WRITE OUTPUT TAPE NOUT,10,{TITLE(I),I=1,12},XCOR
10  FORMAT(1H1,47X26HLISTING OF HERESY GEOMETRY/1H0,23X12A6/1H0,7HXCOR
      1 = F10.3)
      DO 40 I=1,NTYPE
      WRITE OUTPUT TAPE NOUT,20,I
20  FORMAT(1H0,10HROD TYPE  I3/1H0,5(8X1HX,9X1HY,4X))
      L=K+1
      M=K+NRPRT(I)
      WRITE OUTPUT TAPE NOUT,30,((X(N),Y(N)),N=L,M)
30  FORMAT(5(4XF9.4,1XF9.4))
      K=K+NRPRT(I)
40  CONTINUE
      RETURN
      END(2,1,0,0,0)

```

```

C          SRL HERESY 1  SUBROUTINE PUNCH  VERSION OF 10/2/64
C
C          SUBROUTINE PUNCH
C
C          INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
          CALL BFM
          K=5
          NX=0
          N=0
          DO 10 I=1, NTYPE
10      NX=NX+NRPRT(I)
          LOCATE=2500
15      IF(NX-K)40,20,20
20      PUNCH 30,LOCATE,AA(LOCATE),AA(LOCATE+1),AA(LOCATE+2),AA(LOCATE+3),
1AA(LOCATE+4)
30      FORMAT(I12,5F12.6)
          LOCATE=LOCATE+5
          K=K+5
          GO TO 15
40      K=K-5
          M=NX-K
          DO 50 I=1,5
50      BB(I)=0.0
          DO 60 I=1,M
          MX=LOCATE-1+I
60      BB(I)=AA(MX)
          PUNCH 30,LOCATE,(BB(I),I=1,M)
          IF(N)70,70,80
70      LOCATE=9000
          K=5
          N=1
          GO TO 15
80      RETURN
          END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE PLOT      VERSION OF 10/2/64
C
C      SUBROUTINE PLOT
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NOUT=10
      K=0
      KMAX=0
      DO 10 I=1,NTYPE
10  KMAX=KMAX+NRPRT(I)
      LPAGE=1
      WRITE OUTPUT TAPE NOUT,20,LPAGE
20  FORMAT(1H145X29HPLOT OF FINITE HERESY LATTICE,38X5HPAGE 12)
      WRITE OUTPUT TAPE NOUT,30,(TITLE(I),I=1,12)
30  FORMAT(1H0,23X12A6)
      WRITE OUTPUT TAPE NOUT,40,XSCALE,YSCALE,NTYPE
40  FORMAT(1H0,35HEACH INCREMENT ALONG X AXIS EQUALS F10.5,2X2HCM,5X
      135HEACH INCREMENT ALONG Y AXIS EQUALS F10.5,2X2HCM/1H0,5X20HNUMBER
      2 OF ROD TYPES 12)
      IF(NTYPE-25)80,80,50
50  WRITE OUTPUT TAPE NOUT,60,(I,I=1,25)
60  FORMAT(1H0,8X8HROD TYPE,25(2XI2))
      WRITE OUTPUT TAPE NOUT,70,(NRPRT(I),I=1,25)
70  FORMAT(1H ,16HNO RODS/ROD TYPE,25(I4))
      LX=26
      LH=NTYPE
      GO TO 90
80  LX=1
      LH=NTYPE
90  WRITE OUTPUT TAPE NOUT,60,(I,I=LX,LH)
      WRITE OUTPUT TAPE NOUT,70,(NRPRT(I),I=LX,LH)
      L=1
100  LPAGE=LPAGE+1
      WRITE OUTPUT TAPE NOUT,110,(TITLE(I),I=1,12),LPAGE
110  FORMAT(1H1,12A6,40X5HPAGE 12)
      GO TO (120,130,140,150),L
120  DO 125 I=1,26
      KK(I)=-26+I
125  NN(I)=26-I
      GO TO 160
130  DO 135 I=1,26
      KK(I)=I-26
135  NN(I)=1-I
      GO TO 160
140  DO 145 I=1,26
      KK(I)=I-1
145  NN(I)=26-I
      GO TO 160
150  DO 155 I=1,26
      KK(I)=I-1
155  NN(I)=1-I
160  WRITE OUTPUT TAPE NOUT,170,(KK(I),I=1,26)
170  FORMAT(1H0,13HX INTERVALS =,26(1XI3)//12H Y INTERVALS)
      DO 330 N=1,26
      FN=FLOATE(NN(N))
      DO 180 LXY=1,26
180  MM(LXY)=0

```

```

      DO 310 K=1,KMAX
      IF(Y(K)-(FN-.5)*YSCALE)310,190,190
190  IF(Y(K)-(FN+.5)*YSCALE)200,200,310
200  DO 300 I=1,26
      FI=FLOATF(KK(I))
      IF(X(K)-(FI+.5)*XSCALE)210,210,300
210  IF(X(K)-(FI-.5)*XSCALE)300,220,220
220  KRAX=NRPRT(1)
      DO 290 KX=1,NTYPE
      IF(K-KRAX)250,250,230
230  KRAX=KRAX+NRPRT(KX+1)
      GO TO 290
250  MM(I)=KX
      GO TO 310
290  CONTINUE
300  CONTINUE
310  CONTINUE
      WRITE OUTPUT TAPE NOUT,320,NV(N),(MM(K),K=1,26)
320  FORMAT(1H0,10X13,26(1X13))
330  CONTINUE
      L=L+1
      IF(L-4)100,100,340
340  RETURN
      END(2,1,0,0,0)

```

```

C          SRL HERESY 1 SUBROUTINE INGEOM          VERSION OF 10/8/64
C
C          SUBROUTINE INGEOM
C
C          INSERT DIMENSION DECK FOR MAIN PROGRAM 1
C
      CALL EFM
      NJUT=10
      IX=1
      RSQ=RADII**2
      DO 80 I=1,NTYPE
      XDX=WDX(I)*XCOR
      YDY=WDY(I)*XCOR
      XX1=WX1(I)*XCOR
      YY1=WY1(I)*XCOR
      NRPRT(I)=0
      MAXX=RADII/XDX
      X(IX)=0.0
      DO 20 J=1,MAXX
      MAXY=(SQRT(RSQ-X(IX)**2))/YDY
      Y(IX)=0.0
      NRPRT(I)=NRPRT(I)+1
      DO 10 K=1,MAXY
      IK=IX+K
      Y(IK)=Y(IK-1)+YDY
      X(IK)=X(IX)
10  NRPRT(I)=NRPRT(I)+1
      KX=IX+MAXY+1
      X(KX)=X(IX)+XDX
20  IX=KX
      IX=IX-1
      K4=IX-NRPRT(I)+1
      K2=0
      DO 40 J=K4,IX
      IF(X(J))90,40,30
30  K2=K2+1
      K3=IX+K2
      Y(K3)=Y(J)
      X(K3)=-X(J)
40  CONTINUE
      IX=IX+K2
      NRPRT(I)=NRPRT(I)+K2
      K2=0
      K4=IX-NRPRT(I)+1
      DO 60 J=K4,IX
      IF(Y(J))90,60,50
50  K2=K2+1
      K3=IX+K2
      Y(K3)=-Y(J)
      X(K3)=X(J)
60  CONTINUE
      NRPRT(I)=NRPRT(I)+K2
      IX=IX+K2
      K4=IX-NRPRT(I)+1
      DO 70 J=K4,IX
      X(J)=X(J)+XX1
70  Y(J)=Y(J)+YY1
      IX=IX+1
80  CONTINUE

```

```

      GO TO 105
90  NERR=1
    WRITE OUTPUT TAPE NOUT,100
100 FORMAT(1H0,81H A MACHINE ERROR HAS OCCURRED IN GENERATING AN INFIN
    IITE LATTICE - RESTART PROBLEM)
    GO TO 140
105 IX=0
    DO 110 I=1,NTYPE
110 IX=IX+NRPRT(I)
    IF (IX-6500) 140,140,120
120 WRITE OUTPUT TAPE NOUT,130
130 FORMAT(1H0,31H THE NUMBER OF RDDS EXCEEDS 6500)
    NERR=1
140 RETURN
    END(2,1,0,0,0)

```


C	SRL HERESY 1 SUBROUTINE DATPRT	VERSION OF 10/15/64	DATP
C			DATP
	SUBROUTINE DATPRT		DATP
C			
C	INSERT DIMENSION DECK FOR MAIN PROGRAM 1		
C			
	CALL EFM		DATP
	NOUT=10		DATP
	WRITE OUTPUT TAPE NOUT,5,(TITLE(I),I=1,12)		DATP
	5 FORMAT(1H1,23X12A6)		DATP
	WRITE OUTPUT TAPE NOUT,10		DATP
	10 FORMAT(1H0,45X28H SUMMARY OF HERESY INPUT DATA/1H0,53X13H INPUT OPTI		DATP
	IONS)		DATP
	WRITE OUTPUT TAPE NOUT,20,INF,KERNEL,KOOR,NTERM		DATP
	20 FORMAT(1H ,25HFINITE/INFINITE OPTION = 12,5X16H KERNEL OPTION = 12,		DATP
	15X20H COORDINATE OPTION = 12,5X29HPURE THERMAL KERNEL OPTION = 12)		DATP
	IF(KOOR-1)62,50,62		DATP
	50 WRITE OUTPUT TAPE NOUT,60,LAT,KPRNT,KPLOT,KSKIP		DATP
	60 FORMAT(1H ,15HLATTICE TYPE = 12,5X17HGEOLOGY PRINT = 12,5X16HGEO		DATP
	1ETRY PLOT = 12,5X17HGEOLOGY PUNCH = 12)		DATP
	62 IF(KOOR-3)70,65,70		DATPRT
	65 WRITE OUTPUT TAPE NOUT,68,NF(50)		DATP
	68 FORMAT(1H0,25HGEOLOGY TAKEN FROM FILE 14,32H OF PERMANENT TAPE ON		DATP
	1TAPE UNIT 3)		DATP
	70 WRITE OUTPUT TAPE NOUT,80,N1,N2,N3,CR1,CR2,CR3		DATP
	80 FORMAT(1H0,49X20H CONVERGENCE CRITERIA/1H ,30H MATRIX INVERSION ITER		DATP
	ATIONS = 14,5X24HEIGENVALUE ITERATIONS = 15,5X32H CRITICALITY SEAR		DATP
	2H ITERATIONS = 14/1H ,24H INVERSION CONVERGENCE = F7.6,5X25HEIGENVA		DATP
	3LUE CONVERGENCE = F7.6,5X26H CRITICALITY CONVERGENCE = F7.6)		DATP
	WRITE OUTPUT TAPE NOUT,90,NOGRID,XMESH,B(1),B(2),B(3),TAU(1),TAU(2		DATP
	1),TAU(3),TAUR(1),TAUR(2),TAU(3),BSQ,XLSQ,SIGA		DATP
	90 FORMAT(1H0,58X11H KERNEL DATA/1H ,24H NUMBER OF GRID POINTS = 15,5X		DATP
	112H MESH SIZE = F8.3,2X2HCM/1H ,10HB(1) = F6.2,5X10HB(2) = F6		DATP
	2.2,5X10HB(3) = F6.2/1H ,10HTAU(1) = F6.2,5X10HTAU(2) = F6.2,		DATP
	35X10HTAU(3) = F6.2/1H ,10HTAUR(1) = F6.2,5X10HTAUR(2) = F6.2,5X10		DATP
	4HTAUR(3) = F6.2/1H ,11HB SQUARE = F7.6,5X11HL SQUARE = F10.4,5X21H		DATP
	5MODERATOR SIGMA(A) = F12.8)		DATP
	WRITE OUTPUT TAPE NOUT,100,RADREFL,ROEMIN		DATP
	100 FORMAT(1H0,19H REFLECTOR RADIUS = F6.2,11H RHO(MIN) = F6.2)		DATP
	IF(INF)108,108,102		DATP
	102 WRITE OUTPUT TAPE NOUT,104,RADII		DATP
	104 FORMAT(1H0,52H RADIUS BEYOND WHICH NO MORE RODS ARE CONSIDERED =		DATP
	1F8.3)		DATP
	108 IF(LCRIT)220,220,110		DATP
	110 IF(LCRIT-1)120,120,140		DATP
	120 WRITE OUTPUT TAPE NOUT,130,EIGEN		DATP
	130 FORMAT(1H0,60H CRITICALITY SEARCH ON ETA - EIGENVALUE TO BE SEARCHE		DATP
	1D FOR = F10.7)		DATP
	GO TO 190		DATP
	140 IF(LCRIT-2)150,150,170		DATP
	150 WRITE OUTPUT TAPE NOUT,160,EIGEN		DATP
	160 FORMAT(1H0,62H CRITICALITY SEARCH ON GAMMA - EIGENVALUE TO BE SEAR		DATP
	1HED FOR = F10.7)		DATP
	GO TO 190		DATP
	170 WRITE OUTPUT TAPE NOUT,180,EIGEN		DATP
	180 FORMAT(1H0,65H CRITICALITY SEARCH ON BUCKLING - EIGENVALUE TO BE SE		DATP
	1ARCHED FOR = F10.7)		DATP

190 IF(LCRIT-2)200,200,220	DATP
200 WRITE OUTPUT TAPE NOUT,210,KSERCH	DATP
210 FORMAT(1H ,29HROD KIND TO BE SEARCHED ON = I3)	DATP
220 WRITE OUTPUT TAPE NOUT,230	DATP
230 FORMAT(1H0,48X22HDATA FOR EACH ROD KIND)	DATP
WRITE OUTRUT TAPE NOUT,240	DATP
240 FORMAT(1H0, 8HROD KIND,4X5HGAMMA,6X3HSMF,5X12H1./ (GAM+SMF),5X3HETA	DATP
1,5X10HRES. COEF.,5X3HLSQ,5X8HSIGMA(A),6X1HF,6X9HCELL VOL.,3X6HRADI	DATP
2US//)	DATP
DO 250 I=1,NKIND	DATP
XXLL=1./ (GAM(I)+SM(I))	DATP
250 WRITE OUTPUT TAPE NOUT,260,I,GAM(I),SM(I),XXLL,ETAI(I),AL(I),XLSQ2	DATP
I(I),SIGC(I),THERU(I),VCELL(I),RADIUS(I)	DATP
260 FORMAT(1H ,3X12,6XF7.5,3XF7.5,5XF8.6,5XF7.5,4XF8.4,4XF7.1,3XF8.7,	DATP
13XF7.5,3XF9.5,3XF6.4)	DATP
WRITE OUTPUT TAPE NOUT,270	DATP
270 FORMAT(1H0,48X22HDATA FOR EACH ROD TYPE)	DATP
WRITE OUTPUT TAPE NOUT,280	DATP
280 FORMAT(1H0,8HROD TYPE,3X8HROD KIND,3X8HNO. RODS,4X5HGAMMA,6X3HSMF,	DATP
15X14H1./ (GAMMA+SMF)//)	DATP
DO 290 I=1,NTYPE	DATP
XXLL=1./ (GAMMA(I)+SMFO(I))	DATP
290 WRITE OUTPUT TAPE NOUT,300,I,KIND(I),NRPRT(I),GAMMA(I),SMFO(I),	DATP
TXLL	DATP
300 FORMAT(1H ,3X12,9X12,7X14,6XF7.5,3XF7.5,6XF8.6)	DATP
RETURN	DATP
END(2,1,0,0,0)	DATP

Link 2

Link 2 of the program tape consists of 13 FORTRAN II subprograms, 2 SAP subprograms, a package of library function subroutines, and a package of FORTRAN II control subroutines. COMMON STORAGE begins at location 14673₈ in this link.

FORTRAN II and SAP Subprograms

<u>Subprogram Name</u>	<u>Function</u>
MAIN PROG 2	Link 2 - executive routine
XYLIB (SAP)	Library tape routine
MATELM	Slowing down matrices
INVERT	Matrix inversion
EIGENL	Eigenvalue power iteration
EDIT	Edit solution and output result
HEAD	Write heading for search summary
COUT	Output search summary
CONVER	Test for criticality convergence
ETASER	Eta search routine
GAMSER	Gamma search routine
BUCK	Buckling search routine
KERGEN	Kernel generator
LEQ(SAP)	Linear equation program
USKER	Bessel's interpolation routine

The following are 10 small matrix routines used in code:

ABMXVC	Finds max. value of a column vector
DGMPMX	Diagonal mx. multiplied by a square mx.
MXADDG	Square mx. added to a diagonal mx.
MXMPCV	Square mx. multiplied by a column vector
MXMPDG	Square mx. multiplied by diagonal mx.
MXMPMX	Square mx. multiplied by square mx.
MXSBMX	Square mx. subtracted from square mx.
RVMPCV	Row vector multiplied by column vector
RVMPDG	Row vector multiplied by diagonal mx.
RVMPMX	Row vector multiplied by square mx.

FORTRAN II Library Function Subroutines

<u>FORTRAN Name</u>	<u>Function</u>
EXPF (arg)	$e(\text{arg})$
LOG (arg)	$\ln(\text{arg})$
EXPIF (arg)	$E_1(\text{arg})$ integral
SQRTF (arg)	$\sqrt{\text{arg}}$

FORTRAN II Control Subroutines

<u>Subroutine Name</u>	<u>Operation</u>
XLOCF (arg)	Locate variable in storage
LFM	Leave floating point trap
EFM	Enter floating point trap
(FIL)	Format interpretive loader
(DBC)	Decimal to binary conversion
(BDC)	Binary to decimal conversion
(IOH)O	Input - Output Hollerith, Output
(IOH)I	Input - Output Hollerith, Input
(RTN)	Return
(LEV)	Leave
(STH)	Storage to tape Hollerith
(SPH)	Storage to printer Hollerith
CHAIN	Chain search routine
CPTG	CHAIN program tape generator

```

C      DIMENSION DECK FOR SRL HERESY 1 MAIN PROGRAM 2
C
C      DIMENSION DECK FOR MAIN PROGRAM 2      VERSION OF 10/2/64
C
      DIMENSION LL(225),AA(15500),AAA(10250)
      DIMENSION NICPRT(50),NRPRT(50),KIND(50),NF(50)
      DIMENSION B(3),TAU(3),TAUR(3),XLSQ2(50),SIGC(50),VCELL(50),GAM(5
1      ,THERU(50),ETAI(50),AL(50),RADIUS(50),XO(50,3),YO(50,3),
2      THETA(50,3),WDX(50),WDY(50),WX1(50),WY1(50),VO(50),W(8),
3      TITLE(12),BGF(250),BGFR(250),SMGR(250),SMF(250),SM(50),D(50)
4      GAMMA(50),ETA(50),A(50),SMFO(50),FKMAX(50),X(6500),Y(6500)
      DIMENSION T1(50,50),T2(50,50),T3(50,50),T4(50,50),T5(50),T6(50),
1      T(50,50),TR(50,50),SR(50,50),Z(50,50),ABSRB(50),RESABS(50),
2      VRATIO(50),V1(50),V2(50)

C      COMMON LL,AA,AAA

C      EQUIVALENCES FOR LL ARRAY
C
      EQUIVALENCE (INF,LL(1)),(KERNEL,LL(2)),(KDOOR,LL(3)),(N1,LL(4)),
1      (N2,LL(5)),(N3,LL(6)),(NOGRID,LL(7)),(NTHET,LL(8)),(LAT,
2      LL(9)),(NICPRT,LL(10)),(KPRNT,LL(60)),(KPLT,LL(61)),(KSKIP,
3      LL(62)),(NTYPE,LL(63)),(NKIND,LL(64)),(NRPRT,LL(65)),(KIND,
4      LL(115)),(LCRIT,LL(165)),(KSERCH,LL(166)),(NSEPR,LL(167)),
5      (KDAT,LL(168)),(KTEST,LL(169)),(IT1,LL(170)),(IT2,LL(171)),
6      (IT3,LL(172)),(IT4,LL(173)),(IMAX1,LL(174)),(IMAX2,LL(175)),
7      (NF,LL(176))

C      EQUIVALENCES FOR AA ARRAY
C
      EQUIVALENCE (CR1,AA(1)),(CR2,AA(2)),(CR3,AA(3)),(XCOR,AA(4)),
1      (XSCALE,AA(5)),(YSCALE,AA(6)),(RADREL,AA(7)),(ROEMIN,AA(8)),
2      (EIGEN,AA(9)),(XMESH,AA(10)),(BSQ,AA(11)),(XLSQ,AA(12)),
3      (SIGA,AA(13)),(B,AA(14)),(TAU,AA(17)),(TAUR,AA(20)),(TAUTOL,
4      AA(23)),(NERR,AA(24)),(XLSQ2,AA(25)),(SIGC,AA(75)),(VCELL,
5      AA(125)),(GAM,AA(175)),(THERU,AA(225)),(ETAI,AA(275)),(AL,
6      AA(325)),(RADIUS,AA(375)),(XO,AA(425)),(YO,AA(575)),(THETA,
7      AA( 725)),(WDX,AA( 875)),(WDY,AA( 925)),(WX1,AA( 975)),(WY1,
8      AA(1025)),(RADII,AA(1075)),(A1,AA(1076)),(A2,AA(1077)),(FK1,
9      AA(1078)),(FK2,AA(1079)),(VO,AA(1080)),(W,AA(2492))
      EQUIVALENCE (BGF,AA(1142)),(BGFR,AA(1392)),(SMGR,AA(1642)),
1      (SMF,AA(1892)),(SM,AA(2142)),(D,AA(2192)),(GAMMA,AA(2242)),
2      (ETA,AA(2292)),(A,AA(2342)),(SMFO,AA(2392)),(FKMAX,AA(2442))
3      (X,AA(2500)),(Y,AA(9000)),(TITLE,AA(1130)),(T,AA(2500)),
4      (TR,AA(5000)),(SR,AA(7500)),(Z,AA(10000)),(T5,AA(12500)),
5      (T6,AA(12550))

C      EQUIVALENCES FOR AAA ARRAY
C
      EQUIVALENCE (T1,AAA(1)),(T2,AAA(2501)),(T3,AAA(5001)),(T4,
1      AAA(7501)),(ABSRB,AAA(10001)),(RESABS,AAA(10051)),(VRATIO,
2      AAA(10101)),(V1,AAA(10151)),(V2,AAA(10201))
C

```

```
SRL-HERESY 1    MAIN PROGRAM 2    VERSION OF 10/19/64
SUBROUTINE MAIN
INSERT DIMENSION DECK FOR MAIN PROGRAM 2
```

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DIME

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REM	FORTAN II SUBROUTINE XYLIB	VERSION OF 10/21/64	XYLB
REM			XYLB
REM	TAPE CONTROL SUBROUTINE FOR SRL-HERESY I GEOMETRY TAPES		XYLB
REM			XYLB
FUL		P C G	XYLB
ORG 0		R A E	XYLB
MZE *,,ENTRY-2		O R N	XYLB
PZE		G D E	XYLB
PZE BREAK,,0		R R	XYLB
PZE 32562		A A	XYLB
BCD 1XYLIB		M T	XYLB
PZE EP		O	XYLB
ENTRY EQU *		R	XYLB
ORG 0			XYLB
REL			XYLB
STORE PZE			XYLB
PZE			XYLB
PZE			XYLB
PZE 0			XYLB
FILE PZE 0			
PZE 0			
SCALE PZE 0			
EP	SXD STORE,1	STORE INDEX REGS.	XYLB
	SXD STORE+1,2		XYLB
	SXD STORE+2,4		XYLB
	SXD EOF,2		XYLB
	TXI *+1,2,1		XYLB
	CLA 1,4		
	STA *+1		
	CLA **		
	STO STORE+3	STORE WRITE INDICATOR	
	CLA 2,4		
	STA *+1		
	CLA **		
	STO STORE+4	STORE FILE NUMBER	
	CLA 3,4		
	STA *+1		
	CLA **		
	STO STORE+5	STORE READ-WRITE INDICATOR	
	CLA 4,4		
	STA *+1		
	CLA **		
	STO STORE+6	STORE SCALE FACTOR	
	CLA STORE+4	GET FILE NUMBER	
	TNZ RDID	BRANCH ON NON ZERO TO LIBRARY TAPE	XYLB
	CLA STORE+5	GET READ-WRITE INDICATOR	
	TNZ WRITE	BRANCH ON NON ZERO TO WRITE ROUTINE	XYLB
	CLA ZERO		XYLB
	LXA ONE,1		XYLB
	RTB 2	READ BINARY TAPE 2	XYLB
	CAD 29839,1		XYLB
	TRA *+3		XYLB
	TRA ERROR	TRANSFER TO HPR 550	XYLB
	TRA ERROR	TRANSFER TO HPR 550	XYLB
	TIX *-4,1,1		XYLB
	SLW CHECK		
	CLA CHECK		
	CPY CHECK		XYLB
	SUB CHECK		XYLB

	TZE **2			XYLB
	HPR *	CHECKSUM ERROR - HPR 424		XYLB
	REW 2	REWIND TAPE 2		XYLB
WRITE	TRA RETN	TRANSFER TO EXIT		XYLB
	LXA ONE,1			XYLB
	CLA ZERO			XYLB
	WTB 2	WRITE BINARY TAPE 2		XYLB
	CAD 29839,1			XYLB
	TIX *-1,1,1			XYLB
	SLW CHECK			
	CPY CHECK			XYLB
	REW 2	REWIND TAPE 2		XYLB
	TRA RETN	TRANSFER TO EXIT		XYLB
RDID	RTB 3	BEGIN FILE SEARCH ON TAPE 3		XYLB
	CPY FILE1			XYLB
	TRA **3			XYLB
	TRA TEST-1	TRANSFER TO MAKE NEW TAPE		
	TRA TEST-1			
	CLA FILE			XYLB
	SUB FILE1			XYLB
	TZE RIGHT	FILE IS CORRECT		
	CLA FOE	FILE IS NOT CORRECT		
	SUB FILE1			
	TNZ RDID	TRANSFER TO READ NEW RECORD		
EOF	TIX REWTP,2,0	REWIND TAPE ONCE ON FOE MARK		
	TRA TEST	FILE NOT ON TAPE,TRANSFER TO TEST		
RIGHT	CLA STORE+5	FILE IS CORRECT,GET READ-WRITE INDICATOR		
	TZE RDTP	TRANSFER TO TAPE 3 READ		XYLB
	CLA STORE+3	GET NEW COORDINATE INDICATOR		
	TNZ **2	TRANSFER TO NEW COORDINATE WRITE		XYLB
	TRA RETN	TRANSFER TO EXIT		XYLB
	CLA ZERO			XYLB
	LXA ONE,1			XYLB
	WTB 3	WRITE NEW COORDINATES IN OLD FILE		
	CAD 29839,1			XYLB
	TIX *-1,1,1			XYLB
	SLW CHECK			
	CPY CHECK			XYLB
	TRA RETN			XYLB
RDTP	LXA ONE,1	READ FILE FROM TAPE 3		XYLB
	CLA ZERO			XYLB
	RTB 3	READ COORDINATES FROM OLD FILE		XYLB
	CAD 29839,1			XYLB
	TRA **3			XYLB
	TRA ERROR	TRANSFER TO HPR 550		
	TRA ERROR	TRANSFER TO HPR 550		
	TIX *-4,1,1			XYLB
	SLW CHECK			
	CLA CHECK			
	CPY CHECK			XYLB
	SUB CHECK			XYLB
	TZE **2			XYLB
	HPR *	CHECKSUM ERROR - HPR 507		XYLB
	LXD ONE,1			XYLB
	EFM			XYLB
	LDQ SCALE	PLACE SCALE FACTOR IN MQ		XYLB
	FMP 29839,1	FLOATING MULTIPLY EACH COORDINATE VALUE		XYLB
	STO 29839,1	STORE BACK IN COORDINATE ARRAY		XYLB
	TIX *-3,1,1			XYLB

LFM			XYLB
TRA RETN			XYLB
REWTP REW 3	REWIND TAPE 3		XYLB
TRA RDID			XYLB
REW 3	TAPE 3 REWIND FOR NEW TAPE		
TEST CLA STORE+3			
TNZ *+5	TRANSFER TO WRITE NEW COORDINATE FILE		XYLB
LXD STORE,1			XYLB
LXD STORE+1,2			XYLB
LXD STORE+2,4			XYLB
HTR 5,4	ERROR - FILE IS NOT ON TAPE HTR 5,4		XYLB
BST 3	BACKSPACE TAPE 3 ONE RECORD		
WTB 3			XYLB
CPY FILE	WRITE NEW FILE NUMBER		XYLB
CLA ZERO			XYLB
LXA ONE,1			XYLB
WTB 3	WRITE NEW COORDINATES		XYLB1
CAD 29839,1			XYLB1
TIX *-1,1,1			XYLB1
SLW CHECK			
CPY CHECK	WRITE CHECKSUM ON TAPE 3		
WTB 3			
CPY FDE	WRITE FDE MARK ON TAPE 3		
WEF 3	END FILE TAPE 3		XYLB1
REW 3			
TRA RETN			XYLB1
ERROR HPR *	READ ERROR ON TAPE 2 OR 3 HPR 550		
RETN LXD STORE,1			XYLB1
LXD STORE+1,2			XYLB1
LXD STORE+2,4			XYLB1
TRA 5,4	EXIT TO MAIN PROGRAM		XYLB1
ONE PZE 13000			XYLB1
FILE1 PZE 0			XYLB1
ZERO PZE 0			XYLB1
CHECK PZE 0			XYLB1
FDE PZE 64625			
BREAK EQU *			XYLB1
END **			XYLB1

```

C      SRL HERESY 1 SUBROUTINE MATELM      VERSION OF 10/2/64
C
C      SUBROUTINE MATELM
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EFM
      PI=3.1415926
      KMN=1
      DO 110 I=1,NTYPE
      N=NRPRT(I)
      KR=1
      KMNN=KMN+N-1
      DO 100 J=1,NTYPE
      T1(J,I)=0.0
      T2(J,I)=0.0
      T3(J,I)=0.0
      T4(J,I)=0.0
      IF(RADRFL)10,10,20
10  IRL=0
      GO TO 40
20  ROERL=SQRTF(X(KMN)**2+Y(KMN)**2)
      IF(ROERL-ROEMLN)10,10,30
30  ROEIM=2.*RADRFL-ROERL
      THET=ROEIM/ROERL
      IRL=1
40  DO 90 K=KMN,KMNN
      IF(KR-K)60,50,60
50  T1(J,I)=T1(J,I)+BGF(1)
      T2(J,I)=T2(J,I)+BGFR(1)
      T3(J,I)=T3(J,I)+SMGR(1)
      T4(J,I)=T4(J,I)+SMFO(I)
      GO TO 70
60  DX=SQRTF((X(KR)-X(K))**2+(Y(KR)-Y(K))**2)
      FF1=1.0
      CALL USKER(DX,FF1,I,J)
70  IF(IRL-1)90,80,90
80  XIMAG=THET*X(K)
      YIMAG=THET*Y(K)
      DX=SQRTF((X(KR)-XIMAG)**2+(Y(KR)-YIMAG)**2)
      FF1=-SQRTF(ROEIM/ROERL)
      CALL USKER(DX,FF1,I,J)
90  CONTINUE
100 KR=KR+NRPRT(J)
110 KMN=KMN+N
      RETURN
      END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE USKER      VERSION OF 10/2/54
C
C      SUBROUTINE USKER(DX,FIMREL,I,J)
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EFM
      NTEMP=2
      NOUT=10
      FNOGR=NOGRID
      IF(DX-XMESH*(FNOGR-3.))30,30,10
10  WRITE OUTPUT TAPE NOUT,20
20  FORMAT(1H0,55HERROR - KERNEL SIZE IS TOO SMALL FOR LATTICE BEING U
      1SED)
      GO TO 70
30  M=DX/XMESH+1.0
      IF(M-2)40,60,60
40  WRITE OUTPUT TAPE NOUT,50
50  FORMAT(1H0,32HERROR - KERNEL MESH IS TOO LARGE)
      GO TO 70
60  U=(MODF(DX,XMESH))/XMESH
      P=U*(U-1.)/4.
      Q=1.-U-P
      R=U-P
      F1F=FIMREL*(P*(BGF(M-1)+BGF(M+2))+Q*BGF(M)+R*BGF(M+1))
      T1(J,I)=T1(J,I)+F1F
      F2F=FIMREL*(P*(BGFR(M-1)+BGFR(M+2))+Q*BGFR(M)+R*BGFR(M+1))
      T2(J,I)=T2(J,I)+F2F
      F3F=FIMREL*(P*(SMGR(M-1)+SMGR(M+2))+Q*SMGR(M)+R*SMGR(M+1))
      T3(J,I)=T3(J,I)+F3F
      F4F=FIMREL*(P*(SMF(M-1)+SMF(M+2))+Q*SMF(M)+R*SMF(M+1))
      T4(J,I)=T4(J,I)+F4F
      RETURN
70  WRITE TAPE NTEMP,(AA(I),I=2500,15500)
      REWIND NTEMP
      WRITE OUTPUT TAPE NOUT,80
80  FORMAT(1H ,70HTHIS ERROR OCCURS IN FORMING THE MATRICES FROM THE K
      1ERNEL AND GEOMETRY/1H ,84HAN ATTEMPT WILL BE MADE TO RUN SUCCESSIV
      2E PROBLEMS, BUT IT IS UNLIKELY THEY WILL RUN)
      PRINT 90
90  FORMAT(1H ,26HPROBLEM DID NOT RUN TO END)
      CALL CHAIN (1,1)
      END(2,1,0,0,0)

```

```

C          SRL HERESY 1  SUBROUTINE INVERT  VERSION OF 10/2/64
C
C          SUBROUTINE INVERT
C
C          INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
          CALL EFM
          NOUT = 10
          IT1=0
          NXN=0
          DO 70 I=1,NTYPE
          DO 70 J=1,NTYPE
          T2(I,J)=T1(I,J)
          IF(I-J)60,50,60
50  T3(I,J)=1.0
          GO TO 70
60  T3(I,J)=0.0
70  CONTINUE
          KIAX=XLOC(T2(1,1))-XLOC(T2(1,2))
          CALL LEQ(T2,T3,NTYPE,NTYPE,KIAX,KIAX,DET)
75  CALL MXMPMX(NTYPE,T1,T3,T4)
          DO 100 I=1,NTYPE
          DO 100 J=1,NTYPE
          IF(I-J)90,80,90
80  IF(ABS(T4(I,J))-1.0)-CR1)100,100,110
90  IF(ABS(T4(I,J))-CR1)100,100,110
100 CONTINUE
          RETURN
110 IT1=IT1+1
          IF(IT1-N1)180,180,120
120 IF(NXN)125,125,170
125 WRITE OUTPUT TAPE NOUT,130,IT1,CR1
130 FORMAT(1H1,72HTHE SPECIFIED NUMBER OF ITERATIONS ON MATRIX INVERSI
ION HAS BEEN EXCEEDED/1H ,26HNO ITERATIONS COMPLETED = I4,5X24HCINV
2ERGEANCE CRITERION = F10.7/1H0,29HMATRIX TO BE INVERTED FOLLOWS)
          DO 140 I=1,NTYPE
140 WRITE OUTPUT TAPE NOUT,150,(T1(I,J),J=1,NTYPE)
150 FORMAT(1H0,8E14.7/(1H ,8E14.7))
          N1=N1+10
160 NXN=1
          GO TO 180
170 PRINT 175
175 FORMAT(1H ,26HPROBLEM DID NOT RUN TO END)
          CALL CHAIN(1,1)
180 DO 210 I=1,NTYPE
          DO 210 J=1,NTYPE
          T2(I,J)=T3(I,J)
          IF(I-J)200,190,200
190 T4(I,J)=2.-T4(I,J)
          GO TO 210
200 T4(I,J)=-T4(I,J)
210 CONTINUE
          CALL MXMPMX(NTYPE,T2,T4,T3)
          GO TO 75
          END(2,1,0,0,0)

```

```

C          SRL HERESY 1  SUBROUTINE EIGEN          VERSION OF 10/2/64
C
C          SUBROUTINE EIGEN1
C
C          INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EPM
      NOUT=10
      NN=2
      DO 230 I=1,NTYPE
230  V1(I)=V0(I)
      IT3=0
      IT2=0
      CALL ABMXVC(NTYPE,V1,A1,IMAX1)
      FK2=1.0
240  CALL MXMPCV(NTYPE,T1,V1,V2)
      CALL ABMXVC(NTYPE,V2,A2,IMAX2)
      IF(V1(IMAX2))250,320,250
250  FK2=V2(IMAX2)/V1(IMAX2)
      GO TO (260,300),NN
260  IT3=IT3+1
      IF(IT3-10)300,300,270
270  IT3=0
      DO 280 I=1,NTYPE
280  VRATIO(I)=V2(I)/A2
      WRITE OUTPUT TAPE NOUT,290,IT2,FK2,((V2(I),VRATIO(I)),I=1,NTYPE)
290  FORMAT(1H1,102HTHIS PROBLEM HAS EXCEEDED THE MAXIMUM NUMBER OF ITE
      RATIONS SPECIFIED FOR THE EIGENVALUE ITERATION (N2)/1H0,27HCURRENT
      2 ITERATION NUMBER = 14,5X21HCURRENT EIGENVALUE = F10.7/1H0,10X11HE
      3IGENVECTOR,9X5HV2/A2/(1H ,10XE12.5,5XE12.4))
      IF(IT2-2*N2)300,300,292
292  PRINT 295
295  FORMAT(1H ,26HPROBLEM DID NOT RUN TO END)
      CALL CHAIN(1,1)
300  DO 310 I=1,NTYPE
      IF(ABS(V1(I)-V2(I))/FK2)-CR2*A1)310,310,320
310  CONTINUE
      RETURN
320  IT2=IT2+1
      IF(IT2-N2)340,340,330
330  NN=1
340  DO 350 I=1,NTYPE
350  V1(I)=V2(I)
      A1=A2
      IMAX1=IMAX2
      FK1=FK2
      GO TO 240
      END(2,1,0,0,0)

```

```

C      SRL-HERESY 1 SUBROUTINE HEAD      VERSION OF 10/2/64
C
C      SUBROUTINE HEAD
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EFM
      NOUT=10
      WRITE OUTPUT TAPE NOUT,10,EIGEN
10  FORMAT(1H0,29HSUMMARY OF CRITICALITY SEARCH/1H ,32HEIGENVALUE BEIN
      16 SEARCHED FOR = F10.7)
      IF(LCRIT-1)20,20,40
20  WRITE OUTPUT TAPE NOUT,30
30  FORMAT(1H ,12HITERATION NO,5X18HCURRENT EIGENVALUE,5X19HPREVIOUS E
      1IGENVALUE,5X17HCURRENT ETA VALUE,5X26HROD KIND BEING SEARCHED ON)
      GO TO 90
40  IF(LCRIT-2)50,50,70
50  WRITE OUTPUT TAPE NOUT,60
60  FORMAT(1H ,12HITERATION NO,5X18HCURRENT EIGENVALUE,5X19HPREVIOUS E
      1IGENVALUE,5X19HCURRENT GAMMA VALUE,5X26HROD KIND BEING SEARCHED ON
      2)
      GO TO 90
70  WRITE OUTPUT TAPE NOUT,80
80  FORMAT(1H ,12HITERATION NO,5X18HCURRENT EIGENVALUE,5X19HPREVIOUS E
      1IGENVALUE,5X22HCURRENT BUCKLING VALUE)
90  RETURN
      END(2,1,0,0,0)

```

C SRL-HERESY 1 SUBROUTINE COUT VERSION OF 10/8/64
 C
 C SUBROUTINE COUT
 C
 C INSERT DIMENSION DECK FOR MAIN PROGRAM 2
 C
 CALL EFM
 NOUT=10
 KSERCH=KSERCH
 W(1)=W(1)
 FFK1=W(1)
 FFK2=FK2*TAUTOL
 IF(LCRIT-1)10,10,30
 10 WRITE OUTPUT TAPE NOUT,20,IT4,FFK2,FFK1,ETAI(KSERCH),KSERCH
 20 FORMAT(1H ,5XI2,14XF10.7,13XF11.7,13XF9.6,21XI2)
 GO TO 80
 30 IF(LCRIT-2)40,40,60
 40 WRITE OUTPUT TAPE NOUT,50,IT4,FFK2,FFK1,GAM(KSERCH),KSERCH
 50 FORMAT(1H ,5XI2,14XF10.7,13XF11.7,14XF9.4,22XI2)
 GO TO 80
 60 WRITE OUTPUT TAPE NOUT,70,IT4,FFK2,FFK1,BSQ
 70 FORMAT(1H ,5XI2,14XF10.7,13XF11.7,15XF10.8)
 80 RETURN
 END(2,1,0,0,0)

```

C      SRL-HERESY 1  SUBROUTINE CONVER      VERSION OF 10/2/64
C
C      SUBROUTINE CONVER
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EPM
      NOUT=10
      IF(ABS(F(EIGEN-FK2*TAUTOL)-CR3)10,10,20
10  IT1=2
      RETURN
20  IF(IT4-N3)50,50,30
30  WRITE OUTPUT TAPE NOUT,40
40  FORMAT(1H ,84H THE MAXIMUM NUMBER OF ITERATIONS HAVE BEEN EXCEEDED
      1IN THE SEARCH WITHOUT CONVERGING/1H ,61H THIS PROBLEM WILL BE SKIPP
      2ED OVER AND SUCCESSIVE PROBLEMS RUN)
      IT1=3
      RETURN
50  IT1=1
      RETURN
      END(2,1,0,0,0)

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```

C          SRL-HERESY 1      SUBROUTINE GAMSER II   VERSION OF 10/2/64
C
SUBROUTINE GAMSER
C  INSERT DIMENSION DECK FOR MAIN PROGRAM 2
CALL EFM
KSERCH=KSERCH
IF(IT4)10,10,30
10 IF(NF(KSERCH))11,11,12
11 XLIMIT=0.0
GO TO 13
12 XLIMIT=(1./(12.56637*D(KSERCH)))*LOGF(XLSQ2(KSERCH)/((1.+BSQ*
1XLSQ2(KSERCH))*VCELL(KSERCH)))+.228714/D(KSERCH)
13 GAMX=1./(GAM(KSERCH)+SM(KSERCH))
W(2)=FK2
FX=(EIGEN-FK2*TAUTOL)/EIGEN
GAM1=GAMX*(1.-FX)
XXXX=1./GAM1
IF(XXXX-XLIMIT)50,50,15
15 GAM(KSERCH)=1./GAM1-SM(KSERCH)
DO 20 I=1,NTYPE
N=KIND(I)
20 GAMMA(I)=GAM(N)
IT4=IT4+1
W(1)=FK2*TAUTOL
RETURN
30 GAM2=1./(GAM(KSERCH)+SM(KSERCH))
GAM1=GAM2+(GAM2-GAMX)*(EIGEN-FK2*TAUTOL)/((FK2-W(2))*TAUTOL)
XXXX=1./GAM1
IF(XXXX-XLIMIT)50,50,35
35 GAM(KSERCH)=1./GAM1-SM(KSERCH)
W(2)=FK2
GAMX=GAM2
DO 40 I=1,NTYPE
N=KIND(I)
40 GAMMA(I)=GAM(N)
IT4=IT4+1
W(1)=FK2*TAUTOL
RETURN
50 GAM1=XLIMIT+.0001
IF(IT4)15,15,35
END(2,1,0,0,0)

```

```

C      SRL HERESY 1 SUBROUTINE ETASER      VERSION OF 10/2/64
C
C      SUBROUTINE ETASER
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EFM
      KSERCH=KSERCH
      IF(IT4)10,10,30
10  ETAIX=ETAI(KSERCH)
      W(3)=FK2
      W(1)=FK2*TAUTOL
      FX=(EIGEN-FK2*TAUTOL)/EIGEN
      ETAI(KSERCH)=ETAI(KSERCH)*(1.+FX)
      DO 20 I=1,NTYPE
      N=KIND(I)
20  ETA(I)=ETAI(N)
      IT4=IT4+1
      RETURN
30  ETMX=ETAI(KSERCH)
      ETAI(KSERCH)=ETMX+(EIGEN-FK2*TAUTOL)*(ETMX-ETAIX)/((FK2-W(3))*
1TAUTOL)
      ETAIX=ETMX
      W(3)=FK2
      DO 40 I=1,NTYPE
      N=KIND(I)
40  ETA(I)=ETAI(N)
      IT4=IT4+1
      W(1)=FK2*TAUTOL
      RETURN
      END(2,1,0,0,0)

```

```

C      SRL HERESY I SUBROUTINE BUCK      VERSIO OF 10/8/64
C
C      SUBROUTINE BUCK
C
C      * INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
      CALL EFM
      IF(IT4)10,10,40
10  BSQ2=BSQ
      IF(FK2*TAUTOL-EIGEN)20,90,30
20  BSQ=BSQ-2.5E-04
      NUM=-1
      GO TO 80
30  BSQ=BSQ+2.5E-04
      NUM=+1
      GO TO 80
40  BSQ2=BSQ
      IF(FK2*TAUTOL-EIGEN)50,90,60
50  IF(NUM)20,70,70
60  IF(NUM)70,70,30
70  BSQ=BSQ2+(BSQ2-BSQ1)*(FK2*TAUTOL-EIGEN)/(FFK1-FK2*TAUTOL)
      NUM=0
80  FFK1=FK2*TAUTOL
      TAUTOL=EXPF(-BSQ*(B(1)*TAU(1)+B(2)*TAU(2)+B(3)*TAU(3)))
      CALL KERGEN
      IT4=IT4+1
      BSQ1=BSQ2
      W(1)=FFK1
      DO 100 I=1,NKIND
      D(I)=XLSQ2(I)*SIGC(I)
      XYLSQ=XLSQ2(I)/(1.+BSQ*XLSQ2(I))
      IF(NF(I))100,100,95
95  GAM(I)=XLSQ2(I)*(1.-THERU(I))/(D(I)*VCELL(I)*THERU(I))+
      1(1./((12.56637*D(I)))*LOGF(XYLSQ/VCELL(I))+.228714/D(I)-SM(I)
100 CONTINUE
      DO 110 I=1,NTYPE
      NX=KIND(I)
      GAMMA(I)=GAM(NX)
110 CONTINUE
90  RETURN
      END(2,1,0,0,0)

```

C SRL HERESY 1 SUBROUTINE KERGEN VERSION OF 10/2/64

C

SUBROUTINE KERGEN

C

C

INSERT DIMENSION DECK FOR MAIN PROGRAM 2

C

CALL EFM

FIZCON(1)=.0045313

FIZCON(2)=.0360768

FIZCON(3)=.2659732

FIZCON(4)=1.2067492

FIZCON(5)=3.0899424

FIZCON(6)=3.5156229

FIZCON(7)=1.0

FKZCN1(1)=.0000074

FKZCN1(2)=.0001075

FKZCN1(3)=.00262698

FKZCN1(4)=.0348859

FKZCN1(5)=.23069756

FKZCN1(6)=.42278420

FKZCN1(7)=-.57721566

FKZCN2(1)=.00053208

FKZCN2(2)=-.0025154

FKZCN2(3)=.00587872

FKZCN2(4)=-.01062446

FKZCN2(5)=.02189568

FKZCN2(6)=-.07832358

FKZCN2(7)=1.25331414

DENL=1.+BSQ*XL SQ

XXLSQ=XL SQ/DENL

SIGMA=SIGA*DENL

ZLSQ=SQRTF(XXLSQ)

CC=1./((6.2831852*XXLSQ*SIGMA)

IF(ABSF(B(2)))260,20,10

10 IF(ABSF(B(3)))260,30,40

20 NN=1

GO TO 50

30 NN=2

GO TO 50

40 NN=3

50 DO 310 I=1,NOGRID

FJM1=I-1

DX=FJM1*XMESH

ARG=DX/ZLSQ

IF(ARG)260,60,70

60 IFPATH=3

GO TO 140

70 IFPATH=1

FKZER=0.0

IF(ARG-2.)80,80,120

80 FIZARG=(ARG/3.75)**2

FIZER=0.0

DO 90 J=1,7

90 FIZER=FIZER*FIZARG+FIZCON(J)

100 FKZARG=(ARG/2.)**2

DO 110 J=1,7

110 FKZER=FKZER*FKZARG+FKZCN1(J)

FKZER=FKZER-LOGF(.5*ARG)*FIZER

GO TO 140

```

120 FKZARG=2./ARG
    DO 130 J=1,7
130 FKZER=FKZER*FKZARG+FKZCN2(J)
    FKZER=FKZER/(SQRTF(ARG)*EXPF(ARG))
140 SMGR(I)=0.0
    D2BY4=DX**2/4.
    DO 150 J=1,NN
150 SMGR(I)=SMGR(I)+B(J)*EXPF(-D2BY4/TAUR(J))/(TAUR(J)*12.5663706)
    DO 160 J=1,NN
160 WTAU(J)=TAU(J)
170 WBG=0.0
    GO TO (210,210,180,180),IFPATH
180 DO 190 J=1,NN
190 WBG=WBG+(B(J)/2.)*CC*EXPF(WTAU(J)/XXLSQ)*(-EXPIF(-WTAU(J)/XXLSQ))
    GO TO (260,260,200,280),IFPATH
200 BGF(I)=WBG
    IFPATH=4
    GO TO 240
210 DO 220 J=1,NN
220 WBG=WBG+CC*EXPF(WTAU(J)/XXLSQ)*B(J)*(FKZER+(WTAU(J)/2./XXLSQ)*
    1EXPF(-D2BY4/WTAU(J))+(.5+D2BY4/(2.*XXLSQ))*EXPIF(-D2BY4/WTAU(J)))
    GO TO (230,280,260,260),IFPATH
230 BGF(I)=WBG
    IFPATH=2
240 DO 250 J=1,NN
250 WTAU(J)=TAU(J)-TAUR(J)
    GO TO (260,170,260,170),IFPATH
260 PRINT 270
270 FORMAT(1H0,54H A MACHINE ERROR HAS OCCURRED IN THE SUBROUTINE KERG
    1EN/1H ,35H RESTART PROBLEM FROM THE BEGINNING)
    NERR=1
    GO TO 330
280 BGFR(I)=WBG
    SMF(I)=CC*FKZER
    IF(NTHER)300,300,290
290 BGFR(I)=SMF(I)
    BGFR(I)=SMF(I)
    SMGR(I)=1.
300 CONTINUE
310 CONTINUE
320 CONTINUE
330 RETURN
    END(2,1,0,0,0)

```

```

C      SRL-HERESY I      SUBROUTINE EDIT      VERSION OF 8/10/65
C
C      SUBROUTINE EDIT
C
C      INSERT DIMENSION DECK FOR MAIN PROGRAM 2
C
C      CALL EFM
C      NOUT=10
C      KSERCH=KSERCH
C      DO 10 I=1,NTYPE
10  ABSRB(I)=V2(I)/A2
C      CALL RVMPDG(NTYPE,ABSRB,ETA,T5)
C      CALL MXMPCV(NTYPE,SR,T5,T6)
C      CALL RVMPDG(NTYPE,T6,A,T5)
C      FINK2=FK2*TAUTOL
C      DO 20 I=1,NTYPE
20  RESABS(I)=T5(I)/FINK2
C      IF(LCRIT)28,28,22
22  IF(NSEPR)28,28,24
24  WRITE OUTPUT TAPE NOUT,26,IT4
26  FORMAT(1H1,32HCRITICALITY SEARCH ITERATION NO I4)
C      GO TO 32
28  WRITE OUTPUT TAPE NOUT,29
29  FORMAT(1H1)
32  WRITE OUTPUT TAPE NOUT,30,IT2,FK2,FK1
30  FORMAT(1H0,24HNO EIGENVALUE ITERATIONS,5X10HEIGENVALUE,5X19HPREVIO
1US EIGENVALUE/1H ,16XI3,10XF10.7,9XF11.7)
C      WRITE OUTPUT TAPE NOUT,35
35  FORMAT(1H0,45X29HABSORPTIONS FOR EACH ROD TYPE)
C      WRITE OUTPUT TAPE NOUT,40
40  FORMAT(1H , 8HROD TYPE,5X12HNO RODS/TYPE, 5X7HROD SMF, 8X5HGAMMA,1
10X3HETA,10X1HA, 8X13HTHERMAL ABSRB, 5X15HRESONANCE ABSRB)
C      DO 50 I=1,NTYPE
50  WRITE OUTPUT TAPE NOUT,60,I,NRPRT(I),SMFO(I),GAMMA(I),ETA(I),A(I),
1ABSRB(I),RESABS(I)
60  FORMAT(1H ,3XI2,11XI4,10XF7.5,5XF11.6,5XF7.5,5XF7.4,7XF9.6,10XF9.6
1)
C      IF(LCRIT-2)550,500,550
500 IF(NF(KSERCH))520,520,510
510 I=KSERCH
C      THERU(I)=1./((1.+VCELL(I)*SIGC(I)*(GAM(I)+SM(I)-.228714/D(I)-
1(1./((12.56637*D(I)))*LOGF(XLSQ2(I)/((1.+BSQ*XLSQ)*VCELL(I))))))
C      GO TO 550
520 THERU(KSERCH)=0.0
550 CALL RVMPDG(NTYPE,FKMAX,ETA,T5)
C      CALL RVMPCV(NTYPE,FKMAX,ABSRB,HI)
C      CALL RVMPCV(NTYPE,T5,ABSRB,HETAI)
C      AVETA=HETAI/HI
C      CALL RVMPDG(NTYPE,FKMAX,A,T5)
C      CALL RVMPMX(NTYPE,T5,SR,T6)
C      DO 70 I=1,NTYPE
70  T6(I)=FKMAX(I)-T6(I)
C      SYSYSY=0.0
C      UYUYUY=0.0
C      VYVYVY=0.0
C      CALL RVMPDG(NTYPE,ETA,T6,T5)
C      CALL RVMPCV(NTYPE,ABSRB,T5,SYSYSY)
C      CALL RVMPCV(NTYPE,ABSRB,FKMAX,UYUYUY)
C      CALL RVMPCV(NTYPE,RESABS,FKMAX,VYVYVY)
C      D2OABS=0.0
C      DO 75 I=1,NKIND

```

```

      RPX=0.0
      DO 73 J=1,NTYPE
        IF(KIND(J)-I)73,72,73
72    RPX=RPX+ABSRB(J)*FKMAX(J)
73    CONTINUE
        IF(THERU(I))430,430,74
74    D2OABS=D2OABS+RPX*(1.-THERU(I))/THERU(I)
75    CONTINUE
76    XMOD2=SYSYSY/FINK2-UYUYUY
        THUT=UYUYUY/(UYUYUY+D2OABS)
        REESPR=(UYUYUY+XMOD2)/(UYUYUY+VYVYVY+XMOD2)
        XLEAK=XMOD2-D2OABS
        XNLEAK=(UYUYUY+D2OABS)/(UYUYUY+XMOD2)
        FKMASB=0.0
        SMFKMX=0.0
        DO 80 I=1,NTYPE
          SMFKMX=SMFKMX+FKMAX(I)
80    FKMASB=FKMAX(I)*ABSRB(I)+FKMASB
          FKMASB=FKMASB/SMFKMX
          WRITE OUTPUT TAPE NOUT,90,FINK2,FKMASB,AVETA,THUT,REESPR,D2OABS,
            XLEAK,XNLEAK
90    FORMAT(1H0,46HEIGENVALUE CORRECTED FOR FAST AXIAL LEAKAGE = F10.8,
            15X25HAVERAGE ROD ABSORPTION = F10.7/1H0,14HAVERAGE ETA = F10.7,5X2
            22HTHERMAL UTILIZATION = F10.8,5X31HRESONANCE ESCAPE PROBABILITY =
            3F10.8/1H0,39HAPPROX MODERATOR THERMAL ABSORPTIONS = F9.5,5X16HTOTA
            4L LEAKAGE = F9.4,5X26HNON-LEAKAGE PROBABILITY = F10.7)
            IF(INF)190,190,120
120    IF(BSQ)190,130,190
130    IF(LCRIT-3)140,190,140
140    BSQ1=.001
            XK01=(1.+XLSQ*(1.-THUT)*BSQ1)*EXP(BSQ1*(B(1)*TAU(1)+B(2)*TAU(2)+
            1 B(3)*TAU(3)))
            BSQ2=.0008
            XK02=(1.+XLSQ*(1.-THUT)*BSQ2)*EXP(BSQ2*(B(1)*TAU(1)+B(2)*TAU(2)+
            1 B(3)*TAU(3)))
150    BSQ=(FK2-XK01)*(BSQ2-BSQ1)/(XK02-XK01)+BSQ1
            XK0=(1.+XLSQ*(1.-THUT)*BSQ)*EXP(BSQ*(B(1)*TAU(1)+B(2)*TAU(2)+
            1 B(3)*TAU(3)))
            IF(ABSF(FK2-XK0)-.00001)170,160,160
160    XK01=XK02
            BSQ1=BSQ2
            XK02=XK0
            BSQ2=BSQ
            GO TO 150
170    WRITE OUTPUT TAPE NOUT,180,BSQ
180    FORMAT(1H0,37HINFINITE LATTICE MATERIAL BUCKLING = F10.8)
            BSQ=0.0
190    WRITE OUTPUT TAPE NOUT,200
200    FORMAT(1H0,41X37HAVERAGE ABSORPTIONS FOR EACH ROD KIND/1H ,27X8HRO
            1D KIND,3X10HTOTAL RODS,4X17HAVG. THERMAL ABS.,3X19HAVG. RESONANCE
            2ABS.)
            DO 230 I=1,NKIND
              N=0
              SUM1=0.0
              SUM2=0.0
              DO 220 J=1,NTYPE
                IF(KIND(J)-I)220,210,220
210    N=N+NRPRT(J)
                FN=NRPRT(J)
                SUM1=SUM1+ABSRB(J)*FN
                SUM2=SUM2+RESABS(J)*FN

```

```

220 CONTINUE
    FN=N
    AVG1=SUM1/FN
    AVG2=SUM2/FN
    WRITE OUTPUT TAPE NOUT,225,I,N,AVG1,AVG2
225 FORMAT(1H ,30X I2,9X I4,11X F9.6,12X F9.6)
230 CONTINUE
    IF(LCRIT-2)239,231,239
231 I=KSERCH
    IF(NF(I))233,233,235
233 WRITE OUTPUT TAPE NOUT,234,I
234 FORMAT(1H0,45H A NEW THERMAL UTILIZATION VALUE FOR ROD KIND I2,44H
1 CANNOT BE COMPUTED DUE TO INSUFFICIENT DATA/1H ,109H THE MODERATOR
2 ABSORPTIONS ABOVE ARE IN ERROR BECAUSE OF THIS, HOWEVER THE PRODUCT
3 OF LATTICE F AND NON-LEAKAGE/1H ,23H PROBABILITY IS CORRECT.)
    GO TO 239
235 WRITE OUTPUT TAPE NOUT,236,GAM(I),I,THERU(I)
236 FORMAT(1H0,67H THE VALUE OF THERMAL UTILIZATION CORRESPONDING TO A
1 GAMMA VALUE OF F10.6,14H FOR ROD TYPE I2,4H IS F10.8)
239 IF(LCRIT)250,250,240
240 IF(IT2-2)290,250,270
250 PRINT 260
260 FORMAT(1H ,26H PROBLEM RAN THROUGH TO END)
    GO TO 290
270 PRINT 280
280 FORMAT(1H ,26H PROBLEM DID NOT RUN TO END)
290 RETURN
430 WRITE OUTPUT TAPE NOUT,440
440 FORMAT(1H0,103H A THERMAL UTILIZATION VALUE FOR A ROD KIND IS ZERO.
1 MODERATOR THERMAL ABSORPTIONS CANNOT BE SOLVED FOR./1H ,63H THE MO
2 DERATOR ABSORPTIONS WILL BE LUMPED INTO THE TOTAL LEAKAGE)
    D2OABS=0.0
    GO TO 78
    END(2,1,0,0,0)

```



```

C   MATRIX SUBROUTINES FOR HERFSY 1
C
SUBROUTINE ABMXVC(JMAX,V,A,IMAX)
  DIMENSION V(50)
  IMAX=1
  A=ABSF(V(1))
  DO 54 I=2,JMAX
    IF(A-ABSF(V(I)))53,54,54
53  A=ABSF(V(I))
    IMAX=I
54  CONTINUE
  RETURN
  END(2,1,0,0,1)

C
SUBROUTINE DGMPMX(JMAX,XS,YS,ZS)
  DIMENSION XS(50),YS(50),ZS(50,50)
  DO 1 I=1,JMAX
    DO 1 J=1,JMAX
1    ZS(I,J)=XS(I)*YS(I,J)
  RETURN
  END(2,1,0,0,1)

C
SUBROUTINE MXMPCV(JMAX,XS,YS,ZS)
  DIMENSION XS(50,50),YS(50),ZS(50)
  DO 1 I=1,JMAX
    ZS(I)=0.0
    DO 1 J=1,JMAX
1    ZS(I)=ZS(I)+XS(I,J)*YS(J)
  RETURN
  END(2,1,0,0,1)

C
SUBROUTINE MXADDG(JMAX,XS,YS,ZS)
  DIMENSION XS(50,50),YS(50),ZS(50,50)
  DO 10 J=1,JMAX
    DO 10 I=1,JMAX
      IF(I-J)11,12,11
11    ZS(I,J)=XS(I,J)
      GO TO 10
12    ZS(I,J)=XS(I,J)+YS(I)
10  CONTINUE
  RETURN
  END(2,1,0,0,1)

C
SUBROUTINE MXMPDG(JMAX,XS,YS,ZS)
  DIMENSION XS(50,50),YS(50),ZS(50,50)
  DO 1 I=1,JMAX
    DO 1 J=1,JMAX
1    ZS(I,J)=XS(I,J)*YS(J)
  RETURN
  END(2,1,0,0,1)

C
SUBROUTINE MXSBMX(JMAX,XS,YS,ZS)
  DIMENSION XS(50,50),YS(50,50),ZS(50,50)
  DO 1 J=1,JMAX
    DO 1 I=1,JMAX
1    ZS(I,J)=XS(I,J)-YS(I,J)
  RETURN
  END(2,1,0,0,1)

C

```

```

C
SUBROUTINE MXMPMX(JMAX,XS,YS,ZS)
DIMENSION XS(50,50),YS(50,50),ZS(50,50)
DO 1 I=1,JMAX
DO 1 J=1,JMAX
ZS(I,J)=0.0
DO 1 K=1,JMAX
1 ZS(I,J)=ZS(I,J)+XS(I,K)*YS(K,J)
RETURN
END(2,1,0,0,1)

```

```

C
SUBROUTINE RVMPGV(JMAX,XS,YS,ZS)
DIMENSION XS(50),YS(50)
ZS=0.0
DO 1 I=1,JMAX
1 ZS=ZS+XS(I)*YS(I)
RETURN
END(2,1,0,0,1)

```

```

C
SUBROUTINE RVMPDG(JMAX,XS,YS,ZS)
DIMENSION XS(50),YS(50),ZS(50)
DO 1 I=1,JMAX
1 ZS(I)=XS(I)*YS(I)
RETURN
END(2,1,0,0,1)

```

```

C
SUBROUTINE RVMPMX(JMAX,XS,YS,ZS)
DIMENSION XS(50),YS(50,50),ZS(50)
DO 1 J=1,JMAX
ZS(J)=0.0
DO 1 I=1,JMAX
1 ZS(J)=ZS(J)+XS(I)*YS(I,J)
RETURN
END(2,1,0,0,1)

```

APPENDIX C - IBM 704 Operating Instructions

1. ROUTINE PROBLEM OPERATION

(a) Machine Requirements

To run a routine problem on the 704 requires the following minimum equipment:

- IBM 704 with 32 k memory
- Five tape units
- Card reader
- On-Line printer

(b) Tape Units and Functions

The five tape units are dialed to numbers 1, 2, 3, 8, and 10, and are assigned the following functions:

Tape 1: Program tape (file protect on)

Tape 2: Geometry temporary storage (must always be mounted)

Tape 3: Geometry permanent storage (mounted only if used)

Tape 8: Input data in BCD card image form

Tape 10: Output of all information

(c) Card Reader Function

The card reader is used only to place data on tape 8 and to initiate the CHAIN program. This initiation is done by first preparing all the tapes in (b) that are used, then placing in the card reader the following two cards:

- Self-Loading Chain Initiator Program card (SLCI card)
- Control Card calling for Link 1 of tape 1 [C.C(1,1) CARD]

Clear and load cards. The program from this point will operate automatically unless an error stop occurs. If stop does occur refer to Section 2 of this appendix for the proper procedure.

(d) On-Line Printer Functions

The alphameric card read as the first card of each problem is printed out, and will be followed by one of four messages which give a summary of what happened to the problem. These messages are listed below with their meaning.

- AN ERROR HAS BEEN DETECTED IN THE INPUT DATA FOR THIS PROBLEM. AN ATTEMPT WILL BE MADE TO CONTINUE WITH SUCCESSIVE PROBLEMS - The error checking routines in the first link of the code have detected an error that cannot allow the problem to run to the end - no action is required of the operator, for restart has been programmed for the next problem.
- PROBLEM RAN THROUGH TO END - Any problem run to the end will have this on the on-line printer - no action is called for by operator.
- PROBLEM DID NOT RUN TO END - This indicates an error condition arose at some point of the calculations in the second link of the program. Information explaining the stop is on tape 10 - no action is called for from operator.
- MACHINE ERROR - RESTART PROBLEM - A branching error has occurred in the machine - Operator should stop machine; rewind tapes 1, 2, 3, and 8; and restart the problem from the beginning using the procedure of paragraph c above.

(e) End of Computations

The signal that the end of the problem has been reached may be one of two transfers that depend on the input data. In some cases a HPR 0,1 will occur, while in others the Card Reader will be selected. More often than otherwise the stop will select the Card Reader; however, the HPR 0,1 is equally valid under the assumption that a part of the input data was left out.

2. RESTART PROCEDURE FOR PROGRAM STOPS

If any standard BSS program stop occurs, other than the HPR 0,1 mentioned in (1e) above, the stop is considered an error stop. Certain steps may and should be taken to interrogate the status of the machine. These are:

(a) Check on-line printer

- If problem in which stop occurred is the first problem, the entire run should be cancelled after the address lights are recorded.
- If at least one problem has been completed, continue with paragraph (b).

(b) The restart procedure following case ii of (a) above is as follows:

- Rewind tape 1
- Place SLCI and C.C.(1,1) cards in card reader
- Load cards WITHOUT CLEARING MACHINE

(c) If further stops occur, the procedure of (b) above should be followed until a HPR 0,1 appears in the STORAGE REGISTER lights.

This procedure assures that problems following the one in which the stop originally occurred will not be affected by errors in previous data if at all possible.

3. PROGRAM TAPE PREPARATION

The method of preparing program tapes is contained in a Special Version of the IBM CHAIN MONITOR program written at SRL. This system is described in Appendix D, and reference to this appendix will explain the operations involved.

4. PERMANENT GEOMETRY TAPE PREPARATION

No special preparations are necessary unless the problem is the first one on the tape. If a new tape is used, the only preparation (other than mounting the tape on unit 3) is to place an END OF FILE mark on the tape in its rewind position. This may be done by placing a WEF 3 instruction into the machine manually from the console keys. The remainder of the tape processing will be controlled by the mnemonics NFILE, NWRITE, and XCOR specified by the user.

5. ADDITIONAL PROGRAM STOPS

In addition to the standard BSS LOADER program stops, the following stops have been programmed.

<u>Storage Register Address Lights</u>	<u>Error Condition</u>
HPR 464	CHECKSUM ERROR in reading geometry scratch tape
HPR 547	CHECKSUM ERROR in reading geometry permanent tape
HPR 610	Misplaced EOF on tape 3 or misread of tape 3

Pressing start allows any of the above three stops to be ignored.

HTR 5, 4	File number called for is not on tape 3
----------	--

Pressing start returns control to calling program.

APPENDIX D - Operation and Preparation of CHAIN Programs

The version of the Chain Monitor system used in the SRL - HERESY I code was written by J. E. Suich of SRL. This system consists of a package of four routines which permit operation on the IBM 704 of programs written within the specifications of the IBM-MONITOR routine, CHAIN. These four routines are described below.

(1) SUBROUTINE CHAIN (LINK, IOUNIT)

This subroutine performs the function of the IBM-FORTRAN MONITOR subroutine of the same name (see, e.g. Reference Manual, 709/7090 FORTRAN (IBM-C28-6054-2)). LINK is the fixed point integer identification number of the next chain link desired, and IOUNIT is the logical tape designation (fixed point integer 1-10) of the tape on which this link is stored. Links may be called in any order from any number of tapes. Execution of the calling sequence causes the desired link to be brought into memory, and control to be passed to it, without destroying COMMON storage.

Termination of any program or the initiation of another CHAIN job is accomplished by calling a tape designation that does not exist (i.e. >10). Execution of this command will initiate a card load sequence which will in turn may cause the SELF-LOADING CHAIN INITIATOR described below to be executed. By this process a sequence of CHAIN programs may be continuously run.

Only one error stop has been programmed: HTR 3,4 indicates that the desired LINK was not found on the designated IOUNIT. Pressing start causes control to return to the calling program.

(2) SELF-LOADING CHAIN INITIATOR (SLCI)

To provide a convenient method of beginning the chain job from a previously prepared chain program tape, a self-loader has been written which reads a single Control Card designating the desired initial values of LINK and IOUNIT. The Control Card format is row binary: LINK in the address field of the 9L row, and IOUNIT in the address field of the 9R row. For example, the Control Card to enter the link with identification number 5 on the chain program tape mounted on tape 9 has 9's punched in columns 34, 36, 69, and 72. The control cards may be punched with any additional characters (except I, R, Z, and 9) in any columns (for identification purposes).

(3) CHAIN PROGRAM TAPE GENERATOR (CPTG)

This routine fulfills the IBM-MONITOR function of preparing chain program tapes, although in substandard fashion. To use this program tape generator, the main program of each link must be converted to a subroutine named MAIN, either by compiling with the header card SUBROUTINE MAIN, or by punching 7's in columns 1, 4, 8, 12, 14, 15, 18, 19, 22, 24, 25, 26, 31, and 32 of the program card for the main program of each link and correcting the check sums. The various links are then stacked in the card hopper, tapes readied, and CLEAR, LOAD CARDS pressed. Figure D-1 shows schematically the deck makeup.

(4) END PROGRAM TAPE CARD (EPTC)

The program tape generated by the above procedure requires an END of FILE mark following the last link. This self-loader reads a Control Card (see discussion under (2) above), ignores the LINK value, and ends the tape designated by IOUNIT. The card should be placed behind the transfer card of the last link when making up a program tape (Figure D-1). There should be an END TAPE loader and Control Card pair for each IOUNIT designated.

FIGURE D-1

EXAMPLE OF THE DECK MAKEUP TO WRITE LINKS
1, 2, AND 5 ON TAPE 1 AND 1, 4, AND 3 ON TAPE 9

BSS LOADER
MAIN
SUBROUTINES (INCLUDING CHAIN)
CPTG
TRANSFER CARD
CONTROL CARD (1,1)

BSS
MAIN
SUBR
CPTG
TRA
C.C. (1,9)

.
:
C.C. (2,1)

:
C.C. (5,1)

:
C.C. (4,9)

:
C.C. (3,9)

EPTC
C.C. (0,1)
EPTC
C.C. (0,9)

Listings of the Four Subroutines

```

REM SUBROUTINE
FUL
ORG 0
MZE *,,ENTRY-2
PZE
PZE BREAK,,0
PZE 32562
BCD 1CHAIN
PZE EP
ENTRY EQU *
ORG 0
REL
EP SXD EOF,2
TXI **1,2,1
ID CLA 1,4
STA **1
CLA **
ARS 18
STD ID
IO CLA 2,4
STA **1
CLA **
ARS 18
ADM BASE
STA RDID
STA REWTP
STA TEMP
CLA TEMP
SUB OUT
TMI RDID
TRA CREAD
RDID RTB 1
CPY 16
TRA **3
EOF TIX REWTP,2,0
HTR 3,4
CLA ID
SUB 16
TNZ RDID
BASE PXD 144,0
CAD 16
CAD 17
LXA 16,4
TIX 16,4,18
REWTP REW 1
TRA RDID
CREAD RCD
CPY 0
CPY 1
TRA 0
OUT PZE 155
TEMP PZE 0
BREAK EQU *
END **

```

CHAIN (ID,IO)

P	C	G
R	A	E
O	R	N
G	D	E
R		R
A		A
M		T
		O
		R

DESIRED LINK IDENT

LOGICAL TAPE DESIGNATION

TAPE UNIT ADDRESS

TAPE LINK IDENT

EOF, REWIND ONCE
LINK NOT ON TAPE, RETURN

NOT THIS LINK
LINK IS FOUND
LOAD
TAPE
SEQUENCE

REM ROUTINE TO MAKE A CHAIN PROGRAM TAPE					
	FUL				G
	ORG 0		P		E
	MZE *,,ENTRY-2		R	C	N
	PZE		Q	A	E
	PZE BREAK,,1		G	R	R
	PZE 32562		R	D	A
	BCD 1000000		A		T
	PZE BP+1		M		O
					R
ENTRY	EQU *				
	ORG 0				
	REL				
EP	BCD 1MAIN				
	CLA *				
	SUB ONE				
	STA LOAD				
	STA WTPGM+5				
	CLA EP				
	STA LOAD+6				
	STA LOAD+7				
	RCD				
	CPY EP				
	CPY EP+1				
	CLA EP+1				
	ADM BASE				
	STA WTPGM				
	LXA EIGHT,4				
	CLA LOAD+8,4				
	STO 24,4				
	TIX *-2,4,1				
WTPGM	WTB 1				
BASE	PXD 144,0				
	LXA **3,4				
	CPY EP				
	TIX *+1,4,16				
	CAD **,4				
	TIX *-1,4,1				
	STO EP				
	CPY EP				
	RCD				
	CPY 0				
	CPY 1				
	TRA 0				
LOAD	CAD **,4				
	TIX 16,4,1				
	STO 1				
	CPY 0				
	CLA 0				
	SUB 1				
	TZE **				
	HTR **				
ONE	PZE 1				
EIGHT	PZE 8				
BREAK	EQU *				
	END **				

MAIN PGM ENTRY POINT
PROGRAM BREAK +1
ESTABLISH
ABSOLUTE
BINARY
PROGRAM
LOADER
READ CONTROL CARD
LINK IDENT
TAPE DESIGNATION
FORM
TAPE ADDRESS
SHIFT TAPE LOADER
WRITE BINARY TAPE
LINK IDENT
PROGRAM
CHECKSUM
LOAD NEXT LINK

REM SELF LOADING CHAIN INITIATOR	
FUL	
ORG 0	SELF
LXD **,4	LOAD
CPY 22274,4	SEQUENCE
TIX *-1,4,1	EOF
HTR **	CONTROL CARD
RCD	LINK
CPY 0	TAPE UNIT
CPY 1	
CLA 1	
ADM BASE	
STA RDTP	
RDTP RTB 1	TAPE ID
CPY 1	
TRA **3	TAPE NOT REWOUND
HTR RDTP	OR WRONG TAPE
HTR RDTP	COMPARE
CLA 0	IDENTS
SUB 1	WRONG LINK
TNZ RDTP	RIGHT LINK
BASE PXD 144,0	LOAD
CAD 16	TAPE
CAD 17	SEQUENCE
LXA 16,4	
TIX 16,4,18	
END **	

REM SBLF LOADER TO END CHAIN TAPE	
FUL	
ORG 0	SELF
LXD **,4	LOAD
CPY 22274,4	SEQ
TIX *-1,4,1	
HTR **	CONTROL CARD
RCD	IGNORE
CPY 0	TAPE UNIT - LOGICAL
CPY 1	
CLA 1	- ADDRESS
ADM BASE	
STA WEFTP	
STA REWTP	
WEFTP WEF 1	END FILE
REWTP REW 1	REWIND
RCD	LOAD
CPY 0	NEXT
CPY 1	END
TRA 0	FILE
BASE PXD 144,0	
END **	