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A Method for Calculating the Steady-State Distribution of Tritium in a Molten-Salt Breeder Reactor Plant

R. B. Briggs
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Printed in the United States of America. Available from
National Technical Information Service
U.S. Department of Commerce
5285 Port Royal Road, Springfield, Virginia 22161
Price: Printed Copy \$5.45; Microfiche \$2.25

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ORNL-TM-4804
UC-76 - Molten Salt
Reactor Technology

Contract No. W-7405-eng-26

Reactor Division

A METHOD FOR CALCULATING THE STEADY-STATE DISTRIBUTION
OF TRITIUM IN A MOLTEN-SALT BREEDER REACTOR PLANT

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APRIL 1975

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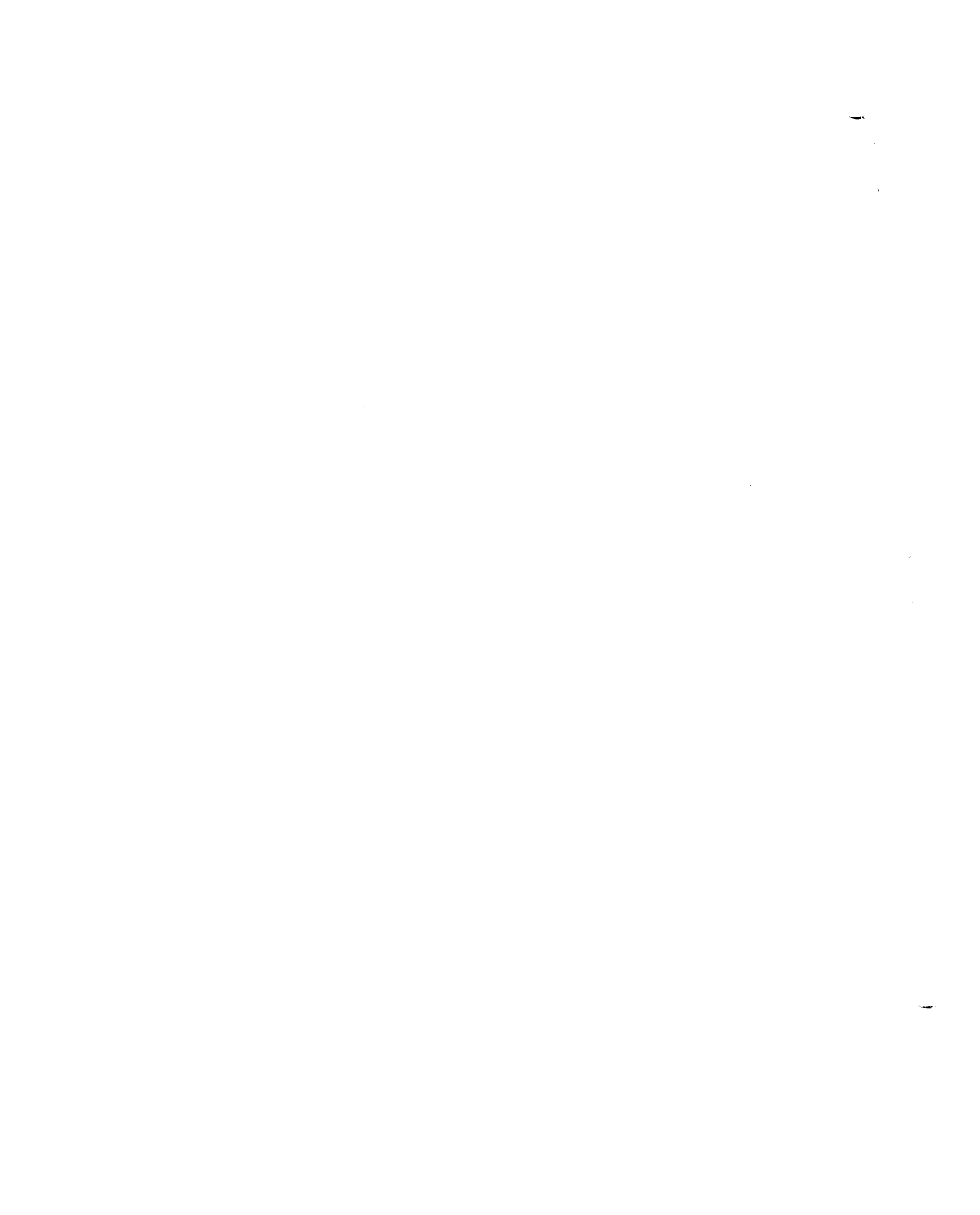
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R. B. Briggs and C. W. Nestor, Jr.

ABSTRACT

Tritium is produced in molten salt reactors primarily by fissioning of uranium and absorption of neutrons by the constituents of the fuel carrier salt. At the operating temperature of a large power reactor, tritium is expected to diffuse from the primary system through pipe and vessel walls to the surroundings and through heat exchanger tubes into the secondary system which contains a coolant salt. Some tritium will pass from the secondary system into the steam power system. This report describes a method for calculating the steady state distribution of tritium in a molten salt reactor plant and a computer program for making the calculations. The method takes into account the effects of various processes for removing tritium, the addition of hydrogen or hydrogenous compounds to the primary and secondary systems, and the chemistry of uranium in the fuel salt. Sample calculations indicate that 30 percent or more of the tritium might reach the steam system in a large power reactor unless special measures are taken to confine the tritium.

I. INTRODUCTION

Conceptual designs of Molten Salt Breeder Reactor (MSBR) power plants usually can be represented by the diagram shown in Fig. 1. The fissioning of uranium in the fuel salt heats the salt as it is pumped through the reactor vessel in the primary system. The heat is transferred to a coolant salt that circulates in the secondary system and, thence, to water, producing steam to drive a turbine-generator in the steam system.

Fission products and other radioactive materials are produced in large amounts in the fuel salt. Much smaller amounts are produced in the coolant salt by the flux of delayed neutrons in the primary heat exchangers. The radioactivity is normally confined by the walls of the piping and vessels. However, tritium is produced in the salts, partly as a fission product, but mostly by absorption of neutrons by lithium in the fuel salt. At the high temperature of an MSBR, tritium diffuses through metals and might escape to the environs in amounts that would be cause for concern.

The purpose of this report is to describe a method for calculating the distribution of tritium in and its escape from an MSBR plant. We assume that the tritium, born as tritium ions, is present in the fuel salt primarily as tritium molecules* and tritium fluoride molecules.** The ions are estimated to be produced at a rate of $2.6 \times 10^{14}/\text{MWsec}$ ***

*Tritium molecules are intended to include HT and H₂ molecules when hydrogen is present.

**Tritium fluoride molecules are intended to include tritium (and hydrogen) ions associated with fluoride ions in the salt.

***2420 Ci/day in a 2250 MW(t), 1000 MW(e) plant.

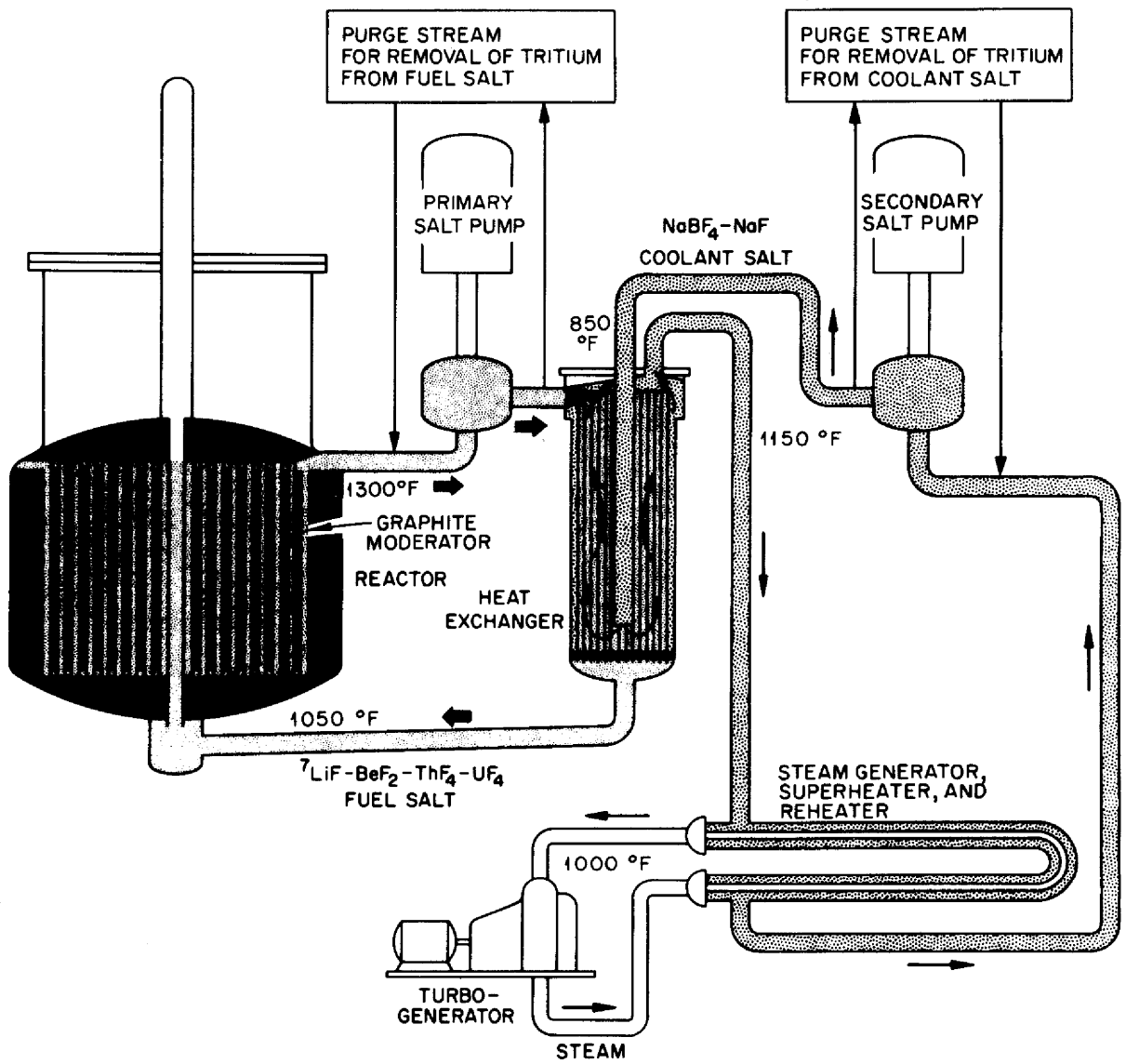
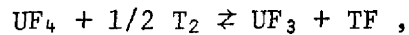


Fig. 1. Molten Salt Breeder Reactor System.

in a typical fuel salt. The relative concentrations of tritium and tritium fluoride in the fuel salt are expected to be governed by the equilibrium relationship for the reaction,



with uranium in the salt. The absolute concentrations are governed by removal processes.

Three types of processes are provided for removing tritium from the primary system: permeation through the metal of the walls of piping and vessels, sorption on materials in contact with the salt, and purging. We assume that tritium molecules that reach a metal surface can sorb on the surface, dissociate into tritium atoms and diffuse through the metal. Tritium in tritium fluoride and other compounds is assumed to be chemically bound and unable to pass through the metal.

Experience with the Molten Salt Reactor Experiment indicated that tritium sorbs on and is tightly bound to graphite. We provide for sorption of tritium and tritium fluoride on the graphite in the reactor core.

Provision is made for purging tritium from the primary system by circulating a stream of salt through an apparatus which extracts gaseous tritium and tritium compounds. A contactor in which tritium and tritium fluoride are transferred to a gas phase by virtue of their vapor pressures would be such an apparatus. Current designs for MSBR's provide for sparging of the fuel salt with helium bubbles in the primary system to remove krypton and xenon. Tritium and tritium fluoride would be removed also. The sparging process can be treated as an equivalent purging process in the calculations.

Tritium will reach the secondary system by diffusion from the primary system through the walls of the tubes in the primary heat exchangers and by neutron capture in the coolant salt. We provide for removal of tritium from the secondary system by diffusion through the metal walls, sorption, and purging. The secondary system would not normally contain a sorber or have an elaborate purging system. Such processes, if incorporated into the plant, would be designed specifically for removing tritium.

The coolant salts do not normally contain constituents that are reducible by tritium and, thereby, able to convert tritium into tritium fluoride and make it unavailable to diffuse through the metal walls. We, therefore, have provided for addition of hydrogen fluoride or other hydrogenous compounds to the secondary system. We assume that tritium will exchange with the hydrogen in the added compound and that the compound will be extracted by the sorption and/or purge process.

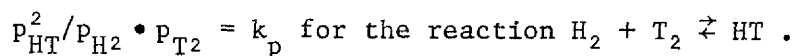
The steam system and the cells around the reactor primary and secondary systems are considered to be sinks for tritium. Tritium reaching the steam system is assumed to exchange with hydrogen in the water, and that reaching the cells is assumed to be oxidized to water. The partial pressure of tritium is effectively zero.

In the calculations we assume that tritium and hydrogen behave identically. The equation used for calculating the diffusion of hydrogen through a metal wall states that the rate of transport per unit of surface area is proportional to the product of a permeability coefficient and the difference between the square roots of the partial pressures of hydrogen at the inner and outer surfaces of the metal.

In this circumstance, addition of hydrogen can reduce the transport of tritium through the metal. Suppose, for example, the partial pressures of tritium and hydrogen at the outer surface of a pipe are zero and the partial pressure of tritium at the inner surface is held constant. If hydrogen were added to increase the total hydrogen partial pressure at the inner surface by a factor of 100, the flow of hydrogen plus tritium through the metal wall would increase by a factor of 10. But the flow of tritium would decrease by a factor of 10 because of the 100-fold dilution of hydrogen. Because of other factors, the effect of adding hydrogen may not be so dramatic, but the calculational method provides for addition of hydrogen to the primary and secondary systems and for hydrogen to be present at a specified concentration in the steam system so that the effects can be studied.*

The calculational model describes the behavior of tritium in an MSBR plant to the extent that it is known or has been inferred at the present time. The removal processes can be included in or eliminated from the calculations by careful choice of the values assigned to coefficients in the equations. The model probably does not include all the chemical reactions and physical processes that will ultimately be

*The calculational procedure might have been developed to treat hydrogen and tritium as separate species. Separate values then could be assigned to important parameters, such as solubility and diffusion coefficients, for each species. Interaction between hydrogen and tritium would be taken into account by the equilibrium relationship

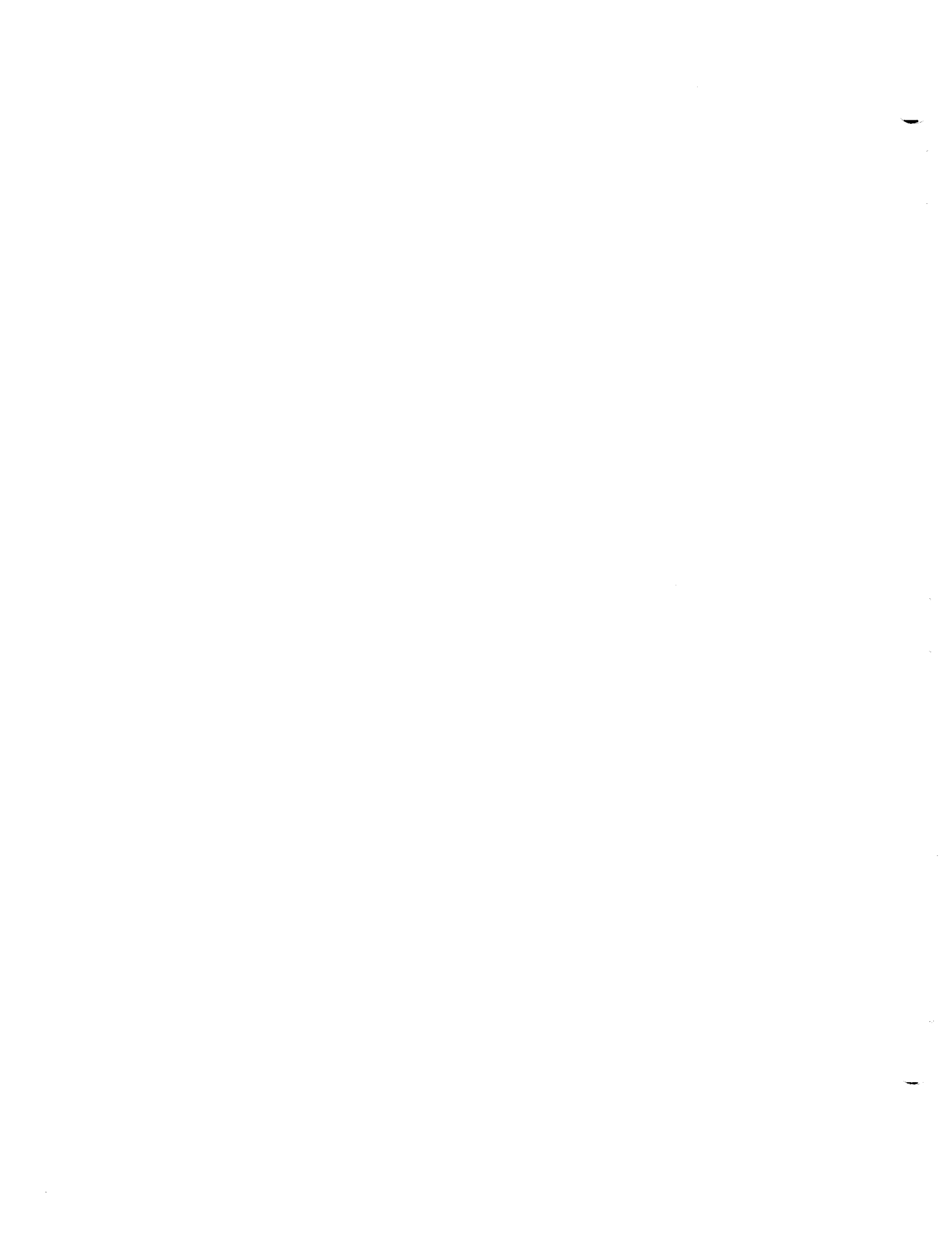


However, k_p has a value near 4 at temperatures of interest, which signifies that hydrogen and tritium interact as though they are the same species. Also, there are substantial uncertainties in the values for most of the parameters. Complicating the procedure to treat hydrogen and tritium separately would not, for the present, improve the accuracy of the results.

shown to affect the distribution of tritium in an MSBR. In some instances these effects can be included, when recognized, simply by adjusting the coefficients in equations for processes presently included. Others may require incorporation of additional processes.

Two assumptions in the calculational procedure should be recognized for their potential for leading to major differences between the calculated distribution of tritium and what would actually occur in a reactor plant. Tritium, present in the salt as tritium fluoride, can react with metal to yield tritium atoms that would dissolve in and diffuse through the metal. Neglect of this reaction could cause the calculations to be greatly in error under circumstances where most of the tritium is present in the salt as tritium fluoride.

Oxide films (and possibly others) that form on metal surfaces reduce the permeability of a metal wall to the passage of hydrogen. They may also cause the transport to vary with pressure to a power in the range of $1/2$ to 1. The reduced permeability appears as a coefficient in the transport equations of the model, but we make no provision for changing the exponent on the pressure terms from $1/2$. The calculated transport of tritium through the metal walls and the effect of the addition of hydrogen in reducing the transport would both be greater than would actually occur if the actual transport were proportional to the pressure to a power in the range $1/2$ to 1. The calculations would not underestimate the transport unless the total pressure of tritium and hydrogen exceeded the reference pressure for the permeability coefficient, which is usually 1 atm.



II. DERIVATION OF EQUATIONS AND COMPUTATIONAL PROCEDURES

In making the calculations, we first calculate the distribution of hydrogen plus tritium in order to establish flows and concentrations of the combined isotopes throughout the plant. Then we calculate the distribution of tritium throughout the plant.

For calculating the distribution, the fluids in the primary and secondary systems and the various parts of the steam system are assumed to be well mixed and to contain uniform bulk concentrations of all constituents. The calculations are for steady-state conditions, and only hydrogen and tritium molecules are assumed to be able to sorb on the metal surfaces, dissociate, and diffuse through the metal walls. The various paths are defined and the distribution is calculated by the use of the following set of equations.*

A. In the primary system:

1. Transport of hydrogen through the salt film to the wall of the piping in the hot leg from the reactor vessel to the heat exchanger:

$$Q_1 = h_1 A_1 (C_F - C_1) . \quad (1a)$$

Transport through the pipe wall to the surroundings where the hydrogen pressure is assumed to be negligible:

$$Q_1 = \frac{p_1 A_1 [(k_1 C_1)^{\frac{1}{2}} - 0]}{t_1} = \frac{p_1 A_1 (k_1 C_1)^{\frac{1}{2}}}{t_1} . \quad (1b)$$

2. Transport of hydrogen to and through the walls of the cold-leg piping from the heat exchanger to the reactor vessel:

*Symbols are defined in Section IV, Nomenclature.

$$Q_2 = h_2 A_2 (C_F - C_2) \quad (2a)$$

$$= \frac{p_2 A_2 (k_2 C_2)^{\frac{1}{2}}}{t_2} . \quad (2b)$$

3. Transport of hydrogen to and through the walls of the reactor vessel and the shells of the heat exchangers in the primary system:

$$Q_3 = h_3 A_3 (C_F - C_3) \quad (3a)$$

$$= \frac{p_3 A_3 (k_3 C_3)^{\frac{1}{2}}}{t_3} \quad (3b)$$

4. Transport of hydrogen to and through the walls of the tubes in the primary heat exchangers into the secondary system:

$$Q_4 = h_4 A_4 (C_F - C_4) \quad (4a)$$

$$= \frac{p_4 A_4}{t_4} \left[(k_4 C_4)^{\frac{1}{2}} - (k_{12} C_{12})^{\frac{1}{2}} \right] . \quad (4b)$$

5. Transport of hydrogen to the surfaces of the graphite in the reactor vessel or to other sorber:

$$Q_5 = h_5 A_5 (C_F - C_5) . \quad (5a)$$

Sorption by the graphite or other sorber assuming that the sorbing surface is replaced continuously and that the concentration of sorbed gas is proportional to the square root of the partial pressure:

$$Q_5 = B_1 W_1 A_5 (k_5 C_5)^{\frac{1}{2}} . \quad (5b)$$

6. Removal of hydrogen by purge:

$$Q_6 = F_1 E_1 C_F . \quad (6)$$

7. Transport of hydrogen fluoride to and removal by sorber:

$$Q_7 = h_7 A_7 (C_{FF} - C_7) \quad (7a)$$

$$= B_2 W_2 A_7 (k_7 C_7)^{\frac{1}{2}} . \quad (7b)$$

8. Removal of hydrogen fluoride by purge:

$$Q_8 = F_2 E_2 C_{FF} . \quad (8)$$

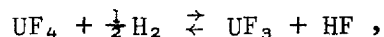
Because the molecular species involved may contain different numbers of hydrogen atoms, all the calculations are done in terms of atoms of hydrogen. This does not mean that the hydrogen necessarily diffuses as single atoms, but only that a transport unit is one hydrogen atom and the parameters are expressed in terms of single hydrogen atoms. A Q value of 1 then represents the transport of one-half molecule of H_2 , one molecule of HF, or one-fourth molecule of a compound like CH_4 , all per unit time. Likewise, a C value of 1 represents a concentration of one-half molecule of H_2 , one molecule of HF, or one-fourth molecule of CH_4 , all per unit volume.

If the rates of inflow of tritium and hydrogen atoms (R_1 and R_2 , respectively) to the primary system are given, a material balance over the primary system gives

$$R_1 + R_2 = \sum_{i=1}^8 Q_i . \quad (9)$$

In our calculations, all flow rates in the sum on the right-hand side of Eq. 9 are positive or zero except for Q_4 , the transport through the

heat exchanger tubes to the secondary system. Q_4 can be positive, negative or zero, depending on the conditions in the various systems. Hydrogen is present in and is removed from the primary system as hydrogen fluoride, but we provide no input of HF. It is produced by the reaction



which has an equilibrium quotient

$$\frac{X(\text{UF}_3)}{X(\text{UF}_4)} \times \frac{P(\text{HF})}{[P(\text{H}_2)]^{\frac{1}{2}}} = M' ,$$

or

$$\frac{X(\text{UF}_3)}{X(\text{UF}_4)} \times \frac{k_7 C_{\text{FF}}}{(k_5 C_{\text{F}})^{\frac{1}{2}}} = M .$$

Corrosion and other chemical considerations make it desirable to maintain the ratio $X(\text{UF}_3)/X(\text{UF}_4) \equiv 1/U$ at a constant value,* so the concentration of HF in the bulk of the salt can be related to the hydrogen concentration by

$$C_{\text{FF}} \approx \frac{MU}{k_7} (k_5 C_{\text{F}})^{\frac{1}{2}} . \quad (10)$$

We replace C_{FF} by the equivalent function of C_{F} in Eqs. 7a and 8 to obtain expressions for Q_7 and Q_8 in terms of C_{F} .

B. Secondary System:

1. Hot-leg piping:

$$Q_{10} = h_{10} A_{10} (C_{\text{C}} - C_{10}) \quad (11a)$$

$$= \frac{P_{10} A_{10}}{t_{10}} (k_{10} C_{10})^{\frac{1}{2}} . \quad (11b)$$

*This might require that hydrogen be added to the primary systems as a mixture of hydrogen and hydrogen fluoride.

2. Cold-leg piping:

$$Q_{11} = h_{11}A_{11} (C_C - C_{11}) \quad (12a)$$

$$= \frac{P_{11}A_{11}}{t_{11}} (k_{11}C_{11})^{\frac{1}{2}} . \quad (12b)$$

3. Transport through the primary heat exchanger tubes into the primary system:

$$Q_{12} = h_{12}A_4 (C_C - C_{12}) \quad (13a)$$

$$= \frac{P_4A_4}{t_4} \left[(k_{12}C_{12})^{\frac{1}{2}} - (k_4C_4)^{\frac{1}{2}} \right] . \quad (13b)$$

4. Transport through the steam generator tubes into the steam system:

$$Q_{13} = h_{13}A_{13} (C_C - C_{13}) \quad (14a)$$

$$= \frac{P_{13}A_{13}}{t_{13}} \left[(k_{13}C_{13})^{\frac{1}{2}} - (k_{21}C_{21})^{\frac{1}{2}} \right] . \quad (14b)$$

5. Transport through the superheater tubes into the steam system:

$$Q_{14} = h_{14}A_{14} (C_C - C_{14}) \quad (15a)$$

$$= \frac{P_{14}A_{14}}{t_{14}} \left[(k_{14}C_{14})^{\frac{1}{2}} - (k_{22}C_{22})^{\frac{1}{2}} \right] . \quad (15b)$$

6. Transport through the reheater tubes into the steam system:

$$Q_{15} = h_{15}A_{15} (C_C - C_{15}) \quad (16a)$$

$$= \frac{P_{15}A_{15}}{t_{15}} \left[(k_{15}C_{15})^{\frac{1}{2}} - (k_{23}C_{23})^{\frac{1}{2}} \right] . \quad (16b)$$

7. Removal by sorber as hydrogen:

$$Q_{16} = h_{16}A_{16} (C_C - C_{16}) \quad (17a)$$

$$= B_3W_3A_{16} (k_{16}C_{16})^{\frac{1}{2}} . \quad (17b)$$

8. Removal by purge as hydrogen:

$$Q_{17} = F_3 E_3 C_C \quad . \quad (18)$$

9. Removal by sorber as HF:

$$Q_{18} = h_{18} A_{18} (C_{CF} - C_{18}) \quad (19a)$$

$$= B_4 W_4 A_{18} (k_{18} C_{18})^{\frac{1}{2}} \quad . \quad (19b)$$

10. Removal by purge as HF:

$$Q_{19} = F_4 E_4 C_{CF} \quad . \quad (20)$$

Since we assume that the hydrogen fluoride does not release hydrogen to diffuse through the metal walls, and that there are no chemical reactions in the secondary system that make the concentrations of hydrogen and hydrogen fluoride interdependent, we write separate material balances for the two species for the distribution of total tritium and hydrogen:

$$R_3 + R_4 = \sum_{i=10}^{17} Q_i \quad (21a)$$

$$R_5 = Q_{18} + Q_{19} \quad . \quad (21b)$$

In these equations all the R's and all the Q's have positive or zero values except for Q_{12} , Q_{13} , Q_{14} and Q_{15} , which can have negative values.

C. Steam generator system:

1. Transport through the steam generator tubes into the secondary system:

$$Q_{21} = h_{21} A_{13} (C_{SG} - C_{21}) \quad (22a)$$

$$= \frac{P_{13} A_{13}}{t_{13}} \left[(k_{21} C_{21})^{\frac{1}{2}} - (k_{13} C_{13})^{\frac{1}{2}} \right] \quad . \quad (22b)$$

2. Transport through superheater tubes into the secondary system:

$$Q_{22} = h_{22} A_{14} (C_{SS} - C_{22}) \quad (23a)$$

$$= \frac{P_{14}A_{14}}{t_{14}} \left[(k_{22}C_{22})^{\frac{1}{2}} - (k_{14}C_{14})^{\frac{1}{2}} \right]. \quad (23b)$$

3. Transport through the reheater tubes into the secondary system:

$$Q_{23} = h_{23}A_{15}(C_{SR} - C_{23}) \quad (24a)$$

$$= \frac{P_{15}A_{15}}{t_{15}} \left[(k_{23}C_{23})^{\frac{1}{2}} - (k_{15}C_{15})^{\frac{1}{2}} \right]. \quad (24b)$$

In the steam system the values for C_{SG} , C_{SS} and C_{SR} will be given. The steam flows will be so large that the diffusion of hydrogen through the metals should not have much effect on the concentration of hydrogen in the steam. Under these assumptions, we do not require a material balance over the steam system. If hydrogen is added to the feed water as hydrazine or in some other manner to give a specified ratio of hydrogen to H_2O , then this ratio, coupled with the steam tables, can be used to calculate the hydrogen concentrations in the water and steam in the steam-raising equipment. Without addition of hydrogen the concentrations are established by the dissociation of water.

We now need to solve the above equations to obtain values for all the flow rates and concentrations. We carry this out in the following sequence, discussed in more detail in Sec. III.

1. Calculate C_{CF} , C_{18} , Q_{18} and Q_{19} from equations 19a, 19b, 20 and 21b.
2. Assume a value for C_C .
3. Calculate Q_{10} , Q_{11} , Q_{16} , Q_{17} and C_{16} from equations 11a, 11b, 12a, 12b, 17a, 17b and 18.
4. Calculate Q_{13} , Q_{14} , Q_{15} , C_{13} , C_{14} and C_{15} from equations 14a, 14b, 15a, 15b, 16a, 16b, 22a, 22b, 23a, 23b, 24a and 24b, noting that the steam system and the secondary system are coupled by the relationships $Q_{13} = -Q_{21}$, $Q_{14} = -Q_{22}$ and $Q_{15} = -Q_{23}$.

5. Calculate Q_{12} from the material balance, Eq. 21a.
6. Calculate C_F , C_{12} and C_4 from Eqs. 4a, 4b, 13a, 13b, the relationship $Q_4 = -Q_{12}$ and the value of Q_{12} obtained in step 5. These concentrations should all be positive. If any one of them is negative, steps 3 through 6 must be repeated with a larger value of C_C .
7. When positive values have been found for C_F , C_{12} and C_4 , calculate Q_1 , Q_2 , Q_3 , Q_5 , Q_6 , Q_7 , Q_8 , C_5 , C_{FF} and C_7 .
8. Calculate R_F from

$$R_F = \sum_{i=1}^8 Q_i - (R_1 + R_2) .$$

If R_F is positive, hydrogen must be added to the primary system in order to maintain a balance. This means that C_F is too large, which in turn means that C_C is too large, and steps 3 through 8 must be repeated with a smaller value of C_C . If R_F is negative, C_C is too small and steps 3 through 8 must be repeated with a larger value of C_C .

When this process has been repeated until the ratio $\left| \frac{R_F}{R_1 + R_2} \right|$ is sufficiently small, the flows and concentrations of hydrogen plus tritium and of hydrogen fluoride plus tritium fluoride have been established throughout the plant and we can proceed with the calculation of the tritium distribution. We ignore the difference in the properties of the two isotopes and assume that they behave identically. Thus, hydrogen and tritium compounds have the same solubilities and diffusivities, and if a hydrogenous compound, such as HF, is added to a mixture of hydrogen and tritium, exchange will occur to give a ratio of tritium to hydrogen that is the same in hydrogen* and the added compound.

*H₂, HT and T₂.

We now proceed with the calculation of the tritium distribution.

D. Primary system:

1. Transport through walls of hot-leg piping:

$$Q_{31} = \frac{C_{FT}}{C_F} Q_1 . \quad (25)$$

2. Transport through walls of cold-leg piping:

$$Q_{32} = \frac{C_{FT}}{C_F} Q_2 . \quad (26)$$

3. Transport through wall of reactor vessel and shells of heat exchangers in primary system:

$$Q_{33} = \frac{C_{FT}}{C_F} Q_3 . \quad (27)$$

4. Transport through walls of primary heat-exchanger tubes into the secondary system:

$$Q_{34} = h_4 A_4 (C_{FT} - C_{34}) \quad (28a)$$

$$= \frac{p_4 A_4}{t_4} \left[\frac{k_4 C_{34}}{(k_4 C_4)^{\frac{1}{2}}} - \frac{k_{12} C_{42}}{(k_{12} C_{12})^{\frac{1}{2}}} \right] . \quad (28b)$$

Equations 25 through 27 are straightforward, simply indicating that the amount of tritium flowing with hydrogen is proportional to the fraction of the concentration that is tritium when the flow of both is into a sink with a zero concentration of both. Equation 28a is straightforward, indicating that the flow of tritium from the bulk salt to the wall is proportional to the difference between the concentrations of tritium in the bulk fluid and the wall. Equation 28b, however, requires some additional explanation.

The rate of transport of hydrogen through a metal wall can be expressed as

$$Q = \frac{DA}{t} (C'_I - C'_O) ,$$

where D is the diffusivity of hydrogen atoms in the metal, the C's are the concentrations of hydrogen atoms dissolved in the metal at the inner (I) and outer (O) surfaces, t is the metal thickness and A is the surface area. Assuming no interaction of tritium and hydrogen atoms as they diffuse through the metal, the rate of transport of tritium is

$$Q_T = \frac{DA}{t} (C'_{TI} - C'_{TO}) .$$

The concentration of hydrogen + tritium atoms in the metal at the surface is

$$C' = SP^{\frac{1}{2}} = S(kC)^{\frac{1}{2}} ,$$

where S is a solubility coefficient and P is the partial pressure of hydrogen + tritium and is equal to the product of Henry's law coefficient and the concentration of hydrogen + tritium in the salt at the surface. Assuming that the ratio of tritium to hydrogen + tritium in the metal at the surface is the same as that in the salt at the surface, we can write

$$C'_{TI} = C'_T \frac{C_{TI}}{C_I} = S(k_I C_I)^{\frac{1}{2}} \frac{C_{TI}}{C_I} = S \frac{k_I C_{TI}}{(k_I C_I)^{\frac{1}{2}}}$$

and a similar expression for the outer surface. Then,

$$Q_T = \frac{DSA}{t} \left[\frac{k_I C_{TI}}{(k_I C_I)^{\frac{1}{2}}} - \frac{k_O C_{TO}}{(k_O C_O)^{\frac{1}{2}}} \right] ,$$

and by substituting the permeability coefficient, p , for the product, DS, we obtain Eq. 28b. This treatment is necessary here because the net flows of hydrogen and tritium may be in opposite directions. The equations provide a means for taking into account the effect of the mass action laws on the concentrations of tritium in the metal and its transport through the metal.

5. Removal by graphite or other sorber:

$$Q_{35} = \frac{C_{FT}}{C_F} Q_5 \cdot \quad (29)$$

6. Removal by purge:

$$Q_{36} = \frac{C_{FT}}{C_F} Q_6 \cdot \quad (30)$$

7. Removal by graphite or other sorber as tritium fluoride:

$$Q_{37} = \frac{C_{FT}}{C_F} Q_7 \cdot \quad (31)$$

8. Removal by purge as tritium fluoride:

$$Q_{38} = \frac{C_{FT}}{C_F} Q_8 \cdot \quad (32)$$

The tritium balance over the primary system is:

$$R_1 = \sum_{i=31}^{38} Q_i \cdot \quad (33)$$

E. Secondary system:

1. Hot-leg piping:

$$Q_{40} = \frac{C_{CT}}{C_C} Q_{10} \cdot \quad (34)$$

2. Cold-leg piping:

$$Q_{41} = \frac{C_{CT}}{C_C} Q_{11} \quad (35)$$

3. Transport through primary heat exchanger tube walls into primary system:

$$Q_{42} = h_{12}A_4(C_{CT} - C_{42}) \quad (36a)$$

$$= \frac{p_4 A_4}{t_4} \left[\frac{k_{12} C_{42}}{(k_{12} C_{12})^{1/2}} - \frac{k_4 C_{34}}{(k_4 C_4)^{1/2}} \right] \quad (36b)$$

4. Transport through steam generator tube walls into the steam system:

$$Q_{43} = h_{13}A_{13}(C_{CT} - C_{43}) \quad (37a)$$

$$= \frac{p_{13} A_{13}}{t_{13}} \frac{k_{13} C_{43}}{(k_{13} C_{13})^{1/2}} \quad (37b)$$

Calculations of the tritium distribution are based on the assumption that tritium will exchange so rapidly with the hydrogen in the steam to form tritiated water that the tritium concentration will be effectively zero.

5. Transport through the superheater tubes into the steam system:

$$Q_{44} = h_{14}A_{14}(C_{CT} - C_{44}) \quad (38a)$$

$$= \frac{p_{14} A_{14}}{t_{14}} \frac{k_{14} C_{44}}{(k_{14} C_{14})^{1/2}} \quad (38b)$$

6. Transport through the reheater tubes into the steam system:

$$Q_{45} = h_{15}A_{15}(C_{CT} - C_{45}) \quad (39a)$$

$$= \frac{p_{15} A_{15}}{t_{15}} \frac{k_{15} C_{45}}{(k_{15} C_{15})^{1/2}} \quad (39b)$$

7. Removal by sorber as tritium:

$$Q_{46} = \frac{C_{CT}}{C_C} Q_{16} \cdot \quad (40)$$

8. Removal by purge as tritium:

$$Q_{47} = \frac{C_{CT}}{C_C} Q_{17} \cdot \quad (41)$$

9. Removal by sorber as tritium fluoride:

$$Q_{48} = \frac{C_{CT}}{C_C} Q_{18} \cdot \quad (42)$$

10. Removal by purge as tritium fluoride:

$$Q_{49} = \frac{C_{CT}}{C_C} Q_{19} \cdot \quad (43)$$

The balance over the secondary system is:

$$R_3 = \sum_{i=40}^{49} Q_i \cdot \quad (44)$$

Since the tritium concentration in the steam system is assumed to be negligible, no equations are needed for the steam system.

To calculate the distribution of tritium, we solve Eqs. 25-44 in the following sequence, discussed in more detail in Section III.

1. Assume a tritium concentration, C_{CT} , in the secondary system and calculate Q_{40} , Q_{41} , Q_{43} through Q_{49} from Eqs. 34, 35, 37a, 37b, 38a, 38b, 39a, 39b, 40, 41, 42 and 43.
2. Calculate Q_{42} from the material balance, Eq. 44.
3. Calculate C_{FT} from Eqs. 28a, 28b, 36a and 36b, the relationship $Q_{34} = -Q_{42}$ and the value of Q_{42} from step 2. If the value of C_{FT} is negative, increase the estimate for C_{CT} and repeat steps 1 through 3. When we have found a positive C_{FT} , we proceed to step 4.

4. Calculate Q_{31} , Q_{32} , Q_{33} , Q_{35} , Q_{36} , Q_{37} and Q_{38} from Eqs. 25-32.
5. Calculate R_F , where

$$R_F = \sum_{i=31}^{38} Q_i - R_1$$

is the term that must be added to the left side of Eq. 33 in order for the equation to balance. If R_F is positive, tritium must be added to the primary system, so C_{FT} and C_{CT} are too large; if R_F is negative, C_{FT} and C_{CT} are too small. Adjust the value of C_{CT} and repeat steps 1 through 5. When $|R_F/R_1|$ is sufficiently small, the calculations are finished.

III. SOLUTION OF EQUATIONS

In the procedure discussed above, we begin with the calculation of C_{CF} , C_{1s} , Q_{1s} and Q_{19} , with Eqs. 19a, 19b and 20, and the material balance, Eq. 21b:

$$Q_{1s} = h_{1s}A_{1s}(C_{CF} - C_{1s}) \quad (19a)$$

$$= B_4W_4A_{1s}(k_{1s}C_{1s})^{\frac{1}{2}}, \quad (19b)$$

$$Q_{19} = F_4E_4C_{CF}, \quad (20)$$

$$R_5 = Q_{1s} + Q_{19}. \quad (21b)$$

Eq. 19b requires that $Q_{1s} \geq 0$ and Eq. 20 requires that $Q_{19} \geq 0$, so if $R_5 = 0$, 21b requires that $Q_{1s} = Q_{19} = 0$. If $R_5 > 0$, we combine 21b, 20 and 19a to obtain

$$R_5 - Q_{1s} = F_4E_4C_{CF} = R_5 - h_{1s}A_{1s}(C_{CF} - C_{1s}),$$

or

$$C_{CF} = \frac{R_5 + h_{1s}A_{1s}C_{1s}}{F_4E_4 + h_{1s}A_{1s}}. \quad (19c)$$

Substituting 19c into 19a, setting the result equal to 19b and collecting terms we obtain

$$\alpha - C_{1s} = \beta C_{1s}^{\frac{1}{2}}, \quad (19d)$$

where we have defined

$$\alpha = \frac{R_5}{F_4E_4},$$

and

$$\beta = \left[\frac{F_4E_4 + h_{1s}A_{1s}}{F_4E_4} \right] \left[\frac{B_4W_4}{H_{1s}} \right] [k_{1s}]^{\frac{1}{2}}.$$

Squaring both sides of 19d results in a quadratic equation for C_{18} ; since the right-hand side of 19d is positive, we want the root of this quadratic which is less than α . We have

$$C_{18}^2 - (2\alpha + \beta^2)C_{18} + \alpha^2 = 0 ,$$

$$C_{18} = \frac{2\alpha + \beta^2 \pm \sqrt{(2\alpha + \beta^2)^2 - 4\alpha^2}}{2} .$$

To obtain the root less than α , we want the root with the negative sign. To avoid possible loss of significant figures, we note that the product of the roots is α^2 , so that we can write the solution in the form

$$C_{18} = \frac{\alpha^2}{\alpha + \frac{\beta^2}{2} \left(1 + \sqrt{1 + \frac{4\alpha}{\beta^2}} \right)} . \quad (19e)$$

Then we have

$$Q_{18} = B_4 W_4 A_{18} (k_{18} C_{18})^{\frac{1}{2}} , \quad (19b)$$

$$C_{CF} = \frac{R_5 + h_{18} A_{18} C_{18}}{F_4 E_4 + h_{18} A_{18}} , \quad (19c)$$

and

$$Q_{19} = F_4 E_4 C_{CF} . \quad (20)$$

With some value for C_C we proceed to the calculation of Q_{10} , Q_{11} , Q_{16} , Q_{17} and C_{16} . Eqs. 11a, 11b, 12a and 12b read

$$Q_{10} = h_{10} A_{10} (C_C - C_{10}) , \quad (11a)$$

$$Q_{10} = \frac{p_{10} A_{10}}{t_{10}} (k_{10} C_{10})^{\frac{1}{2}} , \quad (11b)$$

$$Q_{11} = h_{11} A_{11} (C_C - C_{11}) , \quad (12a)$$

$$Q_{11} = \frac{P_{11}A_{11}}{t_{11}} (k_{11}C_{11})^{\frac{1}{2}} . \quad (12b)$$

These equations (11 and 12) are identical in structure, as are Eqs. 1, 2, 3, 5, 7, 17 and 19. For Eqs. 11 and 12 we define

$$C_i = C_C, \quad \alpha = k_i \left(\frac{p_i}{t_i h_i} \right)^2, \quad i = 10, 11 ,$$

and Eqs. 11 and 12 then can be written in the form of quadratics in the concentration C_i :

$$C_i^2 - (2C_i + \alpha)C_i + C_i^2 = 0 .$$

From Eqs. 11b and 12b, the flow rates Q_{10} and Q_{11} must be positive, so that the root desired in each case is the smaller one. We have

$$C_i = \frac{C_i^2}{C_i + \frac{\alpha}{2} \left(1 + \sqrt{1 + \frac{4C_i}{\alpha}} \right)}, \quad i = 10, 11 ,$$

and

$$Q_i = \frac{p_i A_i}{t_i} (k_i C_i)^{\frac{1}{2}}, \quad i = 10, 11 .$$

By putting

$$C_i = C_C ,$$

$$\alpha = \left(\frac{B_3 W_3}{h_{16}} \right)^2 k_{16} ,$$

C_{16} can be calculated in the same fashion (Eqs. 17a and 17b) and the flow rates Q_{16} and Q_{17} are

$$Q_{16} = B_3 W_3 A_{16} (k_{16} C_{16})^{\frac{1}{2}} ,$$

$$Q_{17} = F_3 E_3 C_C .$$

We continue with step 4, the calculation of the flow rates Q_{13} , Q_{14} and Q_{15} , and the corresponding concentrations C_{13} , C_{14} and C_{15} , using Eqs. 14a, 14b, 15a, 15b, 16a, 16b, 22a, 22b, 23a, 23b, 24a and 24b. Note that the secondary system and the steam system are coupled by the equations

$$Q_{13} = -Q_{21}, \quad Q_{14} = -Q_{22} \quad \text{and} \quad Q_{15} = -Q_{23} .$$

The three equations 14, 15 and 16 all have the same structure and can be written in the form

$$h_K(C_1 - C_K) = \frac{P_K}{t_K} \left[(k_K C_K)^{\frac{1}{2}} - (k_L C_L)^{\frac{1}{2}} \right] , \quad (a)$$

$$h_L(C_L - C_2) = h_K(C_1 - C_K) , \quad (b)$$

where $K = 13, 14$ and 15 , $C_1 = C_C$, $L = 21, 22$ and 23 , and we identify C_2 as C_{SG} , C_{SS} and C_{SR} for $K = 13, 14$ and 15 , respectively. We can solve Eq. b for C_L :

$$C_L = \frac{h_K(C_1 - C_K) + h_L C_2}{h_L} = \frac{h_K}{h_L} (C_1 - C_K) + C_2 . \quad (c)$$

Since C_L must be non-negative, there is a maximum permissible value $C_K^{(max)}$, which is the value such that

$$\frac{h_K}{h_L} (C_1 - C_K^{(max)}) + C_2 = 0 ,$$

or

$$C_K^{(max)} = C_1 + \frac{h_L}{h_K} C_2 . \quad (d)$$

If we substitute (c) into (a) and rearrange, we have

$$C_K = C_1 + \frac{P_K}{h_K t_K} \left\{ k_L^{\frac{1}{2}} \left[\frac{h_K}{h_L} (C_1 - C_K) + C_2 \right]^{\frac{1}{2}} - [k_K C_K]^{\frac{1}{2}} \right\}, \quad (e)$$

or, more concisely,

$$C_K = F(C_K) .$$

To locate the solutions (if any) of this equation, we need to examine the behavior of $F(C_K)$ for $0 \leq C_K \leq C_K^{(\max)}$. We find that

$$F(0) > 0$$

and

$$F'(C_K) < 0, \quad F'(0) = -\infty$$

$$F''(C_K) \geq 0 .$$

The graph of $F(C_K)$ then looks like the curve in Fig. 2.

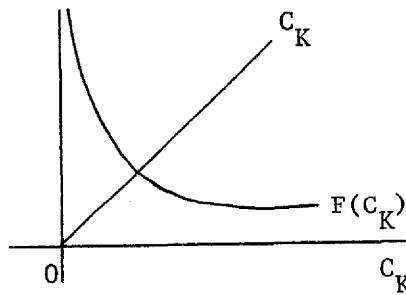


Fig. 2. Sketch of $F(C_K)$ vs C_K .

For there to be a solution between zero and $C_K^{(\max)}$, we must have $C_K^{(\max)} > F(C_K^{(\max)})$ and upon substitution of our expression (d) into $F(C_K)$, we find that this condition is satisfied. We will now examine the function

$$G(C_K) = C_K - F(C_K) .$$

We note that

$$G(0) = -F(0) < 0$$

$$G(C_K^{(\max)}) > 0$$

and

$$G'(C_K) = 1 - F'(C_K) > 0 \quad [\text{since } F'(C_K) < 0] .$$

This insures that $G(C_K)$ has one and only one zero in the range $0 \leq C_K \leq C_K^{(\max)}$. Since $G''(C_K) = -F''(C_K)$, $G''(C_K) \leq 0$, and the graph of $G(C_K)$ looks like the curve shown in Fig. 3.

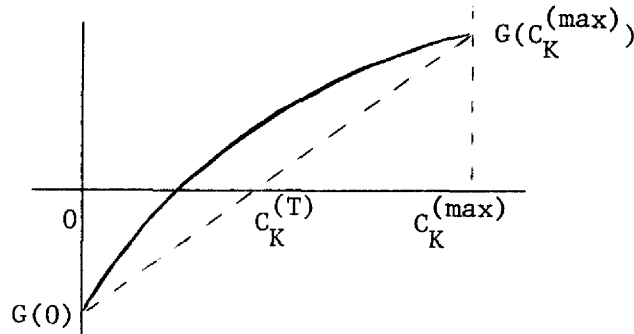


Fig. 3. Sketch of $G(C_K)$ vs C_K

With a suitable $C_K^{(1)}$ we can compute $G_1 = G(C_K^{(1)}) < 0$ (for example, starting with $C_K^{(1)} = 0$) and with a suitable $C_K^{(2)}$, $G_2 = G(C_K^{(2)}) > 0$ ($C_K^{(2)} = C_K^{(\max)}$, to start). An approximation to the solution $C_K^{(T)}$, is derived from the inverse linear interpolation:

$$C_K^{(T)} = \frac{G_2 C_K^{(1)} - G_1 C_K^{(2)}}{G_2 - G_1} ,$$

as shown in Fig. 2. A better approximation can be derived with inverse quadratic interpolation:

$$C_K^{(x)} = \frac{(0-G_T)(0-G_2)}{(G_1-G_T)(G_1-G_2)} C_K^{(1)} + \frac{(0-G_1)(0-G_T)}{(G_2-G_1)(G_2-G_T)} C_K^{(2)} + \frac{(0-G_1)(0-G_2)}{(G_T-G_1)(G_T-G_2)} C_K^{(T)} .$$

With $G''(C_K) \leq 0$ as shown and $G'(C_K) > 0$, $G_T = G(C_K^{(T)})$ will be positive and $C_K^{(T)}$ should be larger than the root. If $C_K^{(x)}$ is larger than $C_K^{(T)}$, we replace $C_K^{(2)}$ by $C_K^{(T)}$, G_2 by G_T , and repeat the inverse linear interpolation. If, however, $C_K^{(x)}$ is smaller than $C_K^{(T)}$, we calculate $G_x = G(C_K^{(x)})$; and if this value is negative, we replace $C_K^{(1)}$ by $C_K^{(x)}$, G_1 by G_x , $C_K^{(2)}$ by $C_K^{(T)}$ and G_2 by G_T , and repeat the inverse linear interpolation. If G_x is positive, we replace $C_K^{(2)}$ by $C_K^{(x)}$ and G_2 by G_x and repeat the inverse linear interpolation. We terminate this process when

$$\left| 1 - \frac{C_K^{(T)}}{C_K^{(x)}} \right| < C_{TOL} ,$$

or when we have done 50 iterations. The tolerance C_{TOL} is defined in a DATA statement in our program. We have found that the procedure converges in about four iterations for $C_{TOL} = 10^{-5}$ and in about six iterations for $C_{TOL} = 10^{-7}$.

The required flow rates Q_{13} , Q_{21} , Q_{14} , Q_{22} , Q_{15} and Q_{23} can now be computed from

$$Q_i = h_i A_i (C_C - C_i)$$

$$Q_{i+8} = -Q_i, \quad i = 13, 14, 15 .$$

The flow rate of hydrogen and tritium through heat exchanger tube walls from the secondary to the primary system, Q_{12} , is

$$Q_{12} = R_3 + R_4 - (Q_{10} + Q_{11} + Q_{13} + Q_{14} + Q_{15} + Q_{16} + Q_{17}) ,$$

and from Eq. 13a,

$$C_{12} = C_C - \frac{Q_{12}}{h_{12} A_4} .$$

If the value for $C_{1,2}$ is negative, we have used too small a value for C_C , so we double our previous guess and start over at step 3. If the computed value is positive, we proceed to calculate (Eq. 13b)

$$C_4 = \frac{1}{k_4} \left[(k_{1,2} C_{1,2})^{\frac{1}{2}} - \frac{Q_{1,2} t_4}{p_4 A_4} \right]^2,$$

and finally,

$$C_F = C_4 - \frac{Q_{1,2}}{h_4 A_4}.$$

If the computed value for C_F is negative, we need a larger value for C_C , so we double our previous guess and return to step 3. If positive, we proceed to step 7, the computation of the remaining flow rates Q_1 , Q_2 , Q_3 , Q_5 , Q_6 , Q_7 and Q_8 and the concentrations C_5 , C_{FF} and C_7 .

We can write Eqs. 1, 2 and 3 in the form

$$Q_i = h_i A_i (C_F - C_i) = \frac{p_i A_i}{t_i} (k_i C_i)^{\frac{1}{2}}, \quad i = 1, 2, 3,$$

and with

$$\alpha = \left(\frac{p_i}{t_i h_i} \right)^2 k_i$$

the resulting quadratic equations can be solved in the same way as those for C_{10} and C_{11} . Eqs. 5 can be manipulated into the same form with

$$\alpha = \left(\frac{B_1 W_1}{h_5} \right)^2 k_5$$

so that we can calculate C_5 , and from it

$$Q_5 = B_1 W_1 A_5 (k_5 C_5)^{\frac{1}{2}}. \quad (5b)$$

Again, Eqs. 7a and 7b can be written as a quadratic for C_7 with

$$\alpha = \left(\frac{B_2 W_2}{h_7} \right)^2 k_7$$

so that we can calculate

$$Q_7 = B_2 W_2 A_7 (k_7 C_7)^{\frac{1}{2}}$$

$$Q_8 = F_2 E_2 C_{FF}$$

and

$$R_F = \sum_{i=1}^8 Q_i - R_1 - R_2$$

where C_{FF} is

$$C_{FF} = \frac{MU}{k_7} \left(k_5 C_F \right)^{\frac{1}{2}} \quad . \quad (10)$$

This is the end of the first part of the procedure if R_F is small enough. We test the condition

$$\left| \frac{R_F}{R_1 + R_2} \right| < T_{TOL}$$

(where the quantity T_{TOL} is defined in a DATA statement in our program) and if it is satisfied, we proceed to the second part. If not, we adjust C_C in a variety of ways, depending on what information we have accumulated so far. We carry out a preliminary search for two values of C_C which bracket the root, i.e., one for which R_F is negative and the other for which R_F is positive. If this is the first iteration or if both our present and previous values of R_F have the same sign, we multiply C_C by a factor m such that

$$m = 10^{-R_F / (R_1 + R_2)}$$

but limited to the range

$$.01 \leq m \leq 100 \quad .$$

When we have bracketed the root, we combine inverse linear and inverse quadratic interpolation in much the same way as we did for the solution of the equations for C_{13} , C_{14} and C_{15} , keeping the root bracketed and attempting to reduce the length of the interval containing the root. When this process has converged, we proceed to the tritium calculation.

With a value for C_{CT} , the concentration of tritium in the secondary salt, we compute

$$Q_{40} = \frac{C_{CT}}{C_C} Q_{10} \quad (34)$$

$$Q_{41} = \frac{C_{CT}}{C_C} Q_{11} \quad (35)$$

and from Eqs. 37a, 37b, 38a, 38b, 39a and 39b we obtain

$$C_{43} = \frac{h_{13}t_{13}(C_{13}/k_{13})^{\frac{1}{2}}/p_{13}}{1+h_{13}t_{13}(C_{13}/k_{13})^{\frac{1}{2}}/p_{13}} C_{CT} \quad (37c)$$

$$Q_{43} = \frac{p_{13}A_{13}}{t_{13}(C_{13}/k_{13})^{\frac{1}{2}}} C_{43} \quad (37b)$$

$$C_{44} = \frac{h_{14}t_{14}(C_{14}/k_{14})^{\frac{1}{2}}/p_{14}}{1+h_{14}t_{14}(C_{14}/k_{14})^{\frac{1}{2}}/p_{14}} C_{CT} \quad (38c)$$

$$Q_{44} = \frac{p_{14}A_{14}}{t_{14}(C_{14}/k_{14})^{\frac{1}{2}}} C_{44} \quad (38b)$$

$$C_{45} = \frac{h_{15}t_{15}(C_{15}/k_{15})^{\frac{1}{2}}/p_{15}}{1+h_{15}t_{15}(C_{15}/k_{15})^{\frac{1}{2}}/p_{15}} C_{CT} \quad (39c)$$

$$Q_{45} = \frac{p_{15}A_{15}}{t_{15}(C_{15}/k_{15})^{\frac{1}{2}}} C_{45} \quad (39b)$$

$$Q_{i+30} = \frac{C_{CT}}{C_C} Q_i, \quad i = 16, 17, 18, 19 \quad (40-43)$$

$$Q_{42} = R_3 - Q_{40} - Q_{41} - Q_{43} - Q_{44} - Q_{45} - Q_{46} - Q_{47} - Q_{48} - Q_{49} \quad (44)$$

and finally

$$C_{42} = C_{CT} - \frac{Q_{42}}{h_{12}A_4} .$$

If this value is negative, we have used too small a value for C_{CT} ; in the same way as before, we double C_{CT} and try again, starting at Eq. 34. When we have found a positive C_{42} , we compute

$$C_{34} = \left(\frac{C_4}{k_4} \right)^{\frac{1}{2}} \left[\frac{C_{42}}{(C_{12}/k_{12})^{\frac{1}{2}}} - \frac{t_4 Q_{42}}{p_4 A_4} \right]$$

Again, if C_{34} is negative, we need to double C_{CT} and try again.

When we have found a positive C_{34} , we compute

$$C_{FT} = C_{34} - \frac{Q_{42}}{h_4 A_4}$$

and continue with the doubling scheme until C_{42} , C_{34} and C_{FT} are all positive. We can now compute the flow rates

$$Q_{30+i} = \frac{C_{FT}}{C_F} Q_i, \quad i = 1, 2, 3, 5, 6, 7, 8$$

and

$$R_F = \sum_{i=31}^{38} Q_i - R_1 .$$

Our test is now on $|R_F/R_1|$, and we use the same adjustment and interpolation procedures as for C_C .



IV. NOMENCLATURE

	<u>Reference Value*</u>	<u>Name**</u>
A = surface area, cm ²		A
A ₁ = hot leg of primary system (piping and pumps)	6 X 10 ⁵	
A ₂ = cold leg of primary system (piping)	5 X 10 ⁵	
A ₃ = reactor vessel and heat exchanger shells	3.5 X 10 ⁶	
A ₄ = tubes of primary heat exchanger	4.9 X 10 ⁷	
A ₅ = core graphite for sorption of hydrogen	5.2 X 10 ⁷	
A ₆ = --	---	
A ₇ = core graphite for sorption of hydrogen fluoride	5.2 X 10 ⁷	
A ₈ = --	---	
A ₉ = --	---	
A ₁₀ = hot leg of secondary system (piping, pumps, half of shells on steam-raising equipment)	1.1 X 10 ⁷	
A ₁₁ = cold leg of secondary system (piping, half of shells on steam-raising equipment)	8.8 X 10 ⁶	
A ₁₂ = A ₄	4.9 X 10 ⁷	
A ₁₃ = tubes of steam generators	3.1 X 10 ⁷	
A ₁₄ = tubes of superheaters	2.7 X 10 ⁷	
A ₁₅ = tubes of reheaters	1.8 X 10 ⁷	
A ₁₆ = sorber of hydrogen	0	
A ₁₇ = --	---	
A ₁₈ = sorber of hydrogen fluoride	0	

*The reference values are based on the design of a 1000 MWe molten salt breeder reactor plant described in ORNL-4541.

**Acronym used in FORTRAN computer program; if no entry appears, the parameter is not used in the program.

	<u>Reference Value</u>	<u>Name</u>
B = sorption factor, atoms/cm ² atm ^{1/2}		B
B ₁ = hydrogen + tritium on core graphite	3 X 10 ²¹	
B ₂ = hydrogen fluoride on core graphite	3 X 10 ²¹	
B ₃ = hydrogen + tritium on sorber in secondary system	1 X 10 ¹⁸	
B ₄ = hydrogen fluoride on sorber in secondary system	1 X 10 ¹⁸	
C = concentration, atoms/cm ³		
C _F = hydrogen + tritium in bulk of primary salt		CF
C _{FF} = hydrogen + tritium as hydrogen fluoride in bulk of primary salt		CFF
C _{FT} = tritium in bulk of primary salt		CFT
C _C = hydrogen + tritium in bulk of secondary salt		CC
C _{CF} = hydrogen + tritium as hydrogen fluoride in bulk of secondary salt		CCF
C _{CT} = tritium in bulk of secondary salt		CCT
C _{SG} = hydrogen in bulk of water in steam generator (672°K)	2 X 10 ¹⁰	CSG
C _{SS} = hydrogen in bulk of steam in superheater (783°K)	9 X 10 ¹¹	CSS
C _{SR} = hydrogen in bulk of steam in reheater (755°K)	1 X 10 ¹¹	CSR
C ₁ = hydrogen + tritium in salt at surface of hot leg of primary system		C
C ₂ = hydrogen + tritium in salt at surface of cold leg of primary system		
C ₃ = hydrogen + tritium in salt at surface of reactor vessel and heat exchanger shells		

	<u>Reference Value</u>	<u>Name</u>
C ₄ = hydrogen + tritium in salt at surfaces of heat exchanger tubes in primary system		
C ₅ = hydrogen + tritium in salt at surfaces of core graphite in primary system		
C ₆ = --		
C ₇ = hydrogen fluoride in salt at surfaces of core graphite in primary system		
C ₈ = --		
C ₉ = --		
C ₁₀ = hydrogen + tritium in salt at surface of hot leg in secondary system		
C ₁₁ = hydrogen + tritium in salt at surface of cold leg in secondary system		
C ₁₂ = hydrogen + tritium in salt at surfaces of heat exchanger tubes in secondary system		
C ₁₃ = hydrogen + tritium in salt at surfaces of steam generator tubes in secondary system		
C ₁₄ = hydrogen + tritium in salt at surfaces of superheater tubes in secondary system		
C ₁₅ = hydrogen + tritium in salt at surfaces of reheater tubes in secondary system		
C ₁₆ = hydrogen + tritium in salt at surfaces of sorber in secondary system		
C ₁₇ = --		
C ₁₈ = hydrogen fluoride in salt at surfaces of sorber in secondary system		
C ₁₉ = --		

	<u>Reference Value</u>	<u>Name</u>
C_{20}		--
C_{21}		hydrogen in steam at surfaces of steam generator tubes in steam system
C_{22}		hydrogen in steam at surfaces of super-heater tubes in steam system
C_{23}		hydrogen in steam surfaces of reheater tubes in steam system
C_{24} - C_{33}		--
C_{34}		tritium in salt at surfaces of heat exchanger tubes in primary system
C_{35} - C_{41}		--
C_{42}		tritium in salt at surfaces of heat exchanger tubes in secondary system
C_{43}		tritium in salt at surfaces of steam generator tubes in secondary system
C_{44}		tritium in salt at surfaces of super-heater tubes in secondary system
C_{45}		tritium in salt at surfaces of reheater tubes in secondary system
E		efficiency
E_1	5×10^{-1}	removal of hydrogen + tritium from purge stream in primary system
E_2	1.7×10^{-2}	removal of hydrogen fluoride from purge stream in primary system
E_3	1.8×10^{-1}	removal of hydrogen + tritium from purge stream in secondary system
E_4	1.8×10^{-3}	removal of hydrogen fluoride from purge stream in secondary system

	<u>Reference Value</u>	<u>Name</u>
F = flow rate, cm ³ /sec		F
F ₁ = purge stream for removal of hydrogen + tritium from primary system	3.6 X 10 ⁵	
F ₂ = purge stream for removal of hydrogen fluoride from primary system	3.6 X 10 ⁵	
F ₃ = purge stream for removal of hydrogen + tritium from secondary system	5.0 X 10 ⁵	
F ₄ = purge stream for removal of hydrogen fluoride from secondary system	5.0 X 10 ⁵	
h = mass transfer coefficient, cm/sec		H
h ₁ = hydrogen through primary salt to surfaces of hot leg in primary system	1.6 X 10 ⁻²	
h ₂ = hydrogen through primary salt to surfaces of cold leg in primary system	6.0 X 10 ⁻³	
h ₃ = hydrogen through primary salt to surfaces of reactor vessel and heat exchanger shells in primary system	9.0 X 10 ⁻⁵	
h ₄ = hydrogen through primary salt to surfaces of heat exchanger tubes in primary system	1.9 X 10 ⁻²	
h ₅ = hydrogen through primary salt to surfaces of core graphite in primary system	3.0 X 10 ⁻³	
h ₆ = --	---	
h ₇ = hydrogen fluoride through primary salt to surfaces of core graphite in primary system	3.0 X 10 ⁻³	
h ₈ = --	---	
h ₉ = --	---	
h ₁₀ = hydrogen through secondary salt to surfaces of hot leg in secondary system	7.4 X 10 ⁻²	

	<u>Reference Value</u>	<u>Name</u>
h_{11} = hydrogen through secondary salt to surfaces of cold leg in secondary system	3.4×10^{-2}	
h_{12} = hydrogen through secondary salt to surfaces of tubes in heat exchangers in secondary system	9.7×10^{-2}	
h_{13} = hydrogen through secondary salt to surfaces of tubes of steam generators in secondary system	4.3×10^{-2}	
h_{14} = hydrogen through secondary salt to surfaces of tubes in superheaters in secondary system	4.7×10^{-2}	
h_{15} = hydrogen through secondary salt to surfaces of tubes in reheaters in secondary system	4.0×10^{-2}	
h_{16} = hydrogen through secondary salt to surfaces of sorber in secondary system	8.0×10^{-1}	
h_{17} = --	---	
h_{18} = hydrogen fluoride through secondary salt to surfaces of sorber in secondary system	8.0×10^{-1}	
h_{19} = --	---	
h_{20} = --	---	
h_{21} = hydrogen through water to surfaces of tubes of steam generators in steam system	5.8	
h_{22} = hydrogen through steam to surfaces of tubes of steam generators in steam system	12	
h_{23} = hydrogen through steam to surfaces of tubes of reheaters in steam system	30	

k = Henry's law coefficient, $\frac{(\text{cm}^3 \text{ melt})(\text{atm.})}{\text{atom H}}$ K

$$= 0.83 \times 10^{-24} \left[k' \frac{\text{moles H}_2}{(\text{cm}^3 \text{ melt})(\text{atm.})} \right]^{-1}$$

	<u>Reference Value</u>	<u>Name</u>
$= 1.7 \times 10^{-24} \left[k' \frac{\text{moles HF}}{(\text{cm}^3 \text{ melt})(\text{atm.})} \right]^{-1}$		
k_1 = hydrogen in primary salt in hot leg in primary system (973°K)	1.2×10^{-17}	
k_2 = hydrogen in primary salt in cold leg in primary system (838°K)	2.0×10^{-17}	
k_3 = hydrogen in primary salt in reactor vessel and heat exchanger shells in primary system (908°K)	1.5×10^{-17}	
k_4 = hydrogen in primary salt in heat exchangers in primary system (908°K)	1.5×10^{-17}	
k_5 = hydrogen in primary salt in reactor core in primary system (923°K)	1.4×10^{-17}	
k_6 = --	---	
k_7 = hydrogen fluoride in primary salt in reactor core in primary system (923°K)	1.5×10^{-19}	
k_8 = --	---	
k_9 = --	---	
k_{10} = hydrogen in secondary salt in hot leg in secondary system (894°K)	3.4×10^{-18}	
k_{11} = hydrogen in secondary salt in cold leg in secondary system (723°K)	5.0×10^{-18}	
k_{12} = hydrogen in secondary salt in heat exchangers in secondary system (809°K)	4.0×10^{-18}	
k_{13} = hydrogen in secondary salt in steam generators in secondary system (783°K)	4.5×10^{-18}	
k_{14} = hydrogen in secondary salt in superheaters in secondary system (866°K)	3.5×10^{-18}	
k_{15} = hydrogen in secondary salt in reheaters in secondary system (810°K)	4.0×10^{-18}	

	<u>Reference Value</u>	<u>Name</u>
k_{16} = hydrogen in secondary salt in contact with sorber in secondary system (773°K)	4.4×10^{-18}	
k_{17} = ---	---	
k_{18} = hydrogen fluoride in secondary salt in contact with sorber in secondary system (773°K)	1.1×10^{-20}	
k_{19} = --	---	
k_{20} = --	---	
k_{21} = hydrogen in steam in steam generators in steam system (660°K)	4.5×10^{-20}	
k_{22} = hydrogen in steam in superheaters in the steam system (755°K)	5.1×10^{-20}	
k_{23} = hydrogen in steam in reheaters in steam system (714°K)	4.8×10^{-20}	
M = equilibrium quotient for reduction of UF_4 by hydrogen, $atm^{1/2}$, (923°K)	1.12×10^{-6}	M
p = permeability coefficient for hydrogen in metal		P
$\frac{(\text{atoms H})(\text{mm})}{(\text{sec})(\text{cm}^2)(\text{atm})^{1/2}} = 1.5 \times 10^{16} p' \frac{(\text{cm}^3 \text{H}_2 \text{ STP})(\text{mm})}{(\text{hr})(\text{cm}^2)(\text{atm})^{1/2}}$		
p_1 = at average temperature of metal in hot leg in primary system (973°K)	2.1×10^{15}	
p_2 = at average temperature of metal in cold leg in primary system (838°K)	6.7×10^{14}	
p_3 = at average temperature of metal in reactor vessel and heat exchanger shells in primary system (873°K)	9.0×10^{14}	
p_4 = at average temperature of metal in tubes in heat exchangers in primary system (873°K)	9.0×10^{14}	

	<u>Reference Value</u>	<u>Name</u>
$P_5 - P_9 = \text{---}$	---	
$P_{10} =$ at average temperature of metal in hot leg in secondary system (893°K)	1.1×10^{15}	
$P_{11} =$ at average temperature of metal in cold leg in secondary system (723°K)	1.8×10^{14}	
$P_{12} = P_4$	9.0×10^{14}	
$P_{13} =$ at average temperature of tubes in steam generators in secondary system (723°K)	1.8×10^{14}	
$P_{14} =$ at average temperature of tubes in super- heaters in secondary system (838°K)	6.7×10^{14}	
$P_{15} =$ at average temperature of tubes in reheaters in secondary system (773°K)	3.5×10^{14}	

P = pressure, atm. or other appropriate units

Q = rate of transport, atoms of hydrogen and/or
tritium per second

Q

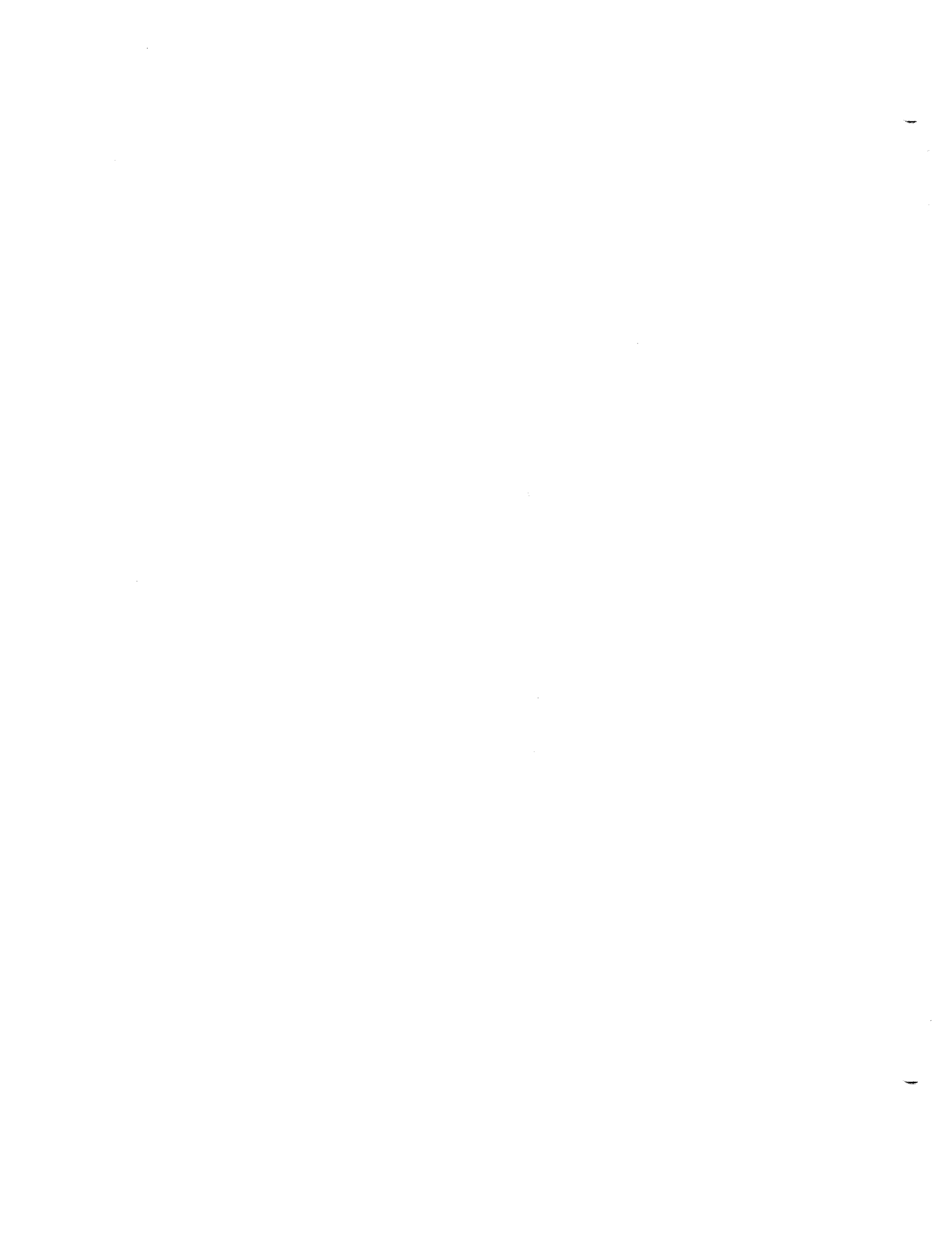
- $Q_1 =$ hydrogen + tritium through walls of hot
leg in primary system
- $Q_2 =$ hydrogen + tritium through walls of cold
leg in primary system
- $Q_3 =$ hydrogen + tritium through wall of reactor
vessel and shells of heat exchangers in
primary system
- $Q_4 =$ hydrogen + tritium through walls of tubes
in heat exchangers from primary system to
secondary system
- $Q_5 =$ hydrogen + tritium to core graphite in
primary system
- $Q_6 =$ hydrogen + tritium to purge in primary
system

	<u>Reference</u> <u>Value</u>	<u>Name</u>
Q_7		= hydrogen fluoride to core graphite in primary system
Q_8		= hydrogen fluoride to purge in primary system
Q_9		= --
Q_{10}		= hydrogen + tritium through walls of hot leg in secondary system
Q_{11}		= hydrogen + tritium through walls of cold leg in secondary system
Q_{12}		= hydrogen + tritium through walls of tubes in heat exchangers from secondary system to primary system = $-Q_4$
Q_{13}		= hydrogen + tritium through walls of the steam generator tubes from the secondary system into the steam system
Q_{14}		= hydrogen + tritium through walls of the superheater tubes from the secondary system into the steam system
Q_{15}		= hydrogen + tritium through walls of the reheater tubes from the secondary system into the steam system
Q_{16}		= hydrogen + tritium to sorber in secondary system
Q_{17}		= hydrogen + tritium to purge in secondary system
Q_{18}		= hydrogen fluoride to sorber in secondary system
Q_{19}		= hydrogen fluoride to purge in secondary system
Q_{20}		= --
Q_{21}		= hydrogen through walls of steam generator tubes from steam system into secondary system = $-Q_{13}$

	<u>Reference Value</u>	<u>Name</u>
Q_{22}	=	hydrogen through walls of superheater tubes from steam system into secondary system = $-Q_{14}$
Q_{23}	=	hydrogen through walls of reheater tubes from steam system into secondary system = $-Q_{15}$
Q_{24}	=	--
Q_{30}	=	--
Q_{31}	=	tritium through walls of hot leg in primary system
Q_{32}	=	tritium through walls of cold leg in primary system
Q_{33}	=	tritium through wall of reactor vessel and shells of heat exchangers in primary system
Q_{34}	=	tritium through walls of heat exchanger tubes from primary system into secondary system
Q_{35}	=	tritium to core graphite in primary system
Q_{36}	=	tritium to purge in primary system
Q_{37}	=	tritium fluoride to core graphite in primary system
Q_{38}	=	tritium fluoride to purge in primary system
Q_{39}	=	--
Q_{40}	=	tritium through walls of hot leg in secondary system
Q_{41}	=	tritium through walls of cold leg in secondary system
Q_{42}	=	tritium through walls of heat exchanger tubes from secondary system into primary system = $-Q_{34}$

	<u>Reference Value</u>	<u>Name</u>
Q ₄₃ = tritium through walls of steam generator tubes from secondary system into steam system		
Q ₄₄ = tritium through walls of superheater tubes from secondary system into steam system		
Q ₄₅ = tritium through walls of reheater tubes from secondary system into steam system		
Q ₄₆ = tritium to sorber in secondary system		
Q ₄₇ = tritium to purge in secondary system		
Q ₄₈ = tritium fluoride to sorber in secondary system		
Q ₄₉ = tritium fluoride to purge in secondary system		
R = rate of production or addition, atoms/sec		
		R
R ₁ = tritium in primary system	5.8 X 10 ¹⁷	
R ₂ = hydrogen to primary system	0	
R ₃ = tritium in secondary system	0	
R ₄ = hydrogen to secondary system	0	
R ₅ = hydrogen fluoride to secondary system	0	
R _F = hydrogen or tritium to primary system in order to obtain overall material balance	---	
T = temperature, °K		
t = wall thickness, mm		
		T
t ₁ = hot leg in primary system	13	
t ₂ = cold leg in primary system	13	

	<u>Reference Value</u>	<u>Name</u>
t_3 = reactor vessel and heat exchanger shells in primary system	50	
t_4 = tubes in heat exchangers in primary system	1	
t_5 - t_9 = --		
t_{10} = hot leg in secondary system	13	
t_{11} = cold leg in secondary system	13	
$t_{12} = t_4$		
t_{13} = tubes in steam generators	2	
t_{14} = tubes in superheaters	2	
t_{15} = tubes in reheaters	1	
U = ratio X_{UF_4} / X_{UF_3}	100	U
W = replacement rate, fraction/sec		W
W_1 = core graphite or other sorber of hydrogen in primary system	1	
W_2 = core graphite or other sorber of hydrogen fluoride in primary system	1	
W_3 = sorber of hydrogen in secondary system	1	
W_4 = sorber of hydrogen fluoride in secondary system	1	
X = mole fraction		



To change various system parameters, the command is

CHANGE__XXX

where XXX is replaced by the appropriate variable name as listed in Sec. IV. If the variable name refers to one of the named concentrations (C_F , C_{FF} , ..., C_{SR}), the next line of input must contain the new parameter value in cols. 1-10. If the variable name refers to any of the subscripted variables in Sec. IV, the next line must contain a starting index, n_1 , a stopping index n_2 and the new values for the variables specified by the subscripts n_1 through n_2 . A maximum of seven consecutive values is allowed; if there are more than seven, put the subsequent values on subsequent lines. End with a line with a starting index of zero. The following example illustrates the format.

CARD COLUMN																																					
1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	25	26	27	28	29	30	31	32	33	34	35	36	37	
C	H	A	N	G	E			A																													
	1				3			1	.	2					+ 6			1	.	0						+ 6			7	.	0					+ 6	
	1	3						6	2	.					+ 6																						
	0																																				

This will insert new values for A_1 , A_2 , A_3 and A_{13} of 1.2×10^6 , 1.0×10^6 , 7.0×10^6 and 62×10^6 , respectively. If only one value is to be changed, the second subscript need not appear.

The user can supply starting estimates for C_C and C_{CT} , the concentrations of hydrogen plus tritium and tritium in the bulk of the secondary salt, with the "CHANGE" command. If no values are supplied the program will use 1×10^{11} for C_C and 1×10^{10} for C_{CT} .

To perform a calculation when all the necessary changes have been made, the command is

```
RUN
```

A calculation will then be done with the parameters specified. For subsequent cases, all parameters will have the values present at the end of the preceding calculation; to change the parameters, the user can supply additional "CHANGE" commands. To reset parameters to their reference values, the command is

```
RESET ___XXX
```

If "XXX" is left blank, all parameters will be reset; if "XXX" is the name of a subscripted variable, all entries with the given name will be reset; and if "XXX" is the name of one of the named concentrations (C_F , C_{FF} , ..., C_{SR}) then just that concentration will be reset. If, for example, after running the case specified by the "CHANGE" command in the example, a user put

```
RESET ___A
```

then all the A's would be reset to their reference values.

The program will stop when an end-of-file condition is detected on the standard input unit, i.e., when it runs out of data.

The input and output for a sample problem are shown in Figs. 4 and 5. Reference values from Section IV were used in the sample calculation. The results indicate that 30 percent or more of the tritium might reach the steam system in a large power reactor unless special measures are taken to confine the tritium.

1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	25	26	27	28	29	30	31	32	33	34	35	36	37	38
---	---	---	---	---	---	---	---	---	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----

```

// XXX1 JØB ( n n n n n ) , ' A D D R E S S ' , C L A S S = A
// EXEC FØRTHLG , GØSIZE = 6 2 K
// LKED . SYSIN DD *
      :
      HEX DECK
      :
/*
// GØ . FT05F001 DD *
ØUTPUT ALL PRINTER
RUN
CHANGE A
      1      3      1 . 2      + 6      1 . 0      + 6      7 . 0      + 6
      13      6 . 2      + 7
      0
RUN
ZILCH
/*
//

```

Fig. 4. Sample Problem Input

VALUES IN ARRAY V

NAME A	DIMENSION	20 USED	18 STARTS AT	1	4	5
1	2	3	4	5		
6.00000D 05	5.00000D 05	3.50000D 06	4.90000D 07	5.20000D 07		
-1.00000D 00	5.20000D 07	-1.00000D 00	-1.00000D 00	1.10000D 07		
8.80000D 06	4.90000D 07	3.10000D 07	2.70000D 07	1.80000D 07		
0.0	-1.00000D 00	0.0				
NAME B	DIMENSION	5 USED	4 STARTS AT	21		
1	2	3	4	5		
3.00000D 21	3.00000D 21	1.00000D 18	1.00000D 18			
NAME C	DIMENSION	50 USED	45 STARTS AT	26		
1	2	3	4	5		
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
0.0	0.0	0.0	0.0	0.0		0.0
NAME CN	DIMENSION	10 USED	9 STARTS AT	76		
1	2	3	4	5		
-1.00000D 00	-1.00000D 00	-1.00000D 00	1.00000D 11	-1.00000D 00		
1.00000D 10	2.00000D 10	9.00000D 11	1.00000D 11			
NAME E	DIMENSION	5 USED	4 STARTS AT	86		
1	2	3	4	5		
5.00000D-01	1.70000D-02	1.80000D-01	1.80000D-03			
NAME F	DIMENSION	5 USED	4 STARTS AT	91		
1	2	3	4	5		
3.60000D 05	3.60000D 05	5.00000D 05	5.00000D 05			
NAME H	DIMENSION	25 USED	23 STARTS AT	96		
1	2	3	4	5		
1.60000D-02	6.00000D-03	9.00000D-05	1.90000D-02	3.00000D-03		
-1.00000D 00	3.00000D-03	-1.00000D 00	-1.00000D 00	7.40000D-02		
3.40000D-02	9.70000D-02	4.30000D-02	4.70000D-02	4.00000D-02		
8.00000D-01	-1.00000D 00	8.00000D-01	-1.00000D 00	-1.00000D 00		
5.80000D 00	1.20000D 01	3.00000D 01				

Fig. 5A. List of Parameter Values Used in Calculation.

VALUES IN ARRAY V

NAME K					
DIMENSION		25 USED	23 STARTS AT 121		
1	2		3	4	5
1.200000-17	2.000000-17		1.500000-17	1.500000-17	1.400000-17
-1.000000 00	1.500000-19		-1.000000 00	-1.000000 00	3.400000-18
5.000000-18	4.000000-18		4.500000-18	3.500000-18	4.000000-18
4.400000-18	-1.000000 00		1.100000-20	-1.000000 00	-1.000000 00
4.500000-20	5.100000-20		4.800000-20		
NAME P					
DIMENSION		20 USED	15 STARTS AT 146		
1	2		3	4	5
2.100000 15	6.700000 14		9.000000 14	9.000000 14	-1.000000 00
-1.000000 00	-1.000000 00		-1.000000 00	-1.000000 00	1.100000 15
1.800000 14	9.000000 14		1.800000 14	6.700000 14	3.500000 14
NAME R					
DIMENSION		10 USED	5 STARTS AT 166		
1	2		3	4	5
5.800000 17	0.0		0.0	0.0	0.0
NAME T					
DIMENSION		20 USED	15 STARTS AT 176		
1	2		3	4	5
1.300000 01	1.300000 01		5.000000 01	1.000000 00	-1.000000 00
-1.000000 00	-1.000000 00		-1.000000 00	-1.000000 00	1.300000 01
1.300000 01	1.000000 00		2.000000 00	2.000000 00	1.000000 00
NAME W					
DIMENSION		5 USED	4 STARTS AT 196		
1	2		3	4	5
1.000000 00	1.000000 00		1.000000 00	1.000000 00	
NAME M					
DIMENSION		1 USED	1 STARTS AT 201		
1	2		3	4	5
1.120000-06					
NAME U					
DIMENSION		1 USED	1 STARTS AT 202		
1	2		3	4	5
1.000000 02					

Fig. 5A. (Continued).

ITERATIVE SOLUTION FOR CC

NCC	CC1	CCL	CCX	RFX	CC2
0				3.47462D 17	1.00000D 11
1	2.52116D 10			-2.91224D 17	
2	2.52116D 10	5.93131D 10		1.62379D 16	1.00000D 11
			5.74238D 10	1.92741D 14	
4	2.52116D 10	5.74025D 10		1.13535D 13	5.74238D 10
			5.74012D 10	1.02747D 08	

ITERATIVE SOLUTION FOR CCT

NCC	CC1	CCL	CCX	RFX	CC2
0	1.00000D 10			-4.71400D 17	
1				4.33784D 16	5.74012D 10

Fig. 5B. Output from Iterative Calculations.

OUTPUT SUMMARY

STEAM SYSTEM

FLOW OF H + T INTO STEAM SYSTEM	1.71072D 17
FLOW OF T INTO STEAM SYSTEM	1.76310D 17
FLOW OF H INTO STEAM SYSTEM	-5.23800D 15
FRACTION OF T INTO STEAM SYSTEM	3.03983D-01

SECONDARY SYSTEM

FLOWS

H + T INTO SECONDARY FROM PRIMARY	2.38501D 17
T INTO SECONDARY FROM PRIMARY	2.39047D 17
H + T THRU PIPE WALLS INTO CELLS	6.22626D 16
T THRU PIPE WALLS INTO CELLS	5.79300D 16

SORPTION BY SINK

H + T	0.0
T	0.0
HF	0.0
TF	0.0

REMOVAL BY PURGE

H + T	5.16611D 15
T	4.80662D 15
HF	0.0
TF	0.0

FRACTION OF T

PASSING THRU PIPE WALLS	9.98793D-02
SORBED BY SINK AS T	0.0
SORBED BY SINK AS TF	0.0
REMOVED BY PURGE AS T	8.28727D-03
REMOVED BY PURGE AS TF	0.0

CONCENTRATIONS IN SECONDARY SALT

H + T (CC)	5.74012D 10
T (CCT)	5.34069D 10
HF (CCF)	0.0

PRIMARY SYSTEM

FLOWS

H + T THRU WALLS INTO CELL	3.68436D 15
T THRU WALLS INTO CELL	3.67847D 15

SORPTION BY SINK

H + T	4.45130D 16
T	4.44419D 16
HF	2.32807D 17
TF	2.32435D 17

REMOVAL BY PURGE

H + T	5.13612D 16
T	5.12791D 16
HF	9.13321D 15
TF	9.11861D 15

FRACTION OF T

PASSING THRU WALLS INTO CELL	6.34219D-03
SORBED BY SINK AS T	7.66239D-02
SORBED BY SINK AS TF	4.00750D-01
REMOVED BY PURGE AS T	8.84122D-02
REMOVED BY PURGE AS TF	1.57217D-02

CONCENTRATIONS IN PRIMARY SALT

H + T (CF)	2.85340D 11
T (CFT)	2.84884D 11
HF (CFF)	1.49235D 12

Fig. 5C. Output Summary.


```
A ( 1. 3)
----- 1.200000 06 1.000000 06 7.000000 06
A (13.13)
        6.200000 07
```

Fig. 5D. Output Produced by "CHANGE" Command.

VALUES IN ARRAY V

NAME A	DIMENSION	20 USED	18 STARTS AT	1	4	5
1	2	3	4	5		
1.20000D 06	1.00000D 06	7.00000D 06	4.90000D 07	5.20000D 07		
-1.00000D 00	5.20000D 07	-1.00000D 00	-1.00000D 00	1.10000D 07		
8.80000D 06	4.90000D 07	6.20000D 07	2.70000D 07	1.80000D 07		
0.0	-1.00000D 00	0.0				
NAME B	DIMENSION	5 USED	4 STARTS AT	21		
1	2	3	4	5		
3.00000D 21	3.00000D 21	1.00000D 18	1.00000D 18			
NAME C	DIMENSION	50 USED	45 STARTS AT	26		
1	2	3	4	5		
7.94414D 07	6.58548D 07	1.62035D 05	7.73273D 10	6.94438D-09		
0.0	1.62245D-05	0.0	0.0	6.13618D 08		
2.99629D 09	2.87318D 11	6.77408D 08	1.37958D 10	1.41313D 09		
4.06149D 02	0.0	0.0	0.0	0.0		
2.03867D 10	9.00153D 11	1.00069D 11	0.0	0.0		
0.0	0.0	0.0	0.0	0.0		
0.0	0.0	0.0	2.86783D 10	0.0		
0.0	0.0	0.0	0.0	0.0		
0.0	1.05570D 11	3.16772D 08	4.74597D 08	1.16510D 08		
NAME CN	DIMENSION	10 USED	9 STARTS AT	76		
1	2	3	4	5		
3.11803D 11	1.56002D 12	2.90143D 11	5.28420D 10	0.0		
5.43552D 10	2.00000D 10	9.00000D 11	1.00000D 11			
NAME E	DIMENSION	5 USED	4 STARTS AT	86		
1	2	3	4	5		
5.00000D-01	1.70000D-02	1.80000D-01	1.80000D-03			
NAME F	DIMENSION	5 USED	4 STARTS AT	91		
1	2	3	4	5		
3.60000D 05	3.60000D 05	5.00000D 05	5.00000D 05			
NAME H	DIMENSION	25 USED	23 STARTS AT	96		
1	2	3	4	5		
1.60000D-02	6.00000D-03	9.00000D-05	1.90000D-02	3.00000D-03		
-1.00000D 00	3.00000D-03	-1.00000D 00	-1.00000D 00	7.40000D-02		
3.40000D-02	9.70000D-02	4.30000D-02	4.70000D-02	4.00000D-02		
8.00000D-01	-1.00000D 00	8.00000D-01	-1.00000D 00	-1.00000D 00		
5.80000D 00	1.20000D 01	3.00000D 01				

Fig. 5E. List of Parameter Values Used in Calculation After "CHANGE" Command.

VALUES IN ARRAY V					
NAME K	DIMENSION	25 USED	23 STARTS AT 121		
1	2		3	4	5
1.200000-17	2.000000-17		1.500000-17	1.500000-17	1.400000-17
-1.000000 00	1.500000-19		-1.000000 00	-1.000000 00	3.400000-18
5.000000-18	4.000000-18		4.500000-18	3.500000-18	4.000000-18
4.400000-18	-1.000000 00		1.100000-20	-1.000000 00	-1.000000 00
4.500000-20	5.100000-20		4.800000-20		
NAME P	DIMENSION	20 USED	15 STARTS AT 146		
1	2		3	4	5
2.100000 15	6.700000 14		9.000000 14	9.000000 14	-1.000000 00
-1.000000 00	-1.000000 00		-1.000000 00	-1.000000 00	1.100000 15
1.800000 14	9.000000 14		1.800000 14	6.700000 14	3.500000 14
NAME R	DIMENSION	10 USED	5 STARTS AT 166		
1	2		3	4	5
5.800000 17	0.0		0.0	0.0	0.0
NAME T	DIMENSION	20 USED	15 STARTS AT 176		
1	2		3	4	5
1.300000 01	1.300000 01		5.000000 01	1.000000 00	-1.000000 00
-1.000000 00	-1.000000 00		-1.000000 00	-1.000000 00	1.300000 01
1.300000 01	1.000000 00		2.000000 00	2.000000 00	1.000000 00
NAME W	DIMENSION	5 USED	4 STARTS AT 196		
1	2		3	4	5
1.000000 00	1.000000 00		1.000000 00	1.000000 00	
NAME M	DIMENSION	1 USED	1 STARTS AT 201		
1	2		3	4	5
1.120000-06					
NAME U	DIMENSION	1 USED	1 STARTS AT 202		
1	2		3	4	5
1.000000 02					

Fig. 5E. (Continued).

ITERATIVE SOLUTION FOR CC

NCC	CC1	CCL	CCX	RFX	CC2
0				1.44089D 17	5.74012D 10
1	3.24168D 10			-1.33923D 17	
2	3.24168D 10	4.44522D 10		2.91408D 15	5.74012D 10
			4.41906D 10	6.33182D 12	
4	3.24168D 10	4.41900D 10		1.41909D 11	4.41906D 10

ITERATIVE SOLUTION FOR CCT

NCC	CC1	CCL	CCX	RFX	CC2
0				1.74565D 17	5.34069D 10
1	2.67277D 10			-2.02375D 17	

Fig. 5F. Output from Iterative Calculations With New Parameters.

OUTPUT SUMMARY

STEAM SYSTEM

FLOW OF H • T INTO STEAM SYSTEM	1.85851D 17
FLOW OF T INTO STEAM SYSTEM	1.89989D 17
FLOW OF H INTO STEAM SYSTEM	-4.13811D 15
FRACTION OF T INTO STEAM SYSTEM	3.27568D-01

SECONDARY SYSTEM

FLOWS

H + T INTO SECONDARY FROM PRIMARY	2.38032D 17
T INTO SECONDARY FROM PRIMARY	2.38464D 17
H + T THRU PIPE WALLS INTO CELLS	4.82035D 16
T THRU PIPE WALLS INTO CELLS	4.47799D 16

SORPTION BY SINK

H + T	0.0
T	0.0
HF	0.0
TF	0.0

REMOVAL BY PURGE

H + T	3.97710D 15
T	3.69463D 15
HF	0.0
TF	0.0

FRACTION OF T

PASSING THRU PIPE WALLS	7.72067D-02
SORBED BY SINK AS T	0.0
SORBED BY SINK AS TF	0.0
REMOVED BY PURGE AS T	6.37006D-03
REMOVED BY PURGE AS TF	0.0

CONCENTRATIONS IN SECONDARY SALT

H • T (CC)	4.41900D 10
T (CCT)	4.10515D 10
HF (CCF)	0.0

PRIMARY SYSTEM

FLOWS

H + T THRU WALLS INTO CELL	7.26328D 15
T THRU WALLS INTO CELL	7.25410D 15

SORPTION BY SINK

H + T	4.38759D 16
T	4.38205D 16
HF	2.31135D 17
TF	2.30843D 17

REMOVAL BY PURGE

H + T	5.06261D 16
T	5.05621D 16
HF	9.06761D 15
TF	9.05616D 15

FRACTION OF T

PASSING THRU WALLS INTO CELL	1.25071D-02
SORBED BY SINK AS T	7.55526D-02
SORBED BY SINK AS TF	3.98006D-01
REMOVED BY PURGE AS T	8.71760D-02
REMOVED BY PURGE AS TF	1.56141D-02

CONCENTRATIONS IN PRIMARY SALT

H + T (CF)	2.81256D 11
T (CFT)	2.80901D 11
HF (CFF)	1.48164D 12

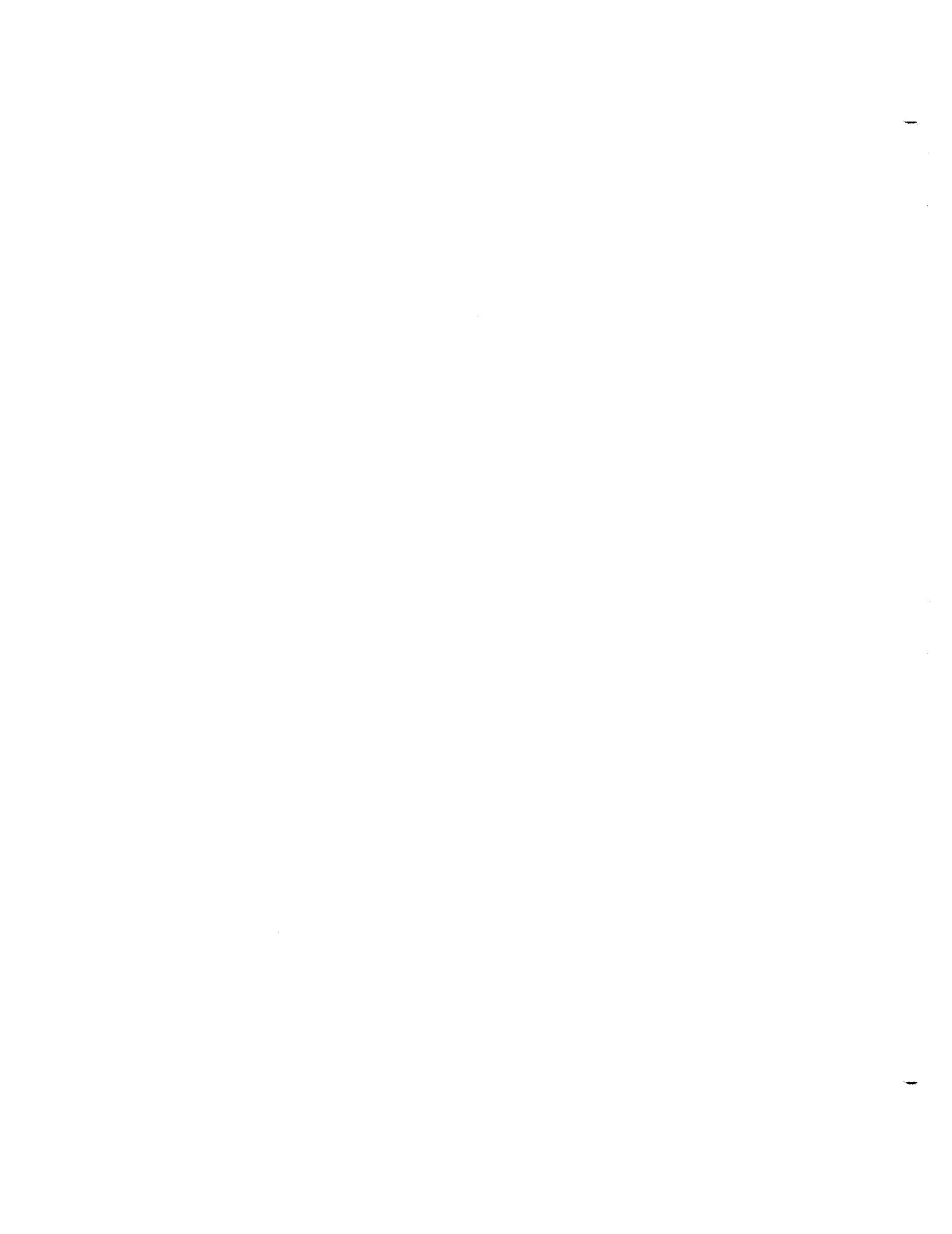
Fig. 5G. Output Summary (New Parameters).

```
ERROR - UNRECOGNIZED INPUT
CARD IMAGE IS 1 2 3 4 5 6 7 8
123456789012345678901234567890123456789012345678901234567890
ZILCH
```

Fig. 5H. Response to Unrecognized Command Card.

```
.....  
NORMAL STOP - ALL DATA PROCESSED  
-----  
IHC002I STOP      0  
-----  
.....
```

Fig. 5I. Normal Ending Message.



APPENDIX
PROGRAM LISTING

LEVEI 21.6 (DFC 72)

OS/360 FORTRAN H

DATE 74.3C4/09.17.44

```

      COMPILER OPTIONS - NAME= MAIN,OPT=U2,LINECNT=95,SIZE=0000K,
      SOURCE,EBLDIC,NULIST,ACDECK,LCAD,NOMAP,NGEDIT,NDIC,NOXREF
ISN 0002      IMPLICIT REAL*8 (A-H, O-Z)
ISN 0003      REAL*8 K,M
ISN 0004      REAL*4 HALL,FCUT,HCRH,HPRT,HCHA,HRES,FRUN,HBLK,FTEE,CARD
ISN 0005      DIMENSION A(20), B(5), C(50), CN(10), E(5), F(5), H(25),
1 K(25), F(20), R(10), T(20), W(5), CARC(20), VALU(7)
ISN 0006      DIMENSION C(50)
ISN 0007      EQUIVALENCE (A(1),V(1)), (R(1),V(21)), (C(1),V(26)),
1 (CN(1),V(76)), (E(1),V(86)), (F(1),V(91)),(H(1),V(96)),
2 (K(1),V(121)), (P(1),V(146)), (R(1),V(166)), (T(1),V(176)),
3 (W(1),V(196)), (M,V(201)), (L,V(202))
ISN 0008      EQUIVALENCE (CN(1),CF), (CN(2),CFF), (CN(3),CFT), (CN(4),CC),
1 (CN(5),CCF), (CN(6),CLT), (CN(7),CSG), (CN(8),CSS), (CN(9),CSR)
ISN 0009      COMMON/BLK2/TA, TOUT, IPR, KCLT, KPR
ISN 0010      COMMON /BLK1/ V(250)
ISN 0011      COMMON/BLK3/ IUIM(20), IUSE(20), NM(20), IBEG(20), NMCN(10),
1 NVAR,NCN
ISN 0012      DATA HALL/4HALL /,HOUT/4HOUTP/,FCRB/4FCRPF/,HPRT/4HPRIN/,
1 HCHA/4HCHAN/,HRES/4HRESF/,FRUN/4FRUN /,HBLK/4H /
ISN 0013      DATA XFCRG/2.300/,HFEE/4HF /
ISN 0014      DATA CTCL/1.0-7/, TTUL/1.0-7/
C
C      CTOL AND TTOL ARE THE CONVERGENCE TOLERANCES FOR CSELVE
C      AND TF RESPECTIVELY.
C
C      MAIN PROGRAM FOR CALCULATION OF MSER TRITIUM FLOW
C
C      SET UP REFERENCE VALUES IN WORKING ARRAY V
C
ISN 0015      CALL SETREF(PLK)
C
C      READ A CARD AND CHECK FOR INSTRUCTIONS
C
ISN 0016      100 READ(IN,END=997) CARD
ISN 0017      1 FORMAT(20A4)
ISN 0018      IF(CARD(1).NE.HOUT) GO TO 120
C
C      SET OUTPUT UNIT NUMBERS
C
ISN 0020      IF(CARD(3).NE.HALL) GO TO 115
ISN 0022      IF(CARD(4).NE.HCRH) GO TO 105
ISN 0024      104 KCUT=ICLT
ISN 0025      KPR=FCUT
ISN 0026      GO TO 100
ISN 0027      105 IF(CARD(4).EQ.HPRT) GO TO 110
ISN 0029      WRITE(KCUT,2)
ISN 0030      2 FORMAT(' CUTPUT NOT SPECIFIED CORRECTLY')
ISN 0031      WRITE(KCUT,20) (I,I=1,8),CARD
ISN 0032      20 FORMAT(' CARD IMAGE IS',1X,B(19X11)/15X,B(10H1234567890)/15X,20A4/
1 1X)
ISN 0033      WRITE(KCUT,21)
ISN 0034      21 FORMAT(' ALL OUTPUT TO SUMMARY UNIT')
ISN 0035      GO TO 104
ISN 0036      110 KCUT=IFP
ISN 0037      114 KPR=IPR
ISN 0038      GO TO 100
ISN 0039      115 KCUT=ICLT
ISN 0040      GO TO 114
C
C      CHECK FOR CHANGES IN WORKING VALUES
C
ISN 0041      120 IF(CARD(1).NE.HCHA) GO TO 135
ISN 0043      CALL MATCH(CARD(3),NM,NVAR,NK)
ISN 0044      IF(NK.NE.0) GO TO 125
ISN 0046      CALL MATCH(CARD(3),NMCN,NCN,NC)
ISN 0047      IF(NC.NE.0) GO TO 121
ISN 0049      WRITE(KCUT,4)
ISN 0050      4 FORMAT(' ERROR IN CHANGE SPECIFICATIONS')
ISN 0051      106 WRITE(KCUT,20) (I,I=1,8),CARD
ISN 0052      GO TO 100
ISN 0053      121 RFAC(IN,7) VALU(1)
ISN 0054      7 FORMAT(E10.0)
ISN 0055      J=IREG(4)+NC-1
ISN 0056      V(J)=VALU(1)
ISN 0057      WRITE(KCUT,12) CARD(3),VALU(1)
ISN 0058      12 FORMAT(1X,B4.1PE14.5)
ISN 0059      GO TO 100
ISN 0060      125 RFAC(TA,3) N1,N2,VALU
ISN 0061      3 FORMAT(2E3.7F10.0)
ISN 0062      IF(N1.FE.0) GO TO 100
ISN 0064      J=IREG(NK)-1+N1
ISN 0065      IF(N2.FE.0) N2=N1
ISN 0067      L=1
ISN 0068      DO 130 N=N1,N2
ISN 0069      V(J)=VALU(L)
ISN 0070      L=L+1
ISN 0071      J=J+1
ISN 0072      130 CONTINUE
ISN 0073      NV=N2-N1+1
ISN 0074      WRITE(KCUT,13) CARD(3),N1,N2,(VALU(L),L=1,NV)
ISN 0075      13 FORMAT(1X,B4,'(,12.,',12.)/'(1X1P5F14.5))

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```

ISN 0076      GC TC 125
C
C      CHECK FOR RESET - PUT REFERENCE VALUES BACK INTO V
C
ISN 0077      135 IF(CARD(1).NE.HRES) GO TO 150
ISN 0079      IF(CARD(3).NE.HBLK) GO TO 137
ISN 0081      136 CALL SETREF(CARD(3))
ISN 0082      GC TC 100
C
C      NAME NOT BLANK - TEST AGAINST NM ARRAY
C
ISN 0083      137 CALL MATCH(CARD(3).NM.NVAR.NK)
ISN 0084      IF(IK.NE.0) GC TO 136
C
C      NO MATCH FOUND IN NM - TRY NMCN
C
ISN 0086      CALL MATCH(CARD(3).NMCN.NCN.AC)
ISN 0087      IF(IC.NE.0) GC TO 136
ISN 0089      WRITE(KCUT,5)
ISN 0090      5 FORMAT(' ERROR IN RESET SPECIFICATIONS')
ISN 0091      GC TC 106
C
C      CHECK FOR PRINT
C
ISN 0092      150 IF(CARD(1).NE.HPRE) GO TO 155
ISN 0094      IF(CARD(3).NE.HBLK) GO TO 152
ISN 0096      151 CALL LCCK(CARD(3))
ISN 0097      GO TO 100
C
C      NAME NOT BLANK - TEST AGAINST NM ARRAY
C
ISN 0098      152 CALL MATCH(CARD(3).NM.NVAR.NK)
ISN 0099      IF(IK.NE.0) GC TO 151
C
C      NO MATCH FOUND IN NM - TRY NMCN
C
ISN 0101      CALL MATCH(CARD(3).NMCN.NCN.AC)
ISN 0102      IF(IC.NE.0) GC TO 151
ISN 0104      WRITE(KCUT,8)
ISN 0105      8 FORMAT(' ERROR IN PRINT SPECIFICATIONS')
ISN 0106      GO TC 106
ISN 0107      155 IF(CARD(1).EC.HRUN) GO TO 160
ISN 0109      WRITE(KCUT,6)
ISN 0110      6 FORMAT(' ERROR - UNRECOGNIZED INPLT')
ISN 0111      GO TC 106
ISN 0112      160 KTRY=1
ISN 0113      ISM=1
ISN 0114      NCC=0
ISN 0115      CCZ=0.00
C
C      CHECK FOR FAILURE TO SET CC AND CCT
C
ISN 0116      IF(CC.LE.0.00) CC=1.011
ISN 0118      IF(CCT.LE.0.00) CCT=1.010
C
C      PRINT WORKING ARRAY V ON LINE PRINTER
C
ISN 0120      KSAV=KCLT
ISN 0121      KOUT=IPR
ISN 0122      CALL NEWPG
ISN 0123      CALL LCCK(+BLK)
ISN 0124      CALL NEWPG
ISN 0125      KOUT=KSAV
ISN 0126      GC TC 200
C
C      GET C(18), CCF, Q(18) AND C(19)
C
C      FCS. 19A,B, 20, 21B
C
ISN 0127      200 C(18)=0.00
ISN 0128      CCF=0.00
ISN 0129      Q(18)=0.00
ISN 0130      Q(19)=0.00
ISN 0131      CALL NEWPG
ISN 0132      IF(R(5).EC.0.00) GO TO 205
ISN 0134      FF=F(4)*E(4)
ISN 0135      HA=H(18)*A(18)
ISN 0136      FEPH=HA+FE
ISN 0137      BW=R(4)*W(4)
ISN 0138      AFSC=(Bb*FEPH/(FE*H(18)))*2*K(18)
ISN 0139      ALFA=R(5)/FE
ISN 0140      C(18)=ALFA**2/(ALFA+.5DU*BESC*(1.00+DSORT(1.00+4.00*ALFA/BESQ)))
ISN 0141      Q(18)=Bb*A(4)*DSORT(C(18)*K(18))
ISN 0142      CCF=(R(5)+A*C(18))/FEPH
ISN 0143      Q(19)=CCF*FE
ISN 0144      205 CONTINUE
C
C      REGIA CALCULATION OF QUANTITIES DEPENDENT ON CC
C
C      GET C(10), C(11), Q(10), C(11)
C
C      FQS. 11A,B, 12A,B
C
ISN 0145      300 NCC=NCC+1
ISN 0146      IF(NCC.LE.50) GO TO 301
ISN 0148      WRITE(KCUT,35) HBLK

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ISN 0149      35 FORMAT(' FAILURE IN SOLUTION FOR CC'.A1)
ISN 0150      GC TO 100
ISN 0151      301 DO 305 I=1,C,11
ISN 0152      ALFA=K(I)*P(I)/(T(I)*H(I))*2
ISN 0153      CALL CCLAD(C,ALFA,C(1))
ISN 0154      Q(I)=P(I)*A(I)*DSQRT(K(I)*C(I))/T(I)
ISN 0155      305 CONTINUE
C
C      GET C(16), Q(16) - EDS. 17A,B
C
ISN 0156      BW=B(3)*H(3)
ISN 0157      ALFA=K(16)*P(H(16))*2
ISN 0158      CALL CCLAD(C,ALFA,C(16))
ISN 0159      Q(16)=BW*A(16)*DSQRT(K(16)*C(16))
C
C      GET C(17) - ED. 18
C
ISN 0160      Q(17)=F(3)*E(3)*CC
C
C      GET C(13), Q(13), C(14), C(14), C(15), Q(15)
C      ECS. 14A,B, 22A, 15A,B, 23A, 16A,B, 24A
C
ISN 0161      IGCFF=0
ISN 0162      CALL CSCOLVE(C,CSG,P(13),T(13),F(13),K(13),K(21),CTCL,
1 IGCFF,C(13),C(21))
ISN 0163      IF(IGCFF.LE.0) GO TO 320
ISN 0165      K1=13
ISN 0166      K2=21
ISN 0167      C2=CSG
ISN 0168      315 WRITE(KCLT,30) K1,K2
ISN 0169      30 FORMAT(' FAILURE IN SOLUTION FOR C('',12,'') AND C('',12,'')')
ISN 0170      IGCFF=-1
ISN 0171      CALL CSCOLVE(C,C2,P(K1),T(K1),F(K1),K(K1),K(K2),CTCL,
1 IGCFF,C(K1),C(K2))
ISN 0172      GO TO 100
ISN 0173      320 CALL CSCOLVE(C,CSS,P(14),T(14),F(14),K(14),K(22),CTCL,
1 IGCFF,C(14),C(22))
ISN 0174      IF(IGCFF.LE.0) GO TO 325
ISN 0176      K1=14
ISN 0177      K2=22
ISN 0178      C2=CSS
ISN 0179      GO TO 315
ISN 0180      325 CALL CSCOLVE(C,CSR,P(15),T(15),F(15),K(15),K(23),CTCL,
1 IGCFF,C(15),C(23))
ISN 0181      IF(IGCFF.LE.0) GO TO 330
ISN 0183      K1=15
ISN 0184      K2=23
ISN 0185      C2=CSR
ISN 0186      GO TO 315
C
C      C(13),Q(21),Q(14),Q(22),Q(15),Q(23)
C
ISN 0187      330 DO 335 I=13,15
ISN 0188      Q(I)=H(I)*A(I)*(CC-C(I))
ISN 0189      Q(18+I)=-C(I)
ISN 0190      335 CONTINUE
C
C      GET C(12), C(4) AND C(12) - EDS. 21A, 13A,B
C
ISN 0191      Q(12)=R(3)+R(4)-Q(10)-Q(11)-C(13)-C(14)-Q(15)-Q(16)-Q(17)
ISN 0192      Q(4)=-C(12)
ISN 0193      C(12)=C-C(12)/(H(12)*A(4))
ISN 0194      IF(C(12).GT.0.00) GO TO 340
ISN 0196      IC=12
ISN 0197      WRITE(KCLT,80) NCC,HBLK,CC,IC,C(FC)
ISN 0198      80 FORMAT(1X,I3,' CC',A1,' =',IPE14.5,' C('',12,'') =',E14.5)
C
C      IF C(12) IS NEGATIVE, ADJUST CC AND TRY AGAIN
C
ISN 0199      336 CC=CC+CC
C
C      SEE IF AN UPPER LIMIT (CC2) WAS FOUND.
C      IF SF, CC NOT EXCEED IT.
C
ISN 0200      IF(CC2.EQ.C.DO) GO TO 300
ISN 0202      IF(CC.LT.CC2) GO TO 300
ISN 0204      CC=0.5DC*(CC2+0.500*CC)
ISN 0205      GC TO 300
C
C      GET C(4) AND CF - EDS. 13A,B, 4A,B
C
ISN 0206      340 C(4)=(DSQRT(K(12)*C(12))-Q(12)*T(4)/(P(4)*A(4)))*2/K(4)
ISN 0207      CF=C(4)-C(12)/(H(4)*A(4))
ISN 0208      IF(CF.GT.0.00) GO TO 500
C
C      IF CF IS NEGATIVE, ADJUST CC AND TRY AGAIN
C
ISN 0210      WRITE(KCLT,81) NCC,HBLK,CC,HBLK,CF
ISN 0211      81 FORMAT(1X,I3,' CC',A1,' =',IPE14.5,' CF',A1,' =',E14.5)
ISN 0212      GO TO 336
ISN 0213      500 IF(KTRY.NE.1) GO TO 501
ISN 0215      RF1=0.00
ISN 0216      RF2=0.00

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ISN 0217      KTRY=2
ISN 0218      NCC=0
ISN 0219      WRITE(NCUT,50)HBLK
ISN 0220      50 FORMAT(' ITERATIVE SOLUTION FOR CC',A1/IX/
              1 ' ACC',6X,'CC1',11X,'CCL',11X,'CCX',11X,'RF1',11X,'CC2',11X)
C
C          BEGIN CALCULATION OF QUANTITIES DEPENDENT ON CF
C          GET C(1), C(2), C(3), Q(1), Q(2), C(3)
C          FOS, 1A,B, 2A,B, 3A,B
C
ISN 0221      501 DO 505 I=1,3
ISN 0222      ALFA=K(I)*P(I)/(T(I)*H(I))**2
ISN 0223      CALL CQUAD(CF,ALFA,C(I))
ISN 0224      Q(I)=P(I)*A(I)*DSORT(K(I)*C(I))/T(I)
ISN 0225      505 CONTINUE
C
C          GET C(5), C(5) - EQS. 5A,B
C
ISN 0226      RH=P(1)*W(1)
ISN 0227      ALFA=K(5)*(B/H(5))**2
ISN 0228      CALL CGLAD(CF,ALFA,C(5))
ISN 0229      Q(5)=RH*A(5)*DSORT(K(5)*C(5))
C
C          GET C(6) - EQ. 6
C
ISN 0230      Q(6)=F(1)*E(1)*CF
C
C          GET CFF - EQ. 10
C
ISN 0231      CFF=M*L*DSORT(K(5)*CF)/K(7)
C
C          GET C(7), C(7) - EQS. 7A,B
C
ISN 0232      RH=P(2)*W(2)
ISN 0233      ALFA=K(7)*(B/H(7))**2
ISN 0234      CALL CGLAD(CFF,ALFA,C(7))
ISN 0235      Q(7)=RH*A(7)*DSORT(K(7)*C(7))
C
C          GET C(8) - EQ. 8
C
ISN 0236      Q(8)=F(2)*E(2)*CFF
C
C          CALCULATE RF FOR EQ. 9
C
ISN 0237      RSUM=R(1)+R(2)
ISN 0238      RF=C(1)+C(2)+C(3)+Q(4)+Q(5)+C(6)+Q(7)+Q(8)-RSUM
ISN 0239      TE=RF/RSUM
C
C          TEST CONVERGENCE
C
ISN 0240      IF(DABS(TE)-LT,TTOL) GO TO 700
C
C          ACT CONVERGED - CHECK KTRY TO SEE WHAT NEXT
C
ISN 0242      IF(KTRY.NE.2) GO TO 540
C
C          KTRY=2 MEANS PRELIMINARY SEARCH FOR CC1 AND CC2
C          WHICH BRACKET THE ANSWER
C
ISN 0244      IF (RF.GT.0.00) GO TO 510
ISN 0246      CC1=CC
ISN 0247      RF1=RF
ISN 0248      WRITE(NCUT,55) NCC,CC1,RF1
ISN 0249      55 FORMAT(1X,13.1PE14.5,2BX,14.5)
ISN 0250      GO TO 515
ISN 0251      510 CC2=CC
ISN 0252      RF2=RF
ISN 0253      WRITE(NCUT,56) NCC,RF2,CC2
ISN 0254      56 FORMAT(1X,13.42X,1P2E14.5)
ISN 0255      515 IF(RF1*RF2) 530,520,516
C
C          RF1*RF2 POSITIVE SHOULD NEVER HAPPEN - SOMETHING WRONG
C
ISN 0256      516 WRITE(NCUT,51) HBLK
ISN 0257      51 FORMAT(' RF1*RF2 POSITIVE - SOMETHING FUELED UP IN CC',A1)
ISN 0258      GO TO 100
C
C          STILL LOCKING FOR ONE LIMIT - ADJUST CC AND TRY AGAIN
C          KEEP ADJUSTMENT FACTOR LESS THAN 100 AND GREATER THAN .01
C
ISN 0259      520 TFXCLG=TE*XCLG
ISN 0260      IF(DABS(TEXCLG).GT.4.600) TEXCLG=DSIGN(4.600,TEXCLG)
ISN 0262      ADJ=DLG(CC)-TEXCLG
ISN 0263      CC=CEXP(ADJ)
ISN 0264      GO TO 300
C
C          RF1*RF2 NEGATIVE - ANSWER BRACKETED
C
ISN 0265      530 KTRY=3
C
C          INVERSE LINEAR INTERPOLATION
C
ISN 0266      535 CCL=(RF1*CC2 - RF2*CC1)/(RF1-RF2)

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ISN 0267      KTRY=4
ISN 0268      IF (ISW.EC.2) GO TO 537
ISN 0270      CC=CCL
ISN 0271      GC TC 300
ISN 0272      537 CCT=CCL
ISN 0273      GC TC 705
ISN 0274      540 IF (KTRY.EC.3) GO TO 535
ISN 0276      IF (KTRY.NE.4) GO TO 555
C
C      KTRY=4 MEANS INVERSE LINEAR INTERPOLATION HAS BEEN
C      COMPLETED AND KF(CCL) CALCULATED
C
ISN 0278      WRITE(KCUT,52) MCC,CCL,CC,RF,CCZ
ISN 0279      52 FORMAT(1X,I3,1P2E14.5,10X,2E14.5)
ISN 0280      542 RFT=RF
C
C      INVERSE QUADRATIC INTERPOLATION
C
ISN 0281      D1=RF1-RFT
ISN 0282      D2=RF2-RFT
ISN 0283      D3=D2-D1
ISN 0284      CCX=CCL*RF1*RF2/(D1*D2)-CCL*RFT*RF2/(D1*D3)+CC2*RF1*RFT/(D2*D3)
ISN 0285      IF (CCX.LT.CC1) GO TO 545
ISN 0287      IF (CCX.GT.CC2) GO TO 545
ISN 0289      KTRY=5
ISN 0290      IF (ISW.EC.2) GO TO 544
ISN 0292      CC=CCX
ISN 0293      GC TC 300
ISN 0294      544 CCT=CCX
ISN 0295      GC TC 705
ISN 0296      545 IF (RFT.LT.D.CC) GO TO 550
ISN 0298      CC2=CCL
ISN 0299      RF2=RFT
ISN 0300      GO TO 530
ISN 0301      550 CC1=CC1
ISN 0302      RF1=RFT
ISN 0303      GC TC 530
C
C      KTRY=5 MEANS INVERSE QUADRATIC INTERPOLATION HAS BEEN
C      COMPLETED AND KF(CCX) CALCULATED
C
ISN 0304      555 IF (KTRY.NE.5) GO TO 585
ISN 0306      556 RFX=RF
ISN 0307      WRITE(KCUT,53) CCX,RFX
ISN 0308      53 FORMAT(32X,1F2E14.5)
C
C      TEST RFT AND RFX TO SEE WHAT NEW LIMITS ARE
C
ISN 0309      IF (RFX.GT.C.CC) GO TO 570
ISN 0311      IF (RFT.GT.D.CC) GO TO 565
ISN 0313      IF (DABS(RFX).GT.DABS(RFT)) GC TC 550
ISN 0315      560 CC1=CCX
ISN 0316      RFX=RFX
ISN 0317      GC TC 530
ISN 0318      545 CC2=CCL
ISN 0319      RF2=RFT
ISN 0320      GC TC 560
ISN 0321      570 IF (RFT.LT.D.CC) GO TO 580
ISN 0323      IF (RFX.GT.C.CC) GO TO 565
ISN 0325      575 CC2=CCX
ISN 0326      RFX=RFX
ISN 0327      GC TC 530
ISN 0328      580 CC1=CCL
ISN 0329      RF1=RFT
ISN 0330      GO TO 575
ISN 0331      585 WRITE(KCUT,54) KTRY
ISN 0332      54 FORMAT(1X,KTRY='*',I4/'* FRRDR*')
ISN 0333      GO TO 100
C
C      CALCULATION OF TRITIUM DISTRIBUTION - CCT SET BY
C      CHANGE INSTRUCTION
C
ISN 0334      700 KTRY=1
ISN 0335      ISW=2
ISN 0336      NCCT=0
C
C      GET C(40) AND U(41) - EQS. 34 AND 35
C
ISN 0337      705 NCCT=NCCT+1
ISN 0338      IF (NCCT.LE.50) GO TO 706
ISN 0340      WRITE(KCUT,35) HTEE
ISN 0341      GC TC 100
ISN 0342      706 RATIC=CCT/CC
ISN 0343      C(40)=RATIC*C(10)
ISN 0344      C(41)=RATIC*C(11)
ISN 0345      QSU=C(41)+Q(40)
C
C      GET C(43),C(44),C(45),U(43),C(44),C(45) - EQS. 37A,B, 38A,B, 39A,B
C
ISN 0346      DC 710 I=43,45
ISN 0347      J=1-30
ISN 0348      TC=T(J)*DCSRT(C(J)/K(J))
ISN 0349      HTCCP=TC*(H(J)/P(J))

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ISN 0350      C(I)=CCT*TCPP/11.00*H1LOP)
ISN 0351      Q(I)=F(I)*A(I)*C(I)/TC
ISN 0352      QSUM=C(I)+CSUM
ISN 0353      710 CONTINUE
C
C      GET C(46) THRU Q(49) - EQS. 40 THRU 43
C
ISN 0354      DO 715 I=46,49
ISN 0355      Q(I)=Q(I-30)*RATIO
ISN 0356      QSUM=C(I)+CSUM
ISN 0357      715 CONTINUE
C
C      GET C(42) AND Q(34) - EQ. 44 - MATERIAL BALANCE
C
ISN 0358      Q(42)=F(3)-CSUM
ISN 0359      Q(34)=-C(42)
C
C      CFT C(42) -EQ. 36A - CHECK FOR POSITIVE
C
ISN 0360      C(42)=CCT-C(42)/(H(4)*A(4))
ISN 0361      IF(C(42).GT.-C.D0) GO TO 725
ISN 0363      IC=42
ISN 0364      721 WRITE(KCUT,80) NCLT,HTEE,CCT,IC,C(I)
ISN 0365      720 CCT=CCT+CCT
ISN 0366      GO TO 705
C
C      C(42) POSITIVE - GET C(34) - EQ. 28B - CHECK FOR POSITIVE
C
ISN 0367      725 C(34)=(C(42)/DSORT(C(12)/K(12))-T(4)*C(42)/(P(4)*A(4)))*DSORT(C(4)
ISN 0368      S/K(4))
ISN 0370      IF(C(34).GT.-C.D0) GO TO 730
ISN 0371      IC=34
ISN 0371      GO TO 721
C
C      C(34) POSITIVE - GET CFT - EQ. 28A - CHECK FOR POSITIVE
C
ISN 0372      730 CFT=C(34)-C(42)/(H(4)*A(4))
ISN 0373      IF(CFT.GT.-C.D0) GO TO 735
ISN 0375      WRITE(KCUT,81) NCLT,HTEE,CCT,+TEE,CFT
ISN 0376      GO TO 720
C
C      CFT, C(34) AND C(42) ALL POSITIVE - CHECK KTRY
C      FOR NEXT STEP
C
ISN 0377      735 IF(KTRY.NE.1) GO TO 750
ISN 0379      KTRY=2
ISN 0380      RF1=0.C0
ISN 0381      RF2=0.C0
ISN 0382      NCC1=0
ISN 0383      WRITE(KCUT,50) HTEE
ISN 0384      750 RATIO=CFT/CF
ISN 0385      QSUM=0.C0
ISN 0386      DO 755 I=31,38
ISN 0387      TF([.EC.34) GO TO 755
ISN 0389      Q(I)=RATIO*Q(I-30)
ISN 0390      755 QSUM=C(I)*Q(I)
ISN 0391      RF=C(I)*R(I)
ISN 0392      TF=RF/R(I)
ISN 0393      IF(CABS(TF).LT.TTOL) GO TO 900
ISN 0395      IF(KTRY.NE.2) GO TO 790
ISN 0397      IF(RF.GT.-C.D0) GO TO 760
ISN 0399      CCI=CCT
ISN 0400      RF1=RF
ISN 0401      WRITE(KCUT,55) NCC1,CCI,RF1
ISN 0402      GO TO 765
ISN 0403      760 CC2=CCT
ISN 0404      RF2=RF
ISN 0405      WRITE(KCUT,56) NCC1,RF2,CC2
ISN 0406      765 IF(RF1*RF2) 530,770,766
ISN 0407      766 WRITE(KCUT,51) HTEE
ISN 0408      GO TO 100
C
C      STILL LOCKING FOR ONE LIMIT - ADJUST CCT AND TRY AGAIN
C      KEEP ADJUSTMENT FACTOR LESS THAN 100 AND GREATER THAN .01
C
ISN 0409      770 TFXCLG=TF*XCLG
ISN 0410      IF(DABS(TFXCLG).GT.4.6D0) TFXCLG=DSIGN(4.6D0,TFXCLG)
ISN 0412      ADJ=CLC(CCT)-TFXCLG
ISN 0413      CCT=DEXP(AEJ)
ISN 0414      GO TO 705
ISN 0415      790 IF(KTRY.EC.3) GO TO 535
ISN 0417      IF(KTRY.NE.4) GO TO 795
ISN 0419      WRITE(KCUT,52) NCC1,CCI,CCT,RF,CC2
ISN 0420      GO TO 542
ISN 0421      795 IF(KTRY.NE.5) GO TO 585
ISN 0423      GO TO 556
C
C      OUTPUT SECTION
C      SUMMARY OUTPUT TO UNIT KOUT
C
ISN 0424      900 CALL NEWPG
ISN 0425      WRITE(KCUT,92)
ISN 0426      52 FORMAT(' OUTPUT SUMMARY',IX)

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ISN 0427      QHTSS=C(13)*C(14)+Q(15)
ISN 0428      QTSS=C(43)+Q(44)+Q(45)
ISN 0429      QHSS=C(TSS-QTSS)
ISN 0430      RSUM=1.00/(R(1)+R(3))
ISN 0431      FRTSS=RSUM*QTSS
ISN 0432      WRITE(KCUT,93) QHTSS,QHSS,QHSS,FRTSS
ISN 0433      93 FORMAT(' STEAM SYSTEM',5X,'FLOW OF H + T INTO STEAM SYSTEM',
1 1PE16.5/5X,'FLOW OF T INTO STEAM SYSTEM',E20.5/5X,'FLOW OF H ',
2 'INTO STEAM SYSTEM',E20.5/5X,'FRACTION OF T INTO STEAM SYSTEM',
3 F16.5/1X)
ISN 0434      QHTFW=C(10)+C(11)
ISN 0435      QTPW=C(40)+C(41)
ISN 0436      FRTPW=CTPW*RSUM
ISN 0437      FTSSK=C(46)*RSUM
ISN 0438      FRTSF=C(48)*RSUM
ISN 0439      FRRFT=C(47)*RSUM
ISN 0440      FRRPF=C(49)*RSUM
ISN 0441      WRITE(KCUT,94) Q(4),Q(34),QHTFW,QTPW,C(16),Q(46),Q(18),C(48),
1 Q(17),C(47),C(19),Q(49),FRTPW,FTSSK,FRTSF,FRRPT,FRRPF
ISN 0442      WRITE(KCUT,95) CC,CCT,CLF
ISN 0443      94 FORMAT(' SECONDARY SYSTEM',5X,'FLOW OF H + T INTO SECONDARY ',
1 'FROM PRIMARY',1PE14.5/5X,'T INTO SECONDARY FROM PRIMARY',E18.5/
2 5X,'H + T THRU PIPE WALLS INTO CELLS',E15.5/5X,'T THRU PIPE ',
3 'WALLS INTO CELLS',E19.5/5X,'SORPTION BY SINK',5X,'H + T',2BX,
4 F14.5/5X,'T',32X,E14.5/5X,'HF',31X,E14.5/5X,'TF',31X,E14.5/
5 'REMOVAL BY PURGE',5X,'H + T',2EX,E14.5/5X,'T',32X,E14.5/5X,
6 'HF',31X,E14.5/5X,'TF',31X,E14.5/5X,'FRACTION OF T',5X,'PASSING',
7 'THRU PIPE WALLS',10X,E14.5/5X,'SORBED BY SINK AS T',14X,E14.5/
8 5X,'SORBED BY SINK AS TF',13X,E14.5/5X,'REMOVED BY PURGE AS T',
9 12X,E14.5/5X,'REMOVED BY PURGE AS TF',11X,E14.5)
ISN 0444      95 FORMAT(' CONCENTRATIONS IN SECONDARY SALT',5X,'H + T (CC)',23X,
1 1PE14.5/5X,'T (CLT)',26X,E14.5/5X,'HF (CCF)',25X,E14.5/1X/1X)
ISN 0445      QHTW=C(11)+C(12)+Q(3)
ISN 0446      QTW=C(21)+C(32)+Q(33)
ISN 0447      FRTWC=CTW*RSUM
ISN 0448      FTSSK=C(35)*RSUM
ISN 0449      FTSSSF=C(37)*RSUM
ISN 0450      FTRPT=C(36)*RSUM
ISN 0451      FTRPTF=C(38)*RSUM
ISN 0452      WRITE(KCUT,96) QHTW,QTW,Q(5),C(35),Q(7),Q(37),Q(6),Q(36),Q(8),
1 C(38),FRTWC,FTSSK,FTSSSF,FTRPT,FTRPTF
ISN 0453      WRITE(KCUT,97) CF,CFT,CFP
ISN 0454      96 FORMAT(' PRIMARY SYSTEM',5X,'H + T THRU WALLS INTO ',
1 'CELL',1PE21.5/5X,'T THRU WALLS INTO CELL',11X,E14.5/5X,'SORPTION
2 BY SINK',5X,'H + T',2BX,E14.5/5X,'T',32X,E14.5/5X,'HF',31X,E14.5/
3 5X,'TF',31X,E14.5/5X,'REMOVAL BY PURGE',5X,'H + T',2BX,E14.5/5X,
4 'T',32X,E14.5/5X,'HF',31X,E14.5/5X,'TF',31X,E14.5/5X,'FRACTION OF
5 T',5X,'PASSING THRU WALLS INTO CELL',E19.5/5X,'SORBED BY SINK AS ',
6 'T',14X,E14.5/5X,'SORBED BY SINK AS TF',13X,E14.5/5X,'REMOVED ',
7 'BY PURGE AS T',12X,E14.5/5X,'REMOVED BY PURGE AS TF',11X,E14.5)
ISN 0455      97 FORMAT(' CONCENTRATIONS IN PRIMARY SALT',5X,'H + T (CF)',23X,
1 1PE14.5/5X,'T (CLT)',27X,E14.5/5X,'HF (CFP)',25X,E14.5)
ISN 0456      CALL NF6PG
ISN 0457      GO TO 100
C
C
C
END OF FILE DETECTED ON INPUT UNIT
ISN 0458      957 CALL NF6PG
ISN 0459      WRITE(KCUT,99)
ISN 0460      99 FORMAT(' NORMAL STOP - ALL DATA PROCESSED')
ISN 0461      STOP
ISN 0462      END

```

OPTICS IN FFFCT NAME= DATA,OPT=02,LINECNT=95,SIZE=CCCCC.

OPTICS IN FFFCT SOURCE=EBCCCIC,NOLIST,NOCHECK,LCAC,NCMAP,NOEDIT,NOID,NOXREF

STATISTICS SOURCE STATEMENTS = 401 ,PROGRAM SIZE = 10188

STATISTICS NO DIAGNOSTICS GENERATED

***** END OF COMPILATION *****

45K BYTES OF CORE NOT USED

LEVEL 21.6 (DFC 72)

OS/360 FORTRAN H

DATE 74.304/09.19.05

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      COMPILER OPTIONS - NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K,
                        SOURCE,EBCDIC,NOLIST,NODECK,LCAC,NOMAP,NOEDIT,NCID,NOXREF
ISN 0007      BLOCK DATA
ISN 0003      COMMON /BLK3/ IDIM(20), IUSE(20), NM(20), IBEG(20), NMCM(10),
& NVAR, NCN
ISN 0004      DATA IDIM / 20,5,50,10,5,5,25,25,20,10,20,5,1,1,6*0/
ISN 0005      DATA IUSE / 18,4,43, 9,4,4,23,23,15, 5,15,4,1,1,6*0/
ISN 0006      DATA IBEG / 1,21,26,76,86,91,56,121,146,166,176,156,201,202,6*0/
ISN 0007      DATA NM/4HA ,4HB ,4HC ,4CN ,4FF ,4HF ,4HH ,
&      4HK ,4HP ,4HK ,4HT ,4TW ,4HM ,4HU /
ISN 0008      DATA NVAR, NCN / 14, 9 /
ISN 0009      DATA NCM / 4HCF ,4HCF ,4HCFT ,4CC ,4HCCF ,4HCCT ,4HCSG ,
&      4HCSS ,4HCSK ,4H /
ISN 0010      COMMON /BLK2/ IN, IOUT, IPR, KCUT, KPR
ISN 0011      DATA IN, IOUT, IPR /5, 20, 6/
      C
      C      IN - INPUT UNIT NO.
      C      IOUT - AUXILIARY OUTPUT UNIT NO.
      C      IPR - LINE PRINTER UNIT NO.
      C
ISN 0012      END

*OPTIONS IN EFFECT*      NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K,
*OPTIONS IN EFFECT*      SOURCE,EBCDIC,NOLIST,NODECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
*STATISTICS*      SOURCE STATEMENTS =      11 ,PROGRAM SIZE =      8
*STATISTICS*      NO DIAGNOSTICS GENERATED
***** END OF COMPILATION *****
125K BYTES OF CORE NOT USED

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LEVEL 71.6 (DEC 72)

OS/360 FORTRAN M

DATE 74.304/09.19.12

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COMPILER OPTIONS - NAME= MAIN,OPT=02,LINECAT=95,SIZE=0000K,
SOURCE,EBODIC,NULIST,ACCECK,LCAC,NOMAP,NOEDIT,NOIG,NOXREF
TSN 0007 SUBROUTINE SETREF(NAME)
C SETS VARIABLES TO THEIR REFERENCE VALUES. IF NAME IS BLANK,
C ALL VARIABLES AS SPECIFIED IN THE ARRAY NM WILL BE SET.
C IF NAME IS PNT, ALL VARIABLES IN THE ARRAY VREF
C WILL BE PRINTED.
C
TSN 0003 IMPLICIT REAL*8 (A-H,O-Z)
TSN 0004 REAL*4 VWR,VW,WDRD
TSN 0005 CCMPCA/BLK3/ I DIM(20), IUSE(20), NM(20), IBEG(10), NMCN(10),
1 NVAR,ACN
TSN 0006 DIMENSION VREF(250), VQ1(155), VQ2(75)
TSN 0007 EQUIVALENCE (VQ1(1),VREF(96)), (VQ2(1),VREF(176))
C
C VQ1 AND VQ2 ARE DUMMY ARRAYS USED IN THE INITIALIZATION OF
C PARTS OF VREF, THE REFERENCE ARRAY.
C
TSN 0008 DATA (RLNK/4H /,IPRT/4HPRNT/
TSN 0009 DATA VWRD/4HVREF/, VW/4HV /
TSN 0010 CCMPCN/BLK2/IN, IOUT, IPR, KCLT, KPR
TSN 0011 CCMPCA /BLK1/ V(250))
C
C THE COMMENT CARDS INTERSPERSED AMONG THE FOLLOWING
C CONTINUATION CARDS CAUSE NO TROUBLE WITH THE ORNL COMPILER.
C THIS IS CONTRARY TO THE RULE ON PG. 12, GC28-6515-8,
C IBM SYSTEM/360 AND SYSTEM/370 FORTRAN IV LANGUAGE.
C
TSN 0012 DATA VREF
C REFERENCE VALUES FOR A(I)
C
1 /,6D6+.5D6+3.5D6+.49D6+.52D6+-.1D0+.52D6+.2*-1D0 +11.D6+.8D6+
2 49.D6+.31.D6+.27.D6+.18.D6+.0.D6+-.1D0+.0D0+.2*-1.DC ,
C
C REFERENCE VALUES FOR B(I)
C
2 3.021,3.021,1.018,1.018+-.1CC,
C
C REFERENCE VALUES FOR C(I) ARE ALL ZERO
C
3 45*0.DC +5*-1.D0 ,
C
C REFERENCE VALUES FOR CN(I)
C
4 6*-1.DC +2.D10+.9.D11+1.D11+-.1CC,
C
C REFERENCE VALUES FOR E(I)
C
5 .500,.(17D0+.18D0+.0018D0+-.1D0,
C
C REFERENCE VALUES FOR F(I)
C
6 3.6D5,3.6C5,5.05,5.05+-.1D0/
C
C REFERENCE VALUES FOR H(I)
C
TSN 0013 DATA VC1/
1 1.6D-2,6.C-3,9.D-5,1.9U-2,3.C-3,-1.CC,3.D-3,2*-1.DC ,
2 7.4D-2,3.4D-2,9.7D-2,4.3D-2,4.7D-2,4.0D-2,.8D0+-.1D0+.8D0,
3 2*-1.DC +5.8D0+12.D0+30.CC+2*-1.DC ,
C
C REFERENCE VALUES FOR K(I)
C
4 1.2D-17,2.D-17,1.5D-17,1.5C-17,1.4D-17,-1.D0,1.5D-19,2*-1.D0,
5 3.4C-18,5.D-18,4.0-18,4.5C-18,3.5D-18,4.0-18,4.4D-18,-1.DC,
X 1.1D-20,2*-1.D0+4.5D-20,5.1C-20,4.8C-20,2*-1.DC,
C
C REFERENCE VALUES FOR P(I)
C
62.1D15,6.7C14,9.0D14,9.0D14,5*-1.CC,1.1D15,1.8D14,9.0D14,1.8D14,
76.7C14,3.5C14,5*-1.D0,
C
C REFERENCE VALUES FOR R(I)
C
8 5.6D17,4*0.CC +5*-1.D0 /
C
C REFERENCE VALUES FOR T(I)
C
TSN 0014 DATA VC2/
1 2*13.D0 +50.D0,1.00,5*-1.00 +2*13.CC +1.D0,2*2.DC +1.CC,5*-1.00,
C
C REFERENCE VALUES FOR W(I)
C
2 4*1.D0 +-.1.CC,
C
C REFERENCE VALUE FOR M
C
31.12D-6+,
C
C REFERENCE VALUE FOR U
C
4 1.D2/

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C
C      CHECK NAME
C
ISN 0015      IF(NAME.EQ.IBLNK) GO TO 102
ISN 0017      IF(NAME.EC.IFRT) GO TO 115
C
C      LCKK FOR MATCH IN NM TABLE
C
ISN 0019      CALL MATCH(NAME,NM,NVAR,NK)
ISN 0020      IF(NK.NE.0) GO TO 101
C
C      NO MATCH FOUND - PRINT MESSAGE
C
C      CHECK NAME AGAINST CN TABLE
C
ISN 0022      CALL MATCH(NAME,NMCN,NCN,KC)
ISN 0023      IF(KC.EC.0) GO TO 300
ISN 0025      J=IBEG(4)+KC-1
ISN 0026      V(J)=VREF(J)
ISN 0027      RETLRA
ISN 0028      300 WRITE(KCLT,1) NAME
ISN 0029      1  FORMAT(1X/' NO MATCH FOR ',A4,' IN SETREF - NO CHANGE IN V')
ISN 0030      RETLRA
C
C      NAME = NM(NK)
C
ISN 0031      101 N1=NK
ISN 0032      N2=NK
ISN 0033      GC TC 103
ISN 0034      102 N1=1
ISN 0035      N2=NVAR
ISN 0036      103 DO 110 N=N1,N2
ISN 0037      IT=ILSE(N)
ISN 0038      J=IBEG(A)
ISN 0039      DO 105 I=1,IT
ISN 0040      V(J)=VREF(J)
ISN 0041      105 J=J+1
ISN 0042      110 CONTINUE
ISN 0043      RETLRA
C
C      PRINT ALL REFERENCE VALUES
C
ISN 0044      115 LINES=5C
ISN 0045      WORD=VWORD
ISN 0046      DO 130 N=1,NVAR
ISN 0047      IT=ILSE(N)/5+4
ISN 0048      IF((LINES+IT).LE.50) GO TO 120
ISN 0050      CALL NEWPG
ISN 0051      WRITE(KCLT,2) WORD
ISN 0052      2  FORMAT(' VALUES IN ARRAY ',A4)
ISN 0053      LINES=C
ISN 0054      120 WRITE(KCLT,3) NM(N),IDIM(N),ILSE(N),IBEG(N),(I,I=1,5)
ISN 0055      3  FORMAT(1X/' NAME ',A4,' DIMENSION',I4,' USED',I4,' STARTS AT',I4/
      1 7X,4(1I,13X),1I/1X)
ISN 0056      J1=IBEG(A)
ISN 0057      I1=J1
ISN 0058      J2=[1+ILSE(N)-1
ISN 0059      DO 125 J=J1,J2,5
ISN 0060      I2=PTNO(I1+J2)
ISN 0061      WRITE(KCLT,4) (VREF(I),I=I1,I2)
ISN 0062      4  FORMAT(1X,1PSE14,5)
ISN 0063      I1=I2+1
ISN 0064      125 CONTINUE
ISN 0065      LINES=LINES+IT
ISN 0066      130 CONTINUE
ISN 0067      RETLRA
ISN 0068      ENTRY LCKK
ISN 0069      WORD=VW
C
C      PRINTS VALUES IN V SELECTED BY NAME. IF NAME
C      IS BLANK, PRINT ALL.
C
C      CHECK NAME
C
ISN 0070      IF(NAME.EC.IBLNK) GO TO 132
C
C      LCKK FOR MATCH IN TABLE
C
ISN 0072      CALL MATCH(NAME,NM,NVAR,NK)
ISN 0073      IF(NK.NE.0) GO TO 201
C
C      CHECK NAME AGAINST CN TABLE
C
ISN 0075      CALL MATCH(NAME,NMCN,NCN,KC)
ISN 0076      IF(KC.EC.0) GO TO 301
ISN 0078      J=IBEG(4)+KC-1
ISN 0079      WRITE(KCLT,6) NMCN(KC), J, V(J)
ISN 0080      6  FORMAT(1X/1X,A4,' =V(',I3,')', VALUE =',1PE14,5)
ISN 0081      RETLRA
ISN 0082      301 WRITE(KCLT,5) NAME
ISN 0083      5  FORMAT(1X/' NO MATCH FOUND FOR ',A4,' IN LOOK - NO PRINT')
ISN 0084      RETLRA

```

```

1SN 0085      2C1 N1=AK
1SN 0086      N2=AK
1SN 0087      GO TC 135
1SN 0088      132 N1=1
1SN 0089      N2=AVAR

C
C          PRINT SELECTED VALUES IN V
C
1SN 0090      135 LINES=50
1SN 0091      DO 150 N=N1,N2
1SN 0092      IT=LUSE(N)/5+4
1SN 0093      IF(LINES+IT.LE.50) GO TO 140
1SN 0094      CALL AFKPG
1SN 0095      WRITE(KCLT,2) WORD
1SN 0096      LINES=0
1SN 0097
1SN 0098      140 WRITE(KCUT,3) NMIN(),IDIM(N),LUSE(N),IBEG(N),(I,I=1,5)
1SN 0099      J1=IBEG(N)
1SN 0100      I1=J1
1SN 0101      J2=I1+LUSE(N)-1
1SN 0102      DO 145 J=J1,J2,5
1SN 0103      I2=I1+5-J2
1SN 0104      WRITE(KCUT,4) (V(I),I=I1,I2)
1SN 0105      I1=I2+1
1SN 0106      145 CONTINUE
1SN 0107      LINES=LINES+IT
1SN 0108      150 CONTINUE
1SN 0109      RETURN
1SN 0110      END

```

OPTIONS IN EFFECT NAME= PA1N.UPT=Q2.LINECNT=55.SIZF=0000K.

OPTIONS IN EFFECT SOURCE.EBCDIC.NOLIST.NUDECK*.LCAD.NCMAP.NOEDIT.NOID.ACXREF

STATISTICS SOURCE STATEMENTS = 109 .PROGRAM SIZE = 3996

STATISTICS NO DIAGNOSTICS GENERATED

***** END OF COMPILATION *****

105K BYTES OF CORE NET USED

LFVFI 21.6 (DFC 72)

05/360 FORTRAN H

DATE 74.304/09.19.44

```

COMPILER OPTIONS - NAME= MAIN,OPT=D2,LINECNT=95,SIZE=0000K,
SOURCE,EBGDI,C,NOLIST,ACDFCK,LCAD,NOMAP,NDEFIT,NOID,NOXREF
ISN 0002 SUBROUTINE CSOLVE(C1,C2,PK,TK,PK,HL,KK,KL,EPS,IGDCF,CK,CL) CSOL 100
ISN 0003 IMPLICIT REAL*8 (A-H,O-Z)
ISN 0004 REAL*8 KK, KL
ISN 0005 TEST=100.00
ISN 0006 ASSIGN 55 TO K CSOL 110
ISN 0007 MKHL=PK/HL CSOL 115
ISN 0008 PCTH=PK/(TK*MK) CSOL 122
ISN 0009 CKMAX=C1+C2/TK*MKHL
C CSOL 130
C CHECK FOR SOLUTION CSOL 131
C CSOL 132
C CSOL 133
ISN 0010 CK2=CKMAX
ISN 0011 G2=CK2-C1+PCTH*DSQRT(KK*CK2)
ISN 0012 IF(G2.GT.0.) GO TO 85 CSOL 135
ISN 0014 ASSIGN 130 TO K CSOL 136
ISN 0015 PRINT 4 CSOL 137
ISN 0016 4 FORMAT('1NC SCLUTION') CSOL 138
ISN 0017 IGCF=1 CSOL 139
ISN 0018 GO TO 90 CSOL 140
ISN 0019 85 IF(IGCF.EQ.0) GO TO 95 CSOL 141
C CSOL 142
C HEADING FOR DEBUG PRINTOUT CSOL 145
C CSOL 150
C CSOL 155
ISN 0021 90 PRINT 1, C1, C2, PK, TK, MK, HL, KK, KL, EPS CSOL 155
ISN 0022 1 FORMAT('1CSOLVE ARGUMENTS*/1C.6X.*C1*.12X.*C2*.12X.*PK*.12X.*TK*. CSOL 160
X 12X.
1 'MK*.12X.*HL*.12X.*KK*.12X*KL*.12X*EPS*/IX,1P9E14.5/'C IT*.6X, CSOL 161
2 'C.K1*.11X.*CKT*.11X.*CX*.12X.*CK2*.10X.*TEST*/1X) CSOL 162
ISN 0023 GO TO K, (95,130) CSOL 163
C CSOL 165
C START ITERATIONS FOR CK CSOL 170
C CSOL 175
C CSOL 176
ISN 0024 55 CK1=0.
ISN 0025 CKCLO=CK1
ISN 0026 CL1=KL*(MKHL*(C1+C2))
ISN 0027 G1=-(C1+PCTH*DSQRT(CL1))
ISN 0028 DO 115 IT=1,50 CSOL 180
C
C INVERSE LINEAR INTERPOLATION AND CHECK FOR CONVERGENCE
C
ISN 0029 CKT=(CK1*G2-CK2*G1)/(G2-G1) CSOL 182
ISN 0030 1F(ABS(CKCLO/CKT-1.00).LT.EPS) GO TO 120 CSOL 183
C
C IF NOT CONVERGED, TRY INVERSE QUADRATIC INTERPOLATION
C
ISN 0032 CLT=KL*(MKHL*(C1-CKT)+C2)
ISN 0033 FF=C1+PCTH*(DSQRT(CL1)-DSQRT(KK*CKT)) CSOL 192
ISN 0034 G=CKT-FF CSOL 193
ISN 0035 DG1=G-G1 CSOL 194
ISN 0036 DG2=G-G2 CSOL 195
ISN 0037 DG3=DG2-DG1 CSOL 196
ISN 0038 CX=-G*G2*CK1/(DG3*DG1)+G1*G2*CKT/(EG1*DG2)+G1*G*CK2/(DG3*DG2) CSOL 240
ISN 0039 1F(IGCF.EQ.0) GO TO 200 CSOL 245
C CSOL 250
C CSOL 255
ISN 0041 PRINT 2,IT,CK1,CKT,CX,CK2,TEST
ISN 0042 2 FORMAT('1X3,1P5F14.5)
ISN 0043 200 IF(CX.LE.CK1) GO TO 102 CSOL 197
ISN 0045 IF(CX.GE.CK2) GO TO 102 CSOL 198
ISN 0047 TEST=DABS(CKCLO/CX-1.00)
ISN 0048 1F(TEST.EQ.EPS) GO TO 201
ISN 0050 CKT=CX
ISN 0051 GO TO 120
ISN 0052 201 CLX=KL*(MKHL*(C1-CX)+C2)
ISN 0053 FX=C1+PCTH*(DSQRT(CLX)-DSQRT(KK*CX)) CSOL 201
ISN 0054 GX=CX-FX CSOL 202
ISN 0055 CKCLO=CX CSOL 203
ISN 0056 1F(GX.LT.0.) GO TO 101 CSOL 204
ISN 0058 CK2=CX CSOL 206
ISN 0059 G2=GX CSOL 207
ISN 0060 GO TO 115 CSOL 208
ISN 0061 101 CK1=CX CSOL 225
ISN 0062 G1=GX CSOL 226
ISN 0063 GO TO 105 CSOL 227
ISN 0064 102 TEST=DABS(CKCLO/CKT-1.00) CSOL 230
ISN 0065 CKCLO = CKT CSOL 231
ISN 0066 1F(TEST.LE.EPS) GO TO 120 CSOL 232
ISN 0068 1F (G .EQ. 0.00) GO TO 120 CSOL 233
ISN 0070 1F(G.GT.0.) GO TO 105
ISN 0072 G1=G
ISN 0073 CK1=CKT
ISN 0074 GO TO 115
ISN 0075 105 G2=G CSOL 235
ISN 0076 CK2=CKT CSOL 236
ISN 0077 115 CONTINUE CSOL 275
C CSOL 280
C PRINT NON-CONVERGENCE ALARM CSOL 285
C CSOL 290
ISN 0078 PRINT 3 CSOL 290
ISN 0079 3 FORMAT('0 CSOLVE UNABLE TO FIND ROOT') CSOL 295

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```
ISN 00A0      TGOCF=PAK0(1,IGUOF+1)      CSOL 300
ISN 00B1      RETURN                      CSOL 305
ISN 00B2      120 CK=CKT                  CSOL 310
ISN 00B3      CL=PKCL*(L1-CK)+C2
ISN 00B4      130 RFLRA                      CSOL 320
ISN 00B5      END                          CSOL 325

*OPTIONS IN EFFECT*      NAME= MAIN,OPT=02,LINFCNT=95,SIZE=CCCOK,
*OPTIONS IN EFFECT*      SCURCE,EBCDIC,NOLIST,NODECK,LOCAC,NCMAP,NOEDIT,NOID,NOXREF
*STATISTICS*      SOURCE STATEMENTS =      84 ,PROGRAM SIZE =      2024
*STATISTICS*      NO DIAGNOSTICS GENERATED
***** END OF COMPILATION *****      121K BYTES OF CORE NOT USED
```

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LEVEL 21.6 (DFC 72)                                CS/360 FORTRAN H                                DATE 74.304/09.20.10

      COMPILER OPTIONS - NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K,
                        SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
ISN 0002      SUBROUTINE NEWPG                                NEWP 100
      C                                                    NEWP 105
      C      IF THE SUMMARY OUTPUT UNIT (KOUT) IS THE LINE PRINTER (IPR), NEWP 110
      C      EJECT TO A NEW PAGE. IF NOT, PRINT 5 BLANK LINES. NEWP 115
      C                                                    NEWP 120
ISN 0003      COMMON /BLK2/ IN, IOUT, IPR, KCUT, KPR        NEWP 125
ISN 0004      IF (KOUT.EC.IPR) GO TO 100                    NEWP 130
ISN 0006      WRITE(KCUT,1)                                  NEWP 135
ISN 0007      1 FORMAT(1X/1X/1X/1X/1X)                      NEWP 140
ISN 0008      RETURN                                         NEWP 145
ISN 0009      100 WRITE(KCUT,2)                              NEWP 150
ISN 0010      2 FORMAT(1H)                                    NEWP 155
ISN 0011      RETURN                                         NEWP 160
ISN 0012      END                                            NEWP 165

*OPTIONS IN FFFFT*      NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K.
*OPTIONS IN FFFFT*      SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
*STATISTICS*      SOURCE STATEMENTS =      11 ,PROGRAM SIZE =      260
*STATISTICS*      NO DIAGNOSTICS GENERATED
***** END OF COMPILATION *****                                133K BYTES OF CORE NOT USED

```

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LEVEL 21.6 (DFC 72)                                CS/360 FORTRAN H                                DATE 74.304/09.20.25

      COMPILER OPTIONS - NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K,
                        SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
ISN 0002      SUBROUTINE CQUAD(C1,ALFA,CK)
      C                                                    CQUA 105
      C      SOLVES THE QUADRATIC EQUATION                    CQUA 110
      C                                                    CQUA 115
      C       $CK**2 - (2*C1 + ALFA)*CK + C1**2 = 0$ 
      C                                                    CQUA 125
      C      FOR THE ROOT CK WHICH IS LESS THAN C1/ALFA.    CQUA 130
      C                                                    CQUA 135
ISN 0003      IMPLICIT REAL*8 (A-H,O-Z)
ISN 0004      REAL*8 KK                                       CQUA 140
ISN 0005      T=C1/ALFA                                       CQUA 145
ISN 0006      CK=C1**2/(C1+.5*ALFA*(1.+DSQRT(4.*C1/ALFA+1.)))
ISN 0007      RETURN                                         CQUA 155
ISN 0008      END                                            CQUA 160

*OPTIONS IN FFFFT*      NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K.
*OPTIONS IN FFFFT*      SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
*STATISTICS*      SOURCE STATEMENTS =      7 ,PROGRAM SIZE =      444
*STATISTICS*      NO DIAGNOSTICS GENERATED
***** END OF COMPILATION *****                                133K BYTES OF CORE NOT USED

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LEVEL 21.6 (DFC 72)                                CS/360 FORTRAN H                                DATE 74.304/09.20.38

      COMPILER OPTIONS - NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K,
                        SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
ISN 0002      SUBROUTINE MATCH (NAME, NTAB, NN, NMAT)
      C
      C      SEARCHES THE ARRAY NTAB WITH NN ITEMS FOR THE FIRST
      C      OCCURRENCE OF NAME. IF NAME IS NOT FOUND IN NTAB, NMAT
      C      IS RETURNED WITH THE VALUE ZERO.
      C
ISN 0003      DIMENSION NTAB (10)
ISN 0004      NMAT = 0
ISN 0005      DO 100 N=1,NN
ISN 0006      IF (ICMPA(NAME, NTAB)(N, 4) .NE. 0) GO TO 100
ISN 0008      NMAT = N
ISN 0009      RETURN
ISN 0010      100 CONTINUE
ISN 0011      RETURN
ISN 0012      END

*OPTIONS IN FFFFT*      NAME= MAIN,OPT=02,LINECNT=95,SIZE=0000K.
*OPTIONS IN FFFFT*      SOURCE,EBCCDIC,NOLIST,ACDECK,LCAC,NOMAP,NOEDIT,NOID,NOXREF
*STATISTICS*      SOURCE STATEMENTS =      11 ,PROGRAM SIZE =      398
*STATISTICS*      NO DIAGNOSTICS GENERATED
***** END OF COMPILATION *****                                133K BYTES OF CORE NOT USED
STATISTICS*      NO DIAGNOSTICS THIS STEP

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