

**WHIRLWIND PROGRAMMING OF S_2
APPROXIMATION FOR FLUX DISTRIBUTION
IN A FINITE CYLINDRICAL
REACTOR**

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42
Cambridge 39, Massachusetts

27 May 1957

Secretary of the Faculty
Massachusetts Institute of Technology
Cambridge 39, Massachusetts

Dear Sirs:

We hereby submit our thesis, Whirlwind Programming of S_2 Approximation for Flux Distribution in a Finite Cylindrical Reactor, in partial fulfillment for the degree of Naval Engineer and the degree of Master of Science in Naval Architecture and Marine Engineering.

Whirlwind Programming
of S_2 Approximation for Flux Distribution
in a Finite Cylindrical Reactor

by

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submitted in partial fulfillment of the requirements for the degree of
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Submitted to the Department of Naval Architecture and Marine Engineering on May 27, 1957 in partial fulfillment of the requirements for the degree of Naval Engineer and the degree of Master of Science in Naval Architecture and Marine Engineering.

ABSTRACT

This thesis attempts to develop a program for the Whirlwind computer operating in the interpretive mode for the flux distribution in a finite cylindrical reactor.

The reactor chosen for investigation was a hypothetical organic moderated, highly enriched, cylindrical reactor having a power output capable of fulfilling one half the propulsive requirements of a nuclear powered naval cruiser.

Heat transfer and basic physics calculations were accomplished to determine preliminary size and material requirement of such a reactor.

The flux equations were derived using numerical integration methods suitable for computer programming.

Conclusion

- 1) The code as presently programmed is incorrect.
- 2) The grid spacing as used is much too coarse. This difficulty is not inherent in the coded program.
- 3) The mode of computer operation is unsuitable for an iterative process such as this because it results in a code that progresses too slowly.

Recommendations

- 1) Prior to the use of any computer a formal or semi-formal course in programming should be completed.
- 2) A grid spacing comparable to the mean free path of the neutron should be selected.
- 3) The fastest possible mode of calculation on the computer available should be used.
- 4) For a coded program of this difficulty a faster machine with a much larger fast memory is very desirable.

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INTRODUCTION

The desirability of a uniform power distribution ($P_{\max} / P_{\text{av}} = 1$) has long been realized since it will not only decrease the mass of the fuel required for the attainment of criticality but also will increase the total power output by allowing the whole core instead of only a small fraction of the reactor to operate at the highest allowable temperature. Operation at the highest possible temperature allowed by metallurgical considerations over the entire reactor will increase the average core life since more uniform burn up is attained.

It is possible to achieve this uniform power distribution by various means. One of the first methods to be used was that of poisons. In this method long lasting poisons are placed in the reactor structure in the anticipated locations of high flux. By placing poison in these regions the peak flux is reduced to such a level that the power density is approximately uniform. This procedure is a very wasteful one from the point of view of neutron economy. The regions of the reactor that are contributing greatly to the overall neutron efficiency of the reactor are penalized to the extent that they can be no more efficient than the most inefficient region. For power producing reactors, particularly mobile ones, this waste of neutrons is very undesirable.

The use of poison is the first of two methods that exploit a nonuniform macroscopic cross section to achieve the uniform distribution of power. The second consists in the use of a nonuniform macroscopic fission cross section. This method allows each region of

the reactor to operate most efficiently. The efficiency is determined by the position of the region within the reactor. The leakage from a reactor causes a reduction in the value of the flux at large distances from the center of the core. To compensate for this reduction in flux the macroscopic fission cross section is increased in the regions far from the center and decreased in the central portions. To this end, the fuel concentration is increased near the edges.

Neutron economy can also be improved by surrounding the core with neutron reflector material. A light substance having a relatively low absorption cross section performs this function very well, i.e., heavy water or beryllium.

Since the reactor to be investigated is intended for a hypothetical naval cruiser, the use of poisons to achieve a uniform power distribution is not desirable, for a compact reactor is required. Though the reactor must be compact, it is desirable to install a reasonably thick reflector for the purpose of improving the neutron economy, attenuating the neutrons leaking out, and cooling of the reactor shell and pressure vessel.

It would be possible to have a nonuniform macroscopic cross section of the reactor in each of its directions. For the spherical reactor the variation would be radial only, but the application of a spherical reactor to shipboard use is impractical. The normal configuration of spaces aboard a ship is rectangular and a spherical reactor does not adapt very well to this arrangement. However, a rectangular reactor allows efficient utilization of space aboard a ship, but from the nuclear point of view this shape is not too desir-

able because leakage depends on the relationship of surface area to volume. It can be seen that the neutron economy decreases in progression from the sphere to the cylinder to the rectangular parallelepiped.

With these considerations in mind a right cylindrical reactor shape will be investigated. With this shape it would be possible to vary the properties in two directions, radially and axially. While this goal is theoretically interesting, it is prohibited when cost is considered. The reactor ~~that is to be~~ investigated will have a uniform composition in the axial direction with finite radial segments differing in properties.

For the shipboard use of a nuclear reactor power plant it is desirable to have a plant as light and safe as can possibly be designed within the space allowed but still without incorporating undue quantities of the "exotic" and expensive materials or coolants. One possible solution is the use of a highly enriched reactor using a coolant that permits high operating temperatures at low pressures and does not become dangerously radioactive while passing thru the reactor core.

A coolant that has a low vapor pressure at high operating temperature will allow a system pressure which is much lower than that employed in the present day water cooled and moderated shipboard reactor installation. The use of such a coolant will not only result in much better steam conditions, but it will also make possible an overall weight reduction, since a low pressure system can be used. The saving in weight is particularly significant in the pressure

vessel.

If the coolant does not become radioactive in its passage thru the reactor core, the shielding can be concentrated primarily around the reactor proper, provided that impurities in the coolant are kept at a very low concentration. Even with a small amount of impurities, the radiation emitted by the coolant would be low enough to permit reduced thicknesses of shielding around the primary loop.

A liquid organic hydrocarbon will fulfill the requirements listed above, but also will present some unsolved problems. It will (1) tend to increase the size of the reactor by decreasing the hydrogen moderation relative to a water cooled and moderated reactor, (2) tend to increase the reactor size due to its less efficient heat transfer capabilities compared with water and (3) exhibit both thermal and radiolytic breakdown in which the original hydrocarbon will, primarily by polymerization, change into longer chain hydrocarbons with the evolution of relatively large amounts of gas and a great increase in viscosity of the coolant.

The latter problems will be investigated at the National Reactor Testing Station on the Organic-Moderated Reactor Experiment (ORME). This experiment is not a prototype for an organic power reactor but is being built to test the feasibility of its use as a reactor coolant. [1] At the West Milton Annex of the Knolls Atomic Power Laboratory the Atomic Energy Commission plans to investigate the feasibility of an organic moderated and cooled reactor for naval propulsion. This will be conducted on the Naval Organic Reactor Experiment, (NORE).

The heat transfer properties of the various liquid organic hydrocarbons that are presently under consideration are not generally available. Some presently under investigation are diphenyl, O-terphenyl, M-terphenyl and P-terphenyl. A proprietary heat transfer medium "Dowtherm" is available commercially and is composed of 73.5% diphenyl oxide and 26.5% diphenyl. It has a working range of 450-750°F. at pressures less than 145 psig. Since "Dowtherm" closely resembles the hydrocarbons that will be investigated on the ORME and its heat transfer properties are readily available^[2] it will be used in this investigation.

The fuel element chosen for this investigation is of the MTR (Materials Testing Reactor) type.^[3] An effective length of element was set at 48". This length was chosen so that the overall height of the reactor, control devices and shield would fit into the present day engineering spaces of a cruiser. Any additional height would have an adverse effect on stability, since it would result in a higher shield deck and thus an increase in the height of the center of gravity of the ship.

Fuel Element Dimensions

$$t_m = 0.020''$$

$$t_c = 0.015''$$

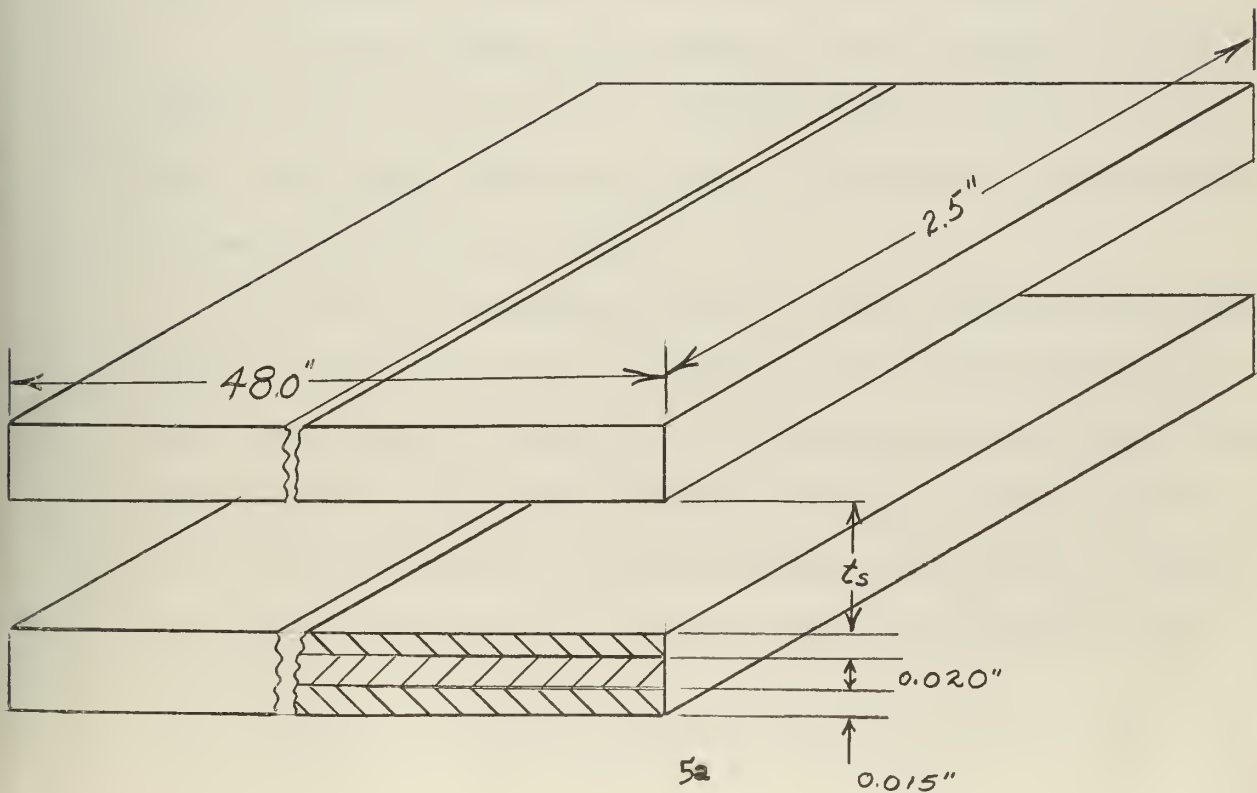
$$W = 2.5''$$

$$L = 48.0''$$

$$t_s^1 = 0.115'' \quad (\text{Case 1})$$

$$t_s^2 = 0.0919'' = 0.8 t_s^1 \quad (\text{Case 2})$$

$$t_s^3 = 0.069'' = 0.6 t_s^1 \quad (\text{Case 3})$$



PROCEDURE

Heat Transfer¹

The controlling variables in this study were (1) the maximum surface temperature of the fuel element or the maximum temperature at the centerline of the fuel element and (2) the maximum coolant speed. The maximum surface temperature of the fuel element may be controlling, since the preliminary investigations on the ORME have shown that above 800° F the thermal breakdown is increased greatly. [4] However, liquid loading requirements of a naval vessel make storage of used coolant a simple matter and in the interest of higher efficiency some breakdown will be accepted. Therefore, a fuel element surface temperature less than 850° F will be accepted. If the maximum fuel temperature at the centerline is allowed to exceed a value of about 1130° F, a phase transformation of uranium is encountered in which an anisotropic expansion and possible rupture of the fuel element may result. Therefore, a second design criterion will be required that the fuel element centerline temperature be less than 1130° F.

Preliminary studies on the ORME have also shown that coolant speeds greater than 15 feet per second result in excessive erosion. As a third design criterion, it will be required that coolant speed be less than 15 feet per second.

The first investigation was with a reactor coolant outlet temperature of 750° F with various temperature rises across the reactor. This investigation was amplified by varying coolant speed while holding the temperature rise across the reactor constant. Each of these situations were further studied by considering the MTR fuel element modified to the extent that the plate spacing was narrowed to 80% and 60%.

¹ Nomenclature for Heat Transfer on pages 22 and 23.

Though it was known in advance that the speed would be kept at or less than 15 feet per second calculations were pursued with higher values. This was done to show that a reduction in size can be accomplished in the future, if the erosion problem at high coolant speeds can be overcome.

A heat rate requirement of 6.06×10^8 Btu/hr was chosen based on the full power steaming rate of a typical cruiser.

With this value of Q (6.06×10^8) a coolant outlet temperature, coolant temperature rise, coolant speed and plate spacing are chosen. Since the inlet and outlet temperatures are known the average values of density and specific heat may be calculated. Formulae 1, 2, 3, and 4 are used respectively to calculate w , A_f , G and n .

$$w = \frac{Q}{\frac{c_p}{p} \Delta T} = G A_f \quad (1)$$

$$A_f = \frac{w}{3600 \bar{\rho} v} \quad (2)$$

$$G = v \bar{\rho} = \frac{w}{3600 A_f} \quad (3)$$

$$n = \frac{A_f}{A_n} \quad (4)$$

Since the dimensions of the fuel element for each case are known, both the hydraulic diameter and the total flow area can be calculated. With the product of the mass flow rate and the hydraulic diameter the product of the heat transfer coefficient and the hydraulic diameter can be read from the graph in the "Dowtherm" booklet. Formulae 5, 6, 7,

and 8 are then used to calculate, respectively, Q_v , ΔT_{fuel} , ΔT_{clad} , and θ .

$$Q_v = \frac{Q}{v_f} \quad (5)$$

$$\Delta T_{\text{fuel}} = T_o - T_1 = \frac{t^2 Q_v}{4k_f} \quad (6)$$

$$\Delta T_{\text{clad}} = T_Q - T_2 = \frac{t Q t}{2k_c} \quad (7)$$

$$\theta = T_2 - T_b = \frac{Q t}{v m} \quad (8)$$

Datum from the calculation of a coolant outlet temperature of 750° F for various values of temperature rise and speed for cases 1, 2 and 3 are in Tables I, II, and III and Figures I, II, and III respectively. The cladding temperature existing for each of these conditions is plotted on Figure IV.

The narrowest plate spacing, case 3, gives a significantly smaller reactor. Since the pumping power of any of these cases proved to be very small compared with the total power output, it was decided to investigate the result of a decrease of coolant outlet temperature for this case only. Coolant outlet temperature was reduced to 700° V and the speed was held at 15 feet per second. Datum from this calculation are in Table IV. These data were plotted as a dotted line on Figure IV.

For these twelve conditions the fraction of uranium in the uranium-zirconium dispersed fuel plate was calculated and plotted on Figure V. It is desirable that the fuel be of such a composition that it contributes

TABLE I

Datum for coolant outlet of 750°F
and wide plate spacing
(case 1)

A. REACTOR $\Delta T = 200^\circ\text{F}$

v (ft/sec)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	A_f (ft ²)	n (no. of plates)
5	236.1	686.	5.36	2,685
15	709.0	1611.	1.79	896
25	1180.0	2340.	1.07	538
50	2360.0	4090.	0.54	269

v (ft/sec)	$Q_v \times 10^8$ (Btu/hr-ft ³)	$T_o - T_1$ (°F)	$T_1 - T_2$ (°F)	$T_2 - T_b$ (°F)	T_2 (°F)	f_u %
15	4.86	20.7	62.2	215.8	1015.8*	25.4

B. REACTOR $\Delta T = 150^\circ\text{F}$

v (ft/sec)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	A_f (ft ²)	n (no. of plates)
5	231.9	682.	7.26	3,639
15	705.0	1591.	2.42	1,211
25	1160.0	2320.	1.45	727
50	2318.0	4040.	0.73	364

v (ft/sec)	$Q_v \times 10^8$ (Btu/hr-ft ³)	$T_o - T_1$ (°F)	$T_1 - T_2$ (°F)	$T_2 - T_b$ (°F)	T_2 (°F)	f_u %
15	3.60	15.4	46.0	197.3	947.3*	19.5

* excessive temperature

Y uranium content too low

Table I (Continued)

C. REACTOR $\Delta T = 100^\circ\text{F}$

v (ft/sec)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	A _f (ft ²)	n (no. of plates)
5	227.	669.	11.01	5,520
15	681.	1569.	3.67	1,839
25	1135.	2295.	2.02	1,103
50	2270	4000.	1.10	552

v (ft/sec)	Q _v x 10 ⁸ (Btu/hr-ft ³)	T ₀ -T ₁ (°F)	T ₁ -T ₂ (°F)	T ₂ -T _b (°F)	T ₂ (°F)	f _u %
15	2.37	10.1	30.3	133.8	883.8*	13.5

D. REACTOR $\Delta T = 50^\circ\text{F}$

v (ft/sec)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	A _f (ft ²)	n (no. of plates)
5	221.8	659.	22.40	11,230
15	667.0	1545.	7.45	3,730
25	1110.0	2275.	4.47	2,240
50	2218.0	3910.	2.24	1,123

v (ft/sec)	Q _v x 10 ⁸ (Btu/hr-ft ³)	T ₀ -T ₁ (°F) ¹	T ₁ -T ₂ (°F)	T ₂ -T _b (°F)	T ₂ (°F)	f _u %
15	1.17	4.97	15.0	66.2	816.2	6.99

* excessive temperature

γ uranium content too low

TABLE II

Datum for coolant outlet of 750°F,
coolant speed of 15 ft/sec, and medium plate spacing
(case 2)

ΔT (°F)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	n (no. of plates)	$Q_v \times 10^8$ (Btu/hr-ft ³)
200.	709.	1671.	1121.	3.89
150.	705.	1660.	1515.	2.88
100.	681.	1639.	2300.	1.89
50.	667.	1631.	4660.	0.94

ΔT (°F)	$T_c - T_1$ (°F)	$T_1 - T_2$ (°F)	$T_2 - T_b$ (°F)	T_2 (°F)	f_u %
200	16.6	49.7	193.9	943.9*	19.6
150	12.3	36.9	144.2	894.0*	16.1
100	8.1	24.3	96.3	846.3	11.0
50	4.0	12.0	47.6	797.6	5.6 ^γ

* temperature excessive

γ uranium content too low

TABLE III

Datum for coolant outlet of 750°F,
coolant speed of 15 ft/sec, and narrow plate spacing
(case 3)

ΔT (°F)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	n (no. of plates)	$Q_v \times 10^8$ (Btu/hr-ft ³)
200	709	1796	1496	2.9
150	705	1775	2020	2.2
100	681	1731	3061	1.4
50	667	1718	6230	0.70

ΔT (°F)	$T_o - T_1$ (°F)	$T_1 - T_2$ (°F)	$T_2 - T_b$ (°F)	T_2 (°F)	f_u %
200	12.4	37.3	135.2	885.2*	16.3
150	9.2	27.6	101.3	851.3	12.4
100	6.1	18.2	68.5	818.5	8.4 ^γ
50	3.0	9.0	33.9	783.9	4.3 ^γ

* temperature excessive

γ uranium content too low

FIGURE 1

Number of plates vs coolant speed
 Coolant contact 750°F
 Wide plate casing
 (Case 1)

Revolutions per minute

- 50°F
- × 100°F
- △ 150°F
- 200°F

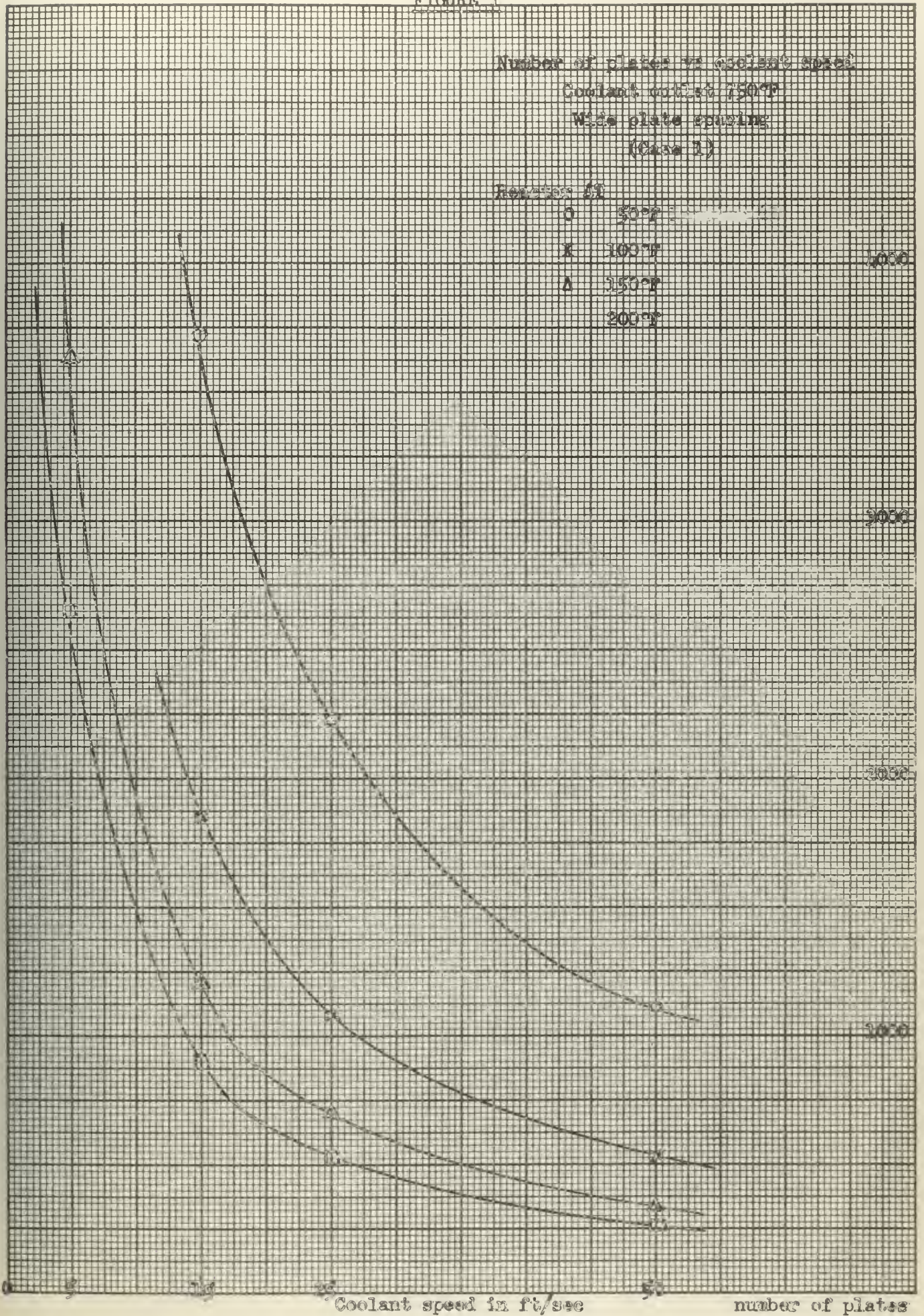
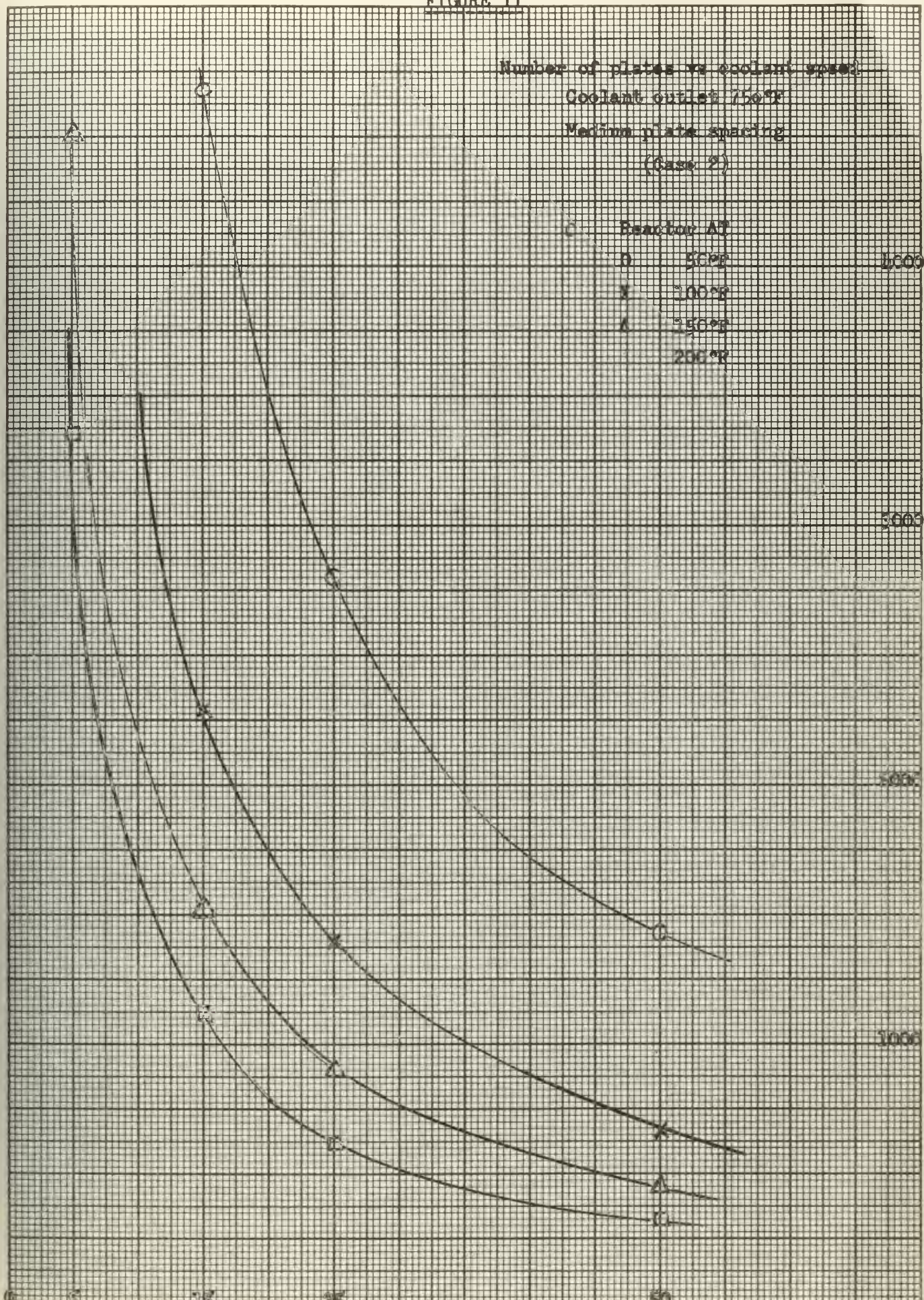


FIGURE II

Number of plates vs coolant speed
 Coolant outlet 75°F
 Medium plate spacing
 (Case 2)

O Reactor AT
 D 50°F
 X 100°F
 Δ 150°F
 □ 200°F



Coolant speed in ft/sec

number of plates

FIGURE III

Number of plates vs coolant speed
 Coolant inlet at 750°F
 Narrow plate spacing
 (Case 3)

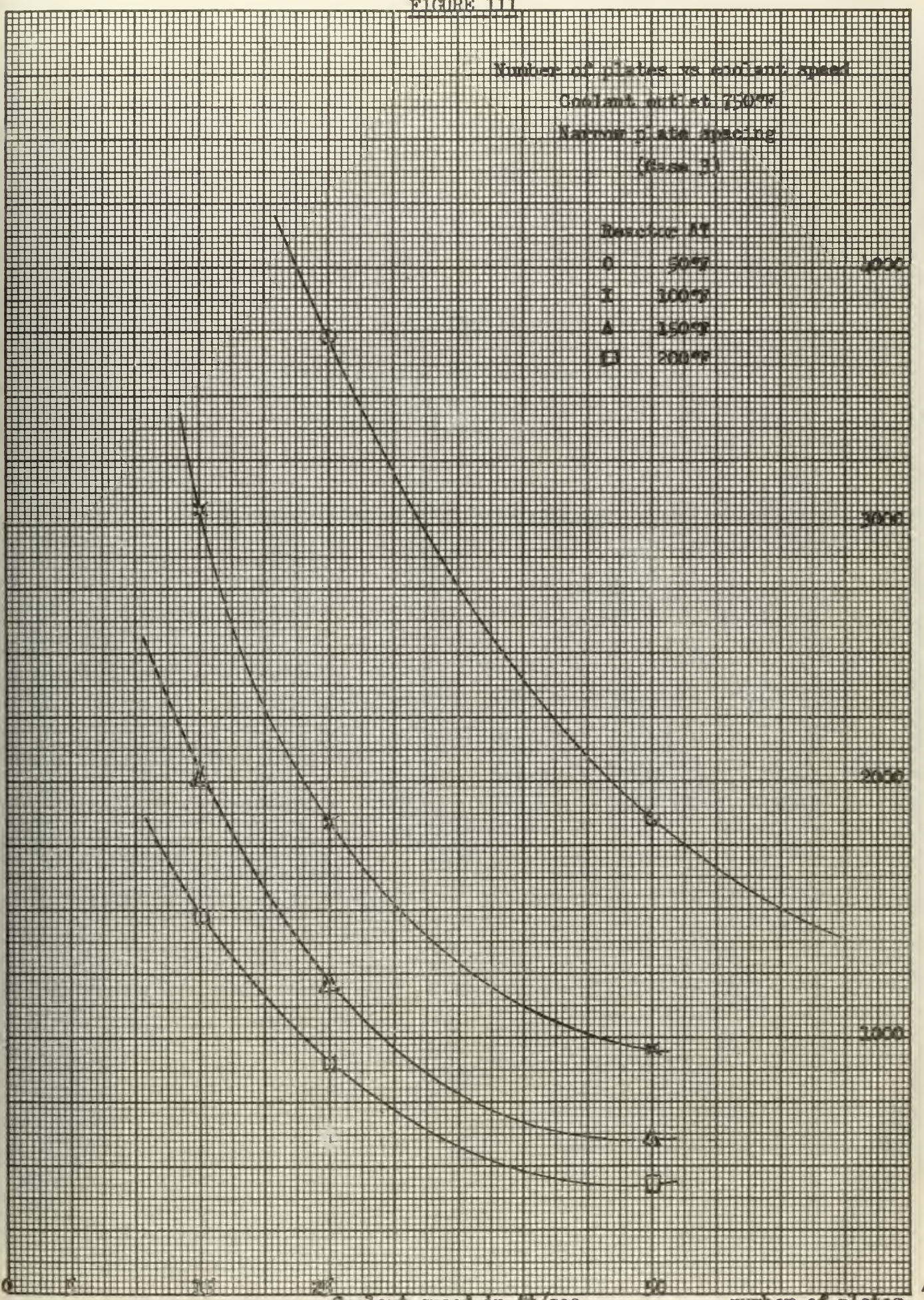
Reactor AT

0 50°F

I 100°F

A 150°F

EJ 200°F



Coolant speed in ft/sec

number of plates ↑

TABLE IV.

Datum for coolant outlet of 700°F,
coolant speed of 15 ft/sec, and narrow plate spacing
(case 3)

ΔT (°F)	G (lb/sec-ft ²)	h (Btu/hr-ft ² -°F)	n (no. of plates)	$Q_v \times 10^8$ (Btu/hr-ft ³)
200	739.5	1826	871	5.0
150	726.8	1812	1164	3.7
100	712.5	1775	1767	2.5

ΔT (°F)	$T_o - T_1$ (°F)	$T_1 - T_2$ (°F)	$T_2 - T_b$ (°F)	T_2 (°F)	f_u %
200	21.3	64.0	228.4	924.4*	26.0
150	15.9	47.9	172.2	872.2	20.3
100	10.5	31.5	115.8	815.8	14.0

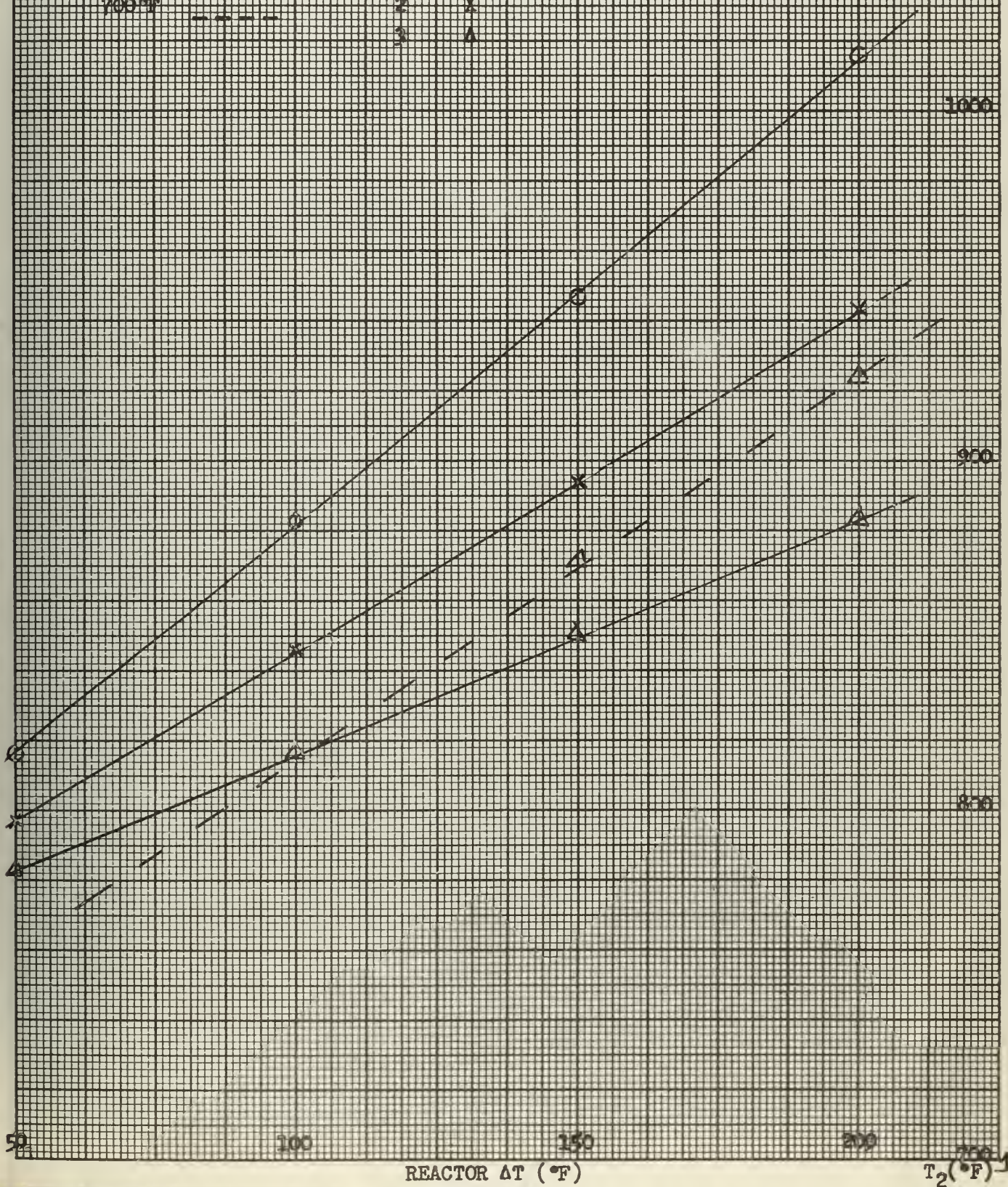
* temperature excessive

γ uranium content too low

FIGURE IV

Fuel element surface temperature
vs reactor ΔT
Coolant speed 15 ft/sec

Coolant outlet T_2	Case
750°F ———	1 O
700°F ———	2 X
650°F - - - -	3 A



REACTOR ΔT (°F)

T₂ (°F)

FIGURE V

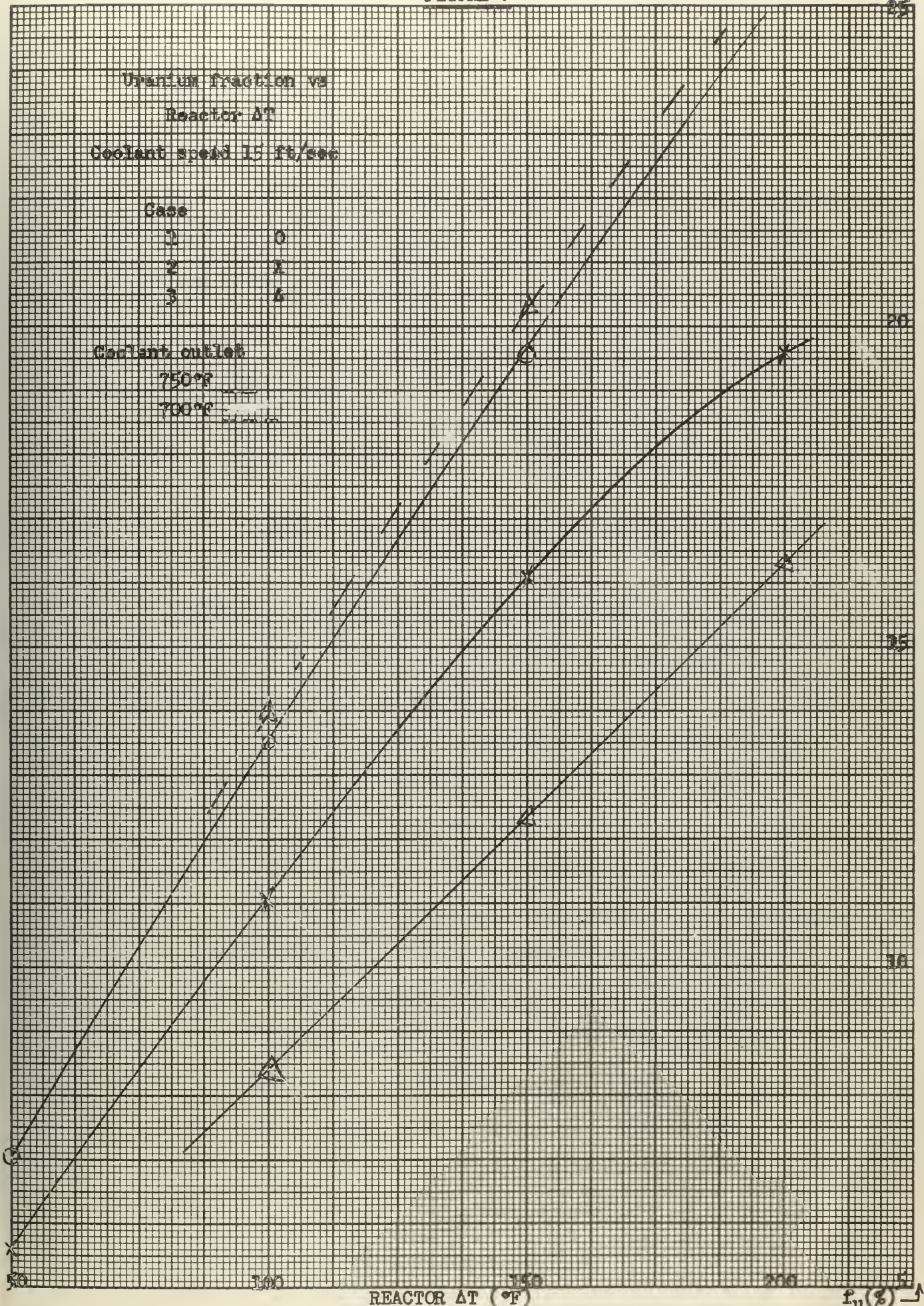
Uranium fraction vs
 Reactor ΔT
 Coolant speed 15 ft/sec

Case	
1	0
2	X
3	A

Coolant outlet

750°F

700°F



REACTOR ΔT (°F)

f_u (%)

to the structural strength of the reactor. A uranium content of 10% by weight or greater has desirable strength characteristics.

Of the six conditions that have a fuel element surface temperature of less than 850° F only two have a uranium fraction of above 10%. These are shown in Table V.

t_{co}	T	case	n	T_2
750	100	3	3061	818
700	100	3	1767	815

Table V Particularly Interesting Cases

Figure VI shows the variation in number of plates for case 3 holding the coolant speed at 15 feet per second and varying the temperature rise across the reactor.

The size resulting from the condition with a coolant outlet temperature of 700° F is better for the following reasons:

- 1) it is by far the smaller,
- 2) the increase of uranium fraction by 2 1/2% gives a significant increase in the strength of the fuel region.

With this condition the fuel element surface temperature will be less than the prescribed value of 850° F. The amount of damage that will result to the coolant from this temperature and the accompanying radiation is uncertain. However, for shipboard applications it does not appear to be a serious penalty.

In the present day cruiser design for torpedo protection there is available a large amount of bunker fuel oil to partially fill the

FIGURE VI

Number of plates vs
 Reactor ΔT ($^{\circ}F$)
 Coolant speed 15 ft/sec
 Narrow plate spacing
 (case 3)

Coolant outlet temperature

750 $^{\circ}F$	X
700 $^{\circ}F$	O

3000

2000

1000

50

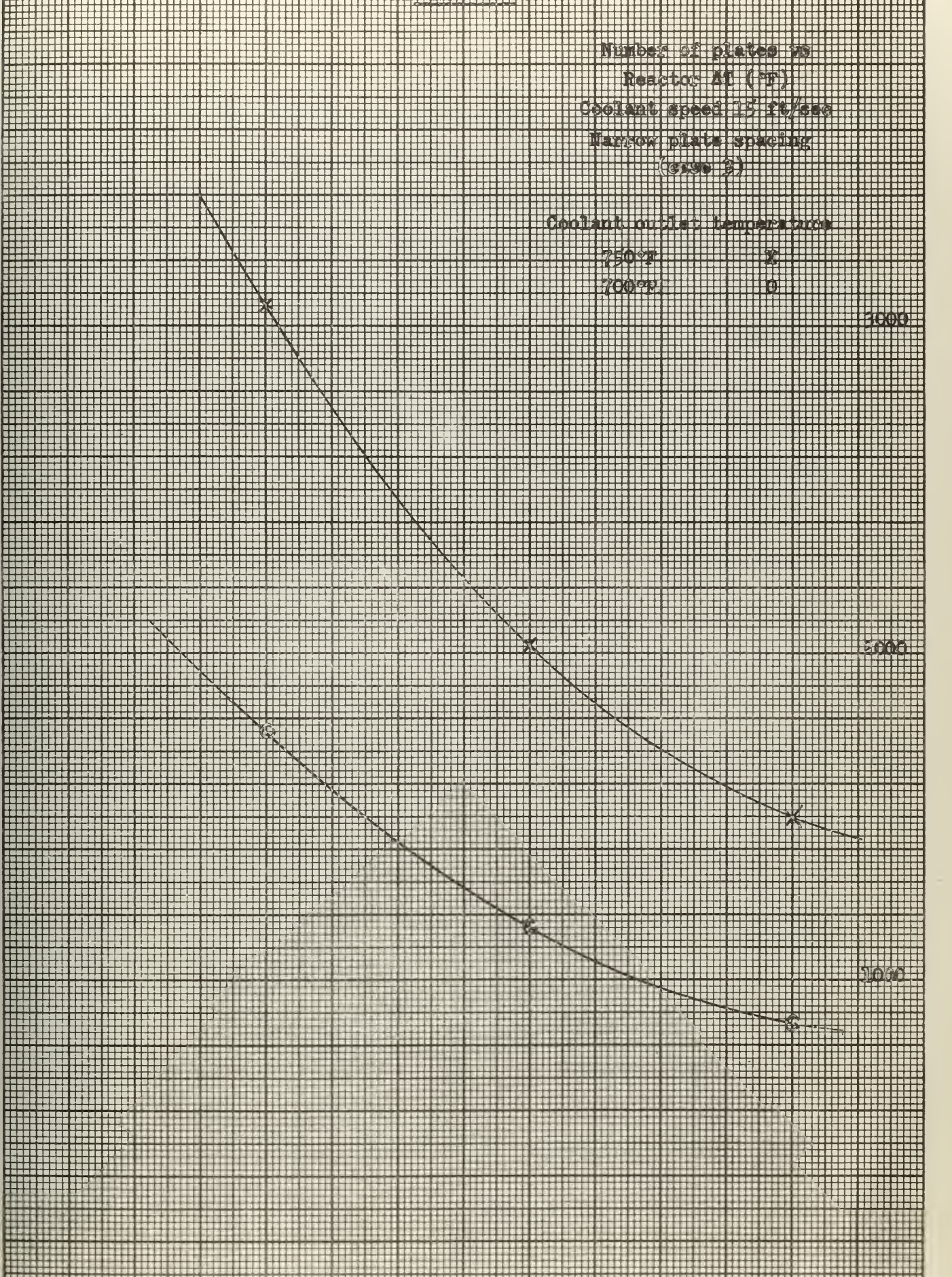
100

150

200

REACTOR ΔT ($^{\circ}F$)

number of plates \uparrow



layers of the torpedo protective system. Certain layers of the torpedo protective system are required to be liquid loaded. This system could be used for the storage of the organic moderator coolant. The unirradiated coolant could be stored in the inner layers of the system and that which had been irradiated would be stored in the outer layers to protect the crew from any induced radioactivity. The long chain hydrocarbons would tend to settle to the lower parts of the tanks where they could be removed during periods of overhaul. With the installation of a purification system this process of separation might be handled continuously aboard ship in the same manner that lubrication oil is purified and recycled.

Nomenclature

A_f	flow total area, (ft ²)
A_f^n	flow area per plate case n, (ft ² /plate)
A_h	heat transfer area, (ft ² /plate)
\bar{c}_p	average specific heat, (Btu/lb °F)
D'	hydraulic diameter, (inches)
f_u	fraction of uranium in fuel region
G	mass flux, (lb/sec-ft ²)
h	film coefficient, (Btu/hr-ft ² -°F)
k_c	thermal conductivity of cladding, (Btu/hr-ft-°F)
k_f	thermal conductivity of cladding, (Btu/hr-ft-°F)
L	fuel length, (inches)
n	number of fuel plates
Q	total heat generation rate, (Btu/ft ³)
Q_v	heat generation rate per unit volume of fuel, (Btu/hr-ft ³)
t_c	fuel cladding thickness, (inches)
t_{co}	coolant outlet temperature, (°F)
t_m	fuel thickness, (inches)
t_s	spacing between plates, (inches)
T_b	bulk temperature of coolant, (°F)

Nomenclature (Continued)

T_0	fuel centerline temperature, ($^{\circ}\text{F}$)
T_1	fuel-clad interface temperature, ($^{\circ}\text{F}$)
T_2	fuel element surface temperature, ($^{\circ}\text{F}$)
U	overall heat transfer coefficient, ($\text{Btu/hr-ft}^2\text{-}^{\circ}\text{F}$)
v	coolant speed, (ft/sec)
w	mass rate of flow, (lb/hr)
W	fuel width, (inches)
ΔT	overall temperature difference across reactor, ($^{\circ}\text{F}$)
$\bar{\rho}$	average density, (lb/ft^3)
θ	film temperature drop, ($^{\circ}\text{F}$)

Basic Reactor Physics

Criticality calculations were made using a one group reflected homogenous model with case 3, narrow fuel elements. Based on heat transfer considerations, inlet and outlet coolant temperature were set at 600° and 700° F, respectively, with a coolant speed of 15 ft/sec.

The assumptions in this calculation were

1. The structural members inside the reactor were of type 347 stainless steel and composed ten percent of the total reactor weight,
2. A fast fission factor of unity,
3. The uranium of fuel is enriched to 90% in U^{235} ,
4. A right circular, cylindrical flux distribution of a cosine axially and a Bessel function radially,
5. A beryllium reflector of 236 cm placed around the reactor.

The reactor was treated as a homogeneous medium. By a one group calculation the volumes of each constituent of the reactor were calculated from the dimensions given previously in the heat transfer section. These volumes were then converted to weights and the volume fraction of the total reactor as shown in Table I. From the cross sections given in BNL-325^[5] effective cross sections were calculated for the operating temperature of the material. With these cross sections the reactor parameters as shown in Table II were calculated. These give an effective multiplication factor of 0.894. This value is sufficiently close to unity that final adjustments are made by the computing machine.

The primary purpose of the calculation is to obtain an approximate

value of the critical radius which is needed as an input to the computer solution.

Table I

Reactor Materials

	Volume (cm ³)	Mass (Kgm)	Volume fraction
Fuel	6.95 x 10 ⁴	495.	
U		69.5	0.0068
Zr		425.5	0.408
Clad (Zr)	15.01 x 10 ⁴	966	
Stainless Steel	1.99 x 10 ⁴	160	
Cr (18% by weight)		28.4	0.007
Ni (11% by weight)		17.35	0.004
Fe (71% by weight)		112.	0.027
Coolant	26.3 x 10 ⁴	219.5	
[73.5 Diphenyl oxide by weight (C ₆ H ₅ OC ₆ H ₅) and 26.5% Diphenyl by weight (C ₆ H ₅) ₂]			
C (06.91 by weight)			0.47043
H (06.03 by weight)			0.03292
O (86.59 by weight)			0.3774

Table II

Reactor Parameters

δ	= Beryllium shield thickness	=	23.6	cm.
Σ_{tr}	= $\sum_i \Sigma_i (1-\mu_0)_i v_i \left(1 - \frac{1}{5} \frac{\Sigma_{a_i}}{\Sigma_{t_i}} \right)$	=	0.28334	cm ⁻¹
D	= $3 \frac{1}{\Sigma_{tr}}$	=	1.17644	cm
d	= $0.71 \lambda_{tr} \frac{1+\beta}{1-\beta} = 33.7D$	=	39.599	cm
Σ_a	= $\sum_i \Sigma_{a_i} v_i$	=	0.1208	cm ⁻¹
L^2	= $\frac{D}{\Sigma_a}$	=	9.7387	cm ²
$p(E)$	= $\exp \left[\frac{\bar{N}_{28}}{\Sigma_s} (3.9) \left(\frac{\Sigma_s \times 10^{24}}{N_{28}} \right)^{0.415} \right] \times 10^{-24}$	=	0.726	
η_f	= $v \frac{\Sigma_f}{\Sigma_{a(\text{total})}}$	=	1.9692	
k_{∞}	= $\eta_f p \epsilon$	=	1.4296	
	= $\frac{\sum_i \Sigma_{s_i} \bar{\nu}_i}{\Sigma_s}$	=	0.17598	
γ	= $\frac{D}{\Sigma_s} \ln \left(\frac{E_1}{E_2} \right)$	=	343	cm ²
B^2	= $\left(\frac{2.405}{R+d} \right)^2 + \left(\frac{\pi}{H+26z} \right)^2$	=	1.332 x 10 ⁻³	cm ²
β	= albedo of Beryllium			
k_{eff}	= $\frac{k_{\infty} e^{-B^2 \gamma}}{1+B^2 L^2}$	=	0.894	

Application of the Flux Equations

The S_n approximation is a method of differencing in which the independent variable is quantized into n intervals, and intermediate values of a function of the variable are approximated by a straight line between points at the boundaries of the interval. This method of numerical integration was developed by Bengt G. Carlson [6].

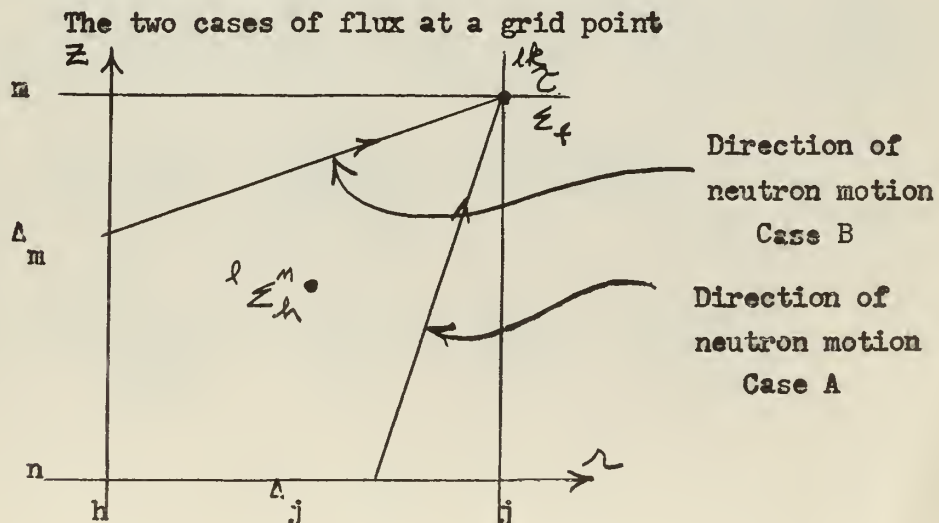
The vector flux at a point within a reactor can be obtained by the use of the S_2 method. The equations necessary are derived in Appendix A for two cases of neutron velocity:

Case A - The component of the vector flux in the axial direction is greater than its component in the radial direction.

Case B - The component of the vector flux in the radial direction is greater than that in the axial direction.

These two cases are illustrated in Figure I using a grid which is used to difference the transport equation in the r and z coordinates, r being the radial coordinate and z the axial coordinate.

Figure I



The general flux equation for case A is:

$$l_{\phi_{j,m}^{i,f}} = \frac{|b_i| (1-\omega)^{i,f} \left(l_{\phi_{j,n}^{i,f}} + l_{\phi_{j,n}^{i-1,f}} - l_{\phi_{j,m}^{i-1,f}} + \omega A_2^{i,f} \right)}{|b_i| + \omega \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_h^n} + \frac{c_i}{r_h} \right)} \quad (1)$$

where h is a value which is one greater or less than j, depending upon the direction of the radial component of vector flux, and n is a value which is one greater or one less than m, depending upon the direction of the axial component of vector flux. Definitions of the other quantities can be found in the nomenclature part of this section.

For Case B the flux at a point is defined by the following equation:

$$l_{\phi_{j,m}^{i,f}} = \frac{|b_i| (1 - \omega^{i,f}) \left(l_{\phi_{h,m}^{i,f}} + l_{\phi_{h,m}^{i-1,f}} - l_{\phi_{h,m}^{i,f}} - l_{\phi_{h,n}^{i-1,f}} \right)}{|b_i| + \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_h^n} + \frac{c_i}{r_h} \right)} \quad (2)$$

where A for Equation (1) and (2) is defined by:

$$A = \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_h^n} + \frac{c_i}{r_h} \right) \right] l_{\phi_{h,n}^{i,f}} - \left[|b_i| + \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_h^n} - \frac{c_i}{r_h} \right) \right] l_{\phi_{j,m}^{i-1,f}} \\ + \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_h^n} - \frac{c_i}{r_h} \right) \right] l_{\phi_{h,n}^{i-1,f}} + \frac{g_i g_f |\Delta_j|}{2} \left[l_{S_{j,m}} + l_{S_{h,n}} \right] \\ + \left[\frac{|b_i| - |d_i|}{2} \right] \left[l_{\phi_{j,m}^{i-1,f}} + l_{\phi_{j,n}^{i-1,f}} - l_{\phi_{h,m}^{i-1,f}} - l_{\phi_{h,n}^{i-1,f}} \right] \quad (3)$$

It should be pointed out that the total macroscopic cross section, Σ , is evaluated using the material properties at the center of the position grid of Figure I, but are designated by the grid values at the corner diagonally opposite the point j,m , i.e.; the point h,n . Further, \bar{r}_h is actually the mean radius and is equal to $\frac{r_j + r_h}{2}$.

The source at a point is defined by the following equations

$$S_{j,m} = \nu_p \sum_{k=1}^3 \frac{\Sigma_k}{\Sigma_f} \phi_{j,m}^k + \sum_{k=1}^3 \lambda_k \tau^k \phi_{j,m}^k$$

where the scalar flux $\phi_{j,m}$ is equal to

$$\sum_{f=0}^1 H_f \sum_{i=0}^2 P_i \phi_{j,m}^{i,f}$$

The fission cross section and transfer probabilities, τ , are evaluated using the material properties at the grid point at which the flux is being calculated. The data necessary for the solution of the flux at a point is summarized in Tables I. and II.

Table I.

Data for the S_2 approximation						
\underline{i}	\underline{i}	\underline{b}_i	\underline{c}_i	\underline{d}_i	\underline{e}_i	\underline{P}_i
0	-1	-1	0	--	1	+0.3183
1	0	-1/3	4/3	-2/3	2	+0.3643
2	1	2/3	4/3	1/3	2	+0.3183

Table II.

Data for Gauss quadrature approximation

<u>f</u>	<u>μ_f</u>	<u>h_f</u>	<u>g_f</u>	<u>H_f</u>
0	-0.8611363	-1.6944625	+1.9677054	+0.3478548
1	-0.3399801	-0.3197284	+1.063340	+0.6521452
2	+0.3399801	+0.3197284	+1.063340	+0.6521452
3	+0.8611363	+1.6944625	+1.9677054	+0.3478548

Not all pairs of values for i and f are relevant for each of the cases A and B as a consequence of the defining inequalities for these cases. The pairs of relevant values are discussed in connection with Figure II. A general point may be defined to be any grid point of the reactor other than a point at the center, on the centerline or in the midplane of the reactor. A point at these other places is called an origin point, a center-line point or a midplane point, respectively. For Case A for which $|b_i| < |h_f|$ and in the case of the first quadrant flux, there is a further requirement (Table I, Appendix A) that b_i and H_f have negative signs. In this quadrant there are three Case A flux equations, one for each of the pairs $(i=0, f=0)$, $(i=1, f=0)$, and $(i=1, f=1)$. Similarly, for Case B, the requirement $|b_i| > |h_f|$ leads to only one relevant combination.

The applicable combinations and those for the remaining three quadrants are summarized in Table III.

Figure II

Flux Configuration
for a particular energy group
at a general grid point

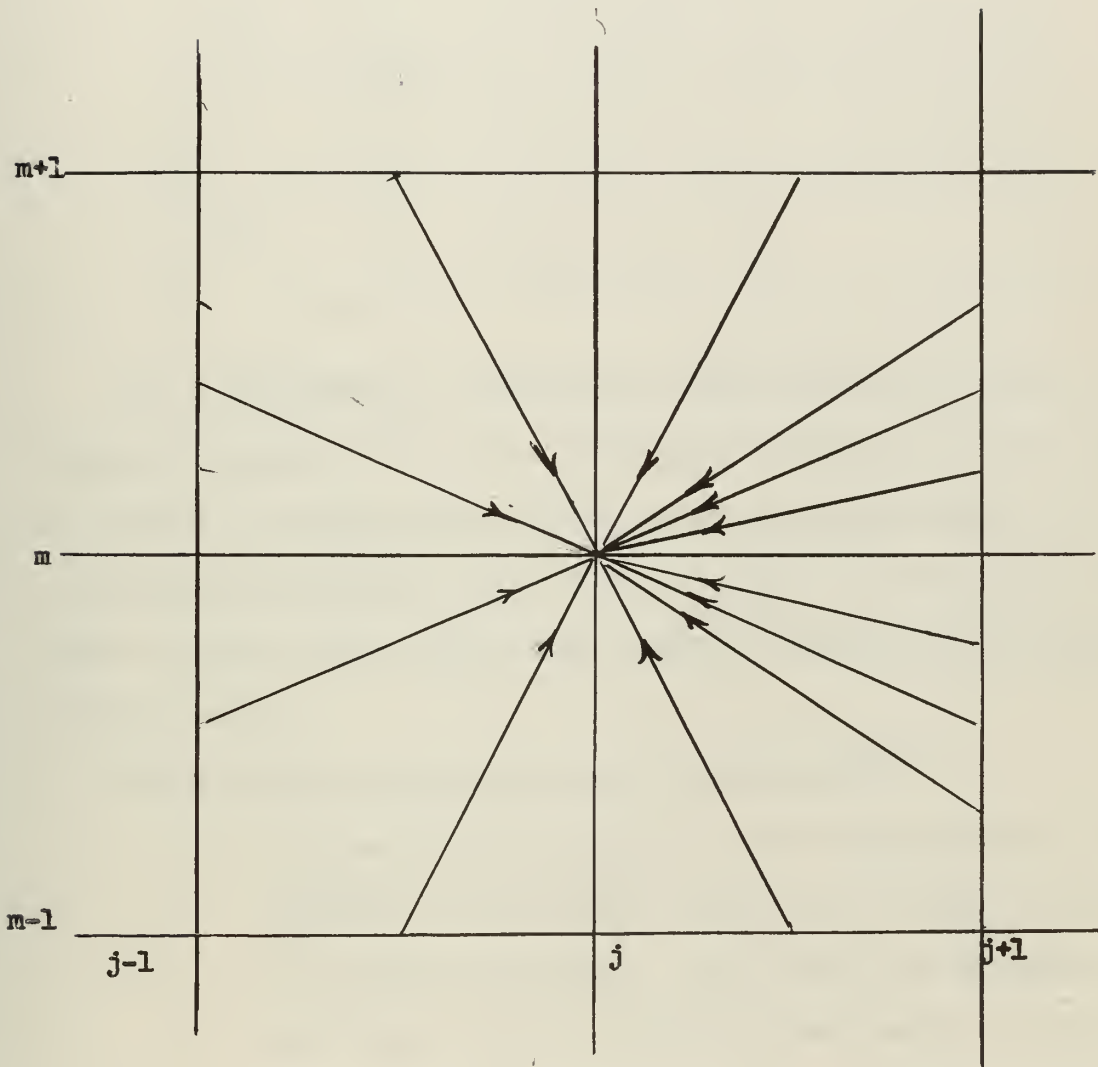


Table III

Allowable combinations of i and f

Quadrant	Case A	Case B
1 st	$i = 0, f = 0$ $i = 1, f = 0$ $i = 1, f = 1$	$i = 0, f = 1$
2 nd	$i = 2, f = 0$	$i = 2, f = 1$
3 rd	$i = 2, f = 3$	$i = 2, f = 2$
4 th	$i = 0, f = 3$ $i = 1, f = 3$ $i = 1, f = 2$	$i = 0, f = 2$

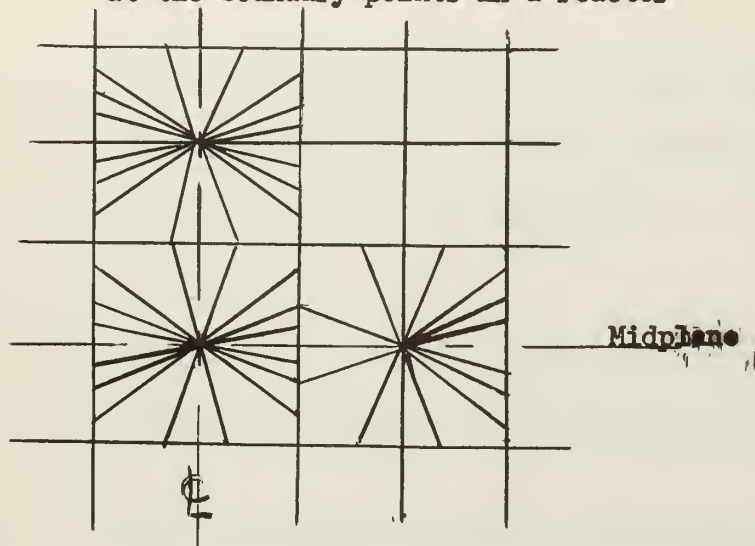
The configuration of the flux at a grid point which results from the allowable combinations of i and f are shown in Figure II. If the flux is assumed to be distributed in $\mathcal{L}+1$ energy groups there will be $12(\mathcal{L}+1)$ flux equation associated with every grid point. For the case considered, there are three energy groups and, therefore, thirty-six flux equations for each point.

The following boundary conditions are assumed: [7]

1. Continuity of flux exists at the centerline and at the midplane of the reactor, and further the flux is symmetrical about these boundaries. This permits the calculation of the flux in only one quarter of the reactor. Figure III shows the flux distribution at these various special points.

Figure III

S_2 approximation to the vector flux distribution
at the boundary points in a reactor



2. A second boundary condition is that at the outside

boundaries of the reactor there exists no inward flux.

In order to satisfy this requirements, an extra set of grid points is placed just outside these boundaries, and at these points the vector flux whose argument describes inwardly moving neutrons is assumed to be zero. Note, however, that leakage does exist at these fictitious points and at the outside boundary points of the reactor.

Nomenclature

b_i
 c_i
 d_i
 e_i

Values obtained by differencing and integration of the transport equation over \mathcal{N} , and defined by Equation (7a to 7d) of Appendix A

g_f
 h_f

Values depending upon the quantized values of μ and defined by Equation (7e and 7f) of Appendix A.

H_p

Values obtained by the use of the Gauss quadrature approximation to sum the flux over μ space, defined by Equation (13) Appendix A.

P_i

Values obtained by the use of the S_2 approximation to sum the flux over \mathcal{N} space, defined by Equation (12) Appendix A.

\bar{r}_h

mean radius between two adjacent grid points

S

Neutron source

l_p

Fraction of fission neutrons which are produced in a particular energy group l .

ν

Number of neutrons produced per fission.

$\lambda_{k\tau}$	The probability per unit length of neutron travel that a neutron is scattered in group k and that a neutron emerges in group ℓ from the scattering event.
Σ_f	Macroscopic fission cross section
Σ	Total Macroscopic cross section
$\phi(r, z, \mu, \nu, v)$	neutron flux which is classified according to speed, v , colatitude and azimuthal direction, cosines μ and ν , and position r and z .
Δ_j	spacing of grid in the radial direction.
Δ_m	spacing of grid in the axial direction.
μ	cosine of the angle Θ between the axis and the direction in which neutrons move. See Figure I Appendix A.
η	Cosine of the angle ψ . See Figure I Appendix A.
ω	See defining equation, page A-12 Appendix A.
ω'	See defining equation, page A-17 Appendix A.

Nomenclature

Superscripts and subscripts

- i index labeling the quantized values of η .
- f index labeling the quantized values of μ .
- j grid position index in the radial direction.
- m grid position index in the axial direction.
- ℓ neutron speed index.
- k neutron speed index of group in which neutron is scattered .
- n axial grid position diagonally opposite grid point at which flux is being calculated.
- h radial grid position.

A computer program is a sequence of logical instructions telling the computer what mathematical operations are to be performed on the data. Therefore, there are two groups of inputs, the data and the instructions. The data fed into the computer are stored in the fast, core memory if there is sufficient room. When the space is not adequate, some of the data are stored in magnetic drums. The only difference in the two storage processes is that the data in the drum takes longer to locate and bring into the core memory. Each number or piece of data is stored in a register or cell the location of which is designated by a number called the address of that register. The instructions are stored in sequence and are performed in the order in which they are stored unless the program directs otherwise. No unused registers are permitted to interrupt the storage as these will stop the computer.

The instructions given to the computer consist of two parts, the operation and the address sections. The operation section is first and uses letters to tell the computer the mathematical operation to be performed. The address section gives the address of the register containing the data on which the operation is to be performed. The instructions recognized by Whirlwind I are listed in the glossary.

For Whirlwind I there are many useful devices that make the programming much simpler. A few of these devices used in this program are floating addresses, preset parameters, and subroutines.

Probably the most useful device is the floating address. This floating address or "flad" is used to eliminate the necessity of referring to the absolute register address in the core memory when

writing the program. The absolute register address of any part of the program must be known primarily so that the initial or final instruction of data is not placed in a register containing part of the subroutines fixed in the memory of the computer.

In the program key instructions on which mathematical operations are to be performed or which will be referred to later in the program are given floating addresses. These addresses consist of a letter and a number, for example a1, a2, etc. Any letter except o and ell may be used and the sum of all the numbers following these letters may not exceed 255. When the program is read into the computer, it will assign absolute addresses or register numbers to these flads. A table of the floating addresses is produced by the computer with the solution unless it is specifically suppressed in the computer. This table gives the absolute address of every assigned floating address and is particularly helpful in trouble shooting a faulty program.

Another useful device is the preset parameter. These are numbers other than data that are used in the program. Preset parameters are frequently used to designate the number of times that a calculation is to be performed. Preset parameters must be assigned a value before they are used, and they will retain this value until it is specifically changed. Preset parameters are designated by two letters and a number. The first letter must be a p, u or z and the second letter and the number may be anything, except the letters o and el may not be used.

A third useful device is the subroutine. Its usefulness comes from the fact that it is a complete entity in itself and may be constructed and trouble shot independently. For the build up of a complete

program it is possible to start with one of the most basic mathematical operational sequences that will be performed repeatedly, program it, perform the hand calculations of the sequence and then place the routine on the computer. Modification of the subroutine or correction of the hand calculations must continue until the results of both are in agreement.

In arranging these subroutines in the main program extreme care must be used. The main program must be interrupted to call the subroutine into use and after the subroutine has performed its function it must direct the computer back to the proper place in the main program.

An example of a subroutine is as follows:

Main program

•		
•		
•		
	isu p 10	
	t1, isp x 4	take the next instruction from register x 4 and continue from there
	imr a 5	forms product of (-m2) and (a5)
•		
•		
•		
	i STOP	end of program
	x4, ita x 5	transfer into the address section of the instruction in register x5, the address that is one more than the address t1.
	ica m 2)	
	} its a 7)	places m 2 in a 7

ics m 2)
its a 8)

places -m 2 in a 8

x5, isp 0

returns to main program one register
after t l; i.e., register containing
imr a 5.

Since the numbers used in this program were over four decimal digits in length, floating point arithmetic was required. In this system two registers are required to express a number. Each of these registers is capable of containing 16 bits of binary number information. The first bit of the first register expresses the sign of the number and the first bit of the second register expresses the sign of the exponent of 2. Six bits are required to locate the binary point of the number leaving 24 bits to express the actual number. In certain stages of the program it is desirable to handle single length register numbers. The instructions for the handling of these numbers are similar to those for double length registers except that the i does not precede the instruction. When operating on double length registers the computer is said to be in the interpretive mode, when on single length registers it is in the Whirlwind mode. To enter the interpretive mode from the Whirlwind mode it is necessary to use the instruction "IN" and to reverse the direction it is necessary to use the instruction "OUT".

GLOSSARY OF INSTRUCTIONS

<u>TERM</u>	<u>DEFINITION</u>
ica x	clear MRA and add contents of register x.
ics x	clear MRA and subtract contents of register x.
iad x	add contents of register x to what is in MRA.
isu x	subtract contents of register x from what is in MRA.
imr x	multiply what is in MRA by the contents of register x and round off product to fifteen digits.
idv x	divide contents of MRA by the contents of register x.
its x	transfer the contents of the MRA to register x losing what was previously in register x. The MRA remains unchanged.
ies x	exchange the contents of the MRA with the contents of register x.
isp x	take the next instruction from register x. Does not affect the contents of the MRA. An unconditional instruction used to break the sequence of operations.

Glossary of Instructions (Continued)

<u>TERM</u>	<u>DEFINITION</u>
icp x	if the contents of the MRA are negative, take the next instruction from register x and continue from there. If the contents of the MRA are positive, take the next instruction in sequence. A conditional instruction used to break the sequence of operations.
isc x	select counter number x. Without this instruction counter zero will be used whenever +C appears.
icr x	cycle reset. Sets index register of counter to zero and criterion register to x.
ict x	cycle transfer. Increases contents of index register by one. If contents of index register greater or equal to contents of criterion register, set index register to zero and do next instruction in sequence. If contents of index register less than contents of criterion register, take next instruction from register x.
ici x	increase the contents of index register by number x.
icd x	decrease the contents of index register by number x.

Glossary of Instructions (Continued)

<u>TERM</u>	<u>DEFINITION</u>
iti x	transfer the right eleven digits of the index register into the right eleven digits of register x.
iat x	add contents of index register to the contents of register x and store the result in the index register and register x.
ita x	replace the address section of the instruction in register x with the address that is one more than the address of the register containing the last <u>isp</u> or <u>icp</u> if the contents of the MRA are negative.
iTOA	record the contents of MRA on direct printer (typewriter).
iMOA	record the contents of MRA on magnetic tape for delayed printing.
iSOA	record the contents of MRA on oscilloscope for photographing.
iFOR	this instruction provides an automatic device for obtaining a suitable layout of output data in columns, lines or blocks.

Note: The above output instructions are usually followed by a series of letters and numbers that indicate the desired form and arrangement of the output.

Computer Logic

The program starts at 1 in the main program by entering the data, program and preset parameters. The initial calculations are with the reactor radius equal to the value placed in register r4. With this radius the program then shifts to the flux iteration and convergence test, 4. In this routine parameters are set to control the number of iterations allowed at one radius and to traverse the computer thru all points in the reactor. With a value of j and m, defining a point in the reactor, the program shifts to Flux, 4.5, where the addresses on the drum of the required data are calculated, the data entered in the fast memory and the routine appropriate to the point is selected and control is shifted to that routine.

The first calculation at any point is the scalar flux for all three groups and the second is the input from fission and scattering from an upper energy group. The first point to be calculated in each cycle is the origin. This is the point of normalization of the inputs to the fast group. The normalization constant, γ , is defined as the reciprocal of the fast group input at the origin. With this constant all other inputs (all groups and all points) are normalized.

As explained in the previous section of the procedure certain quadrants of the flux are required for each point depending on its location. Prior to sequencing to another point all fluxes appropriate are calculated.

The fast source input is then compared to the previous fast source input. If the new value is not within a set specified fractional value of the previous value a negative quantity is introduced into s_{10} .

When the entire grid has traversed a check is made to see if the maximum allowable attempts at source convergence have been made. If this maximum number has been made the computer will stop. If not a check will be made of switch s_{10} . As indicated in routine 4, another cycle will be made thru the reactor if s_{10} is negative, and if s_{10} is positive control will return to the main program to repeat the above procedure based on the radius placed in r_3 .

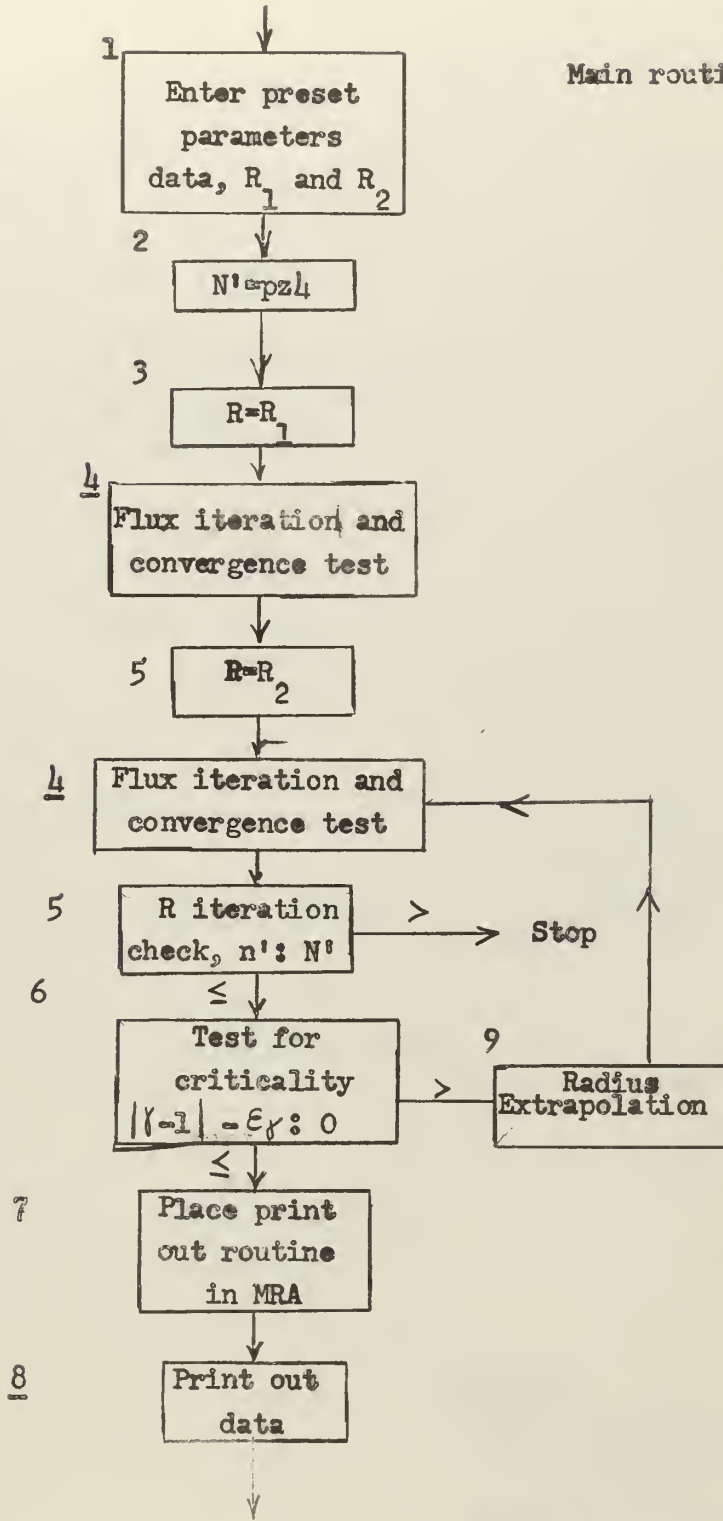
When a gamma is calculated for this radius, control is again returned to the main program but at a different point. All subsequent returns are at this point. A check is made to determine if the maximum allowable number of iterations of radii have been attempted.

A test for criticality is made to see if gamma is within a specified tolerance of unity. If it is not, a new radius is found by extrapolation or interpolation based on the most recent values of R and γ and the best previous value of R and γ . In this determination of radius, the gamma that is closer to unity and the associated radius are preserved for use in later extrapolation. The previously outlined calculations are repeated with this radius. If γ lies within the specified tolerance of unity control remains in the main routine and the print-out routine, which has been stored in the drum, is entered into the fast memory.

The computer is then sequenced thru the grid calculating and printing out scalar fluxes at each point for each group.

If the flux pattern is not as desired the properties of the regions are changed and a new tape of data cut. The program is then run again to determine the flux pattern.

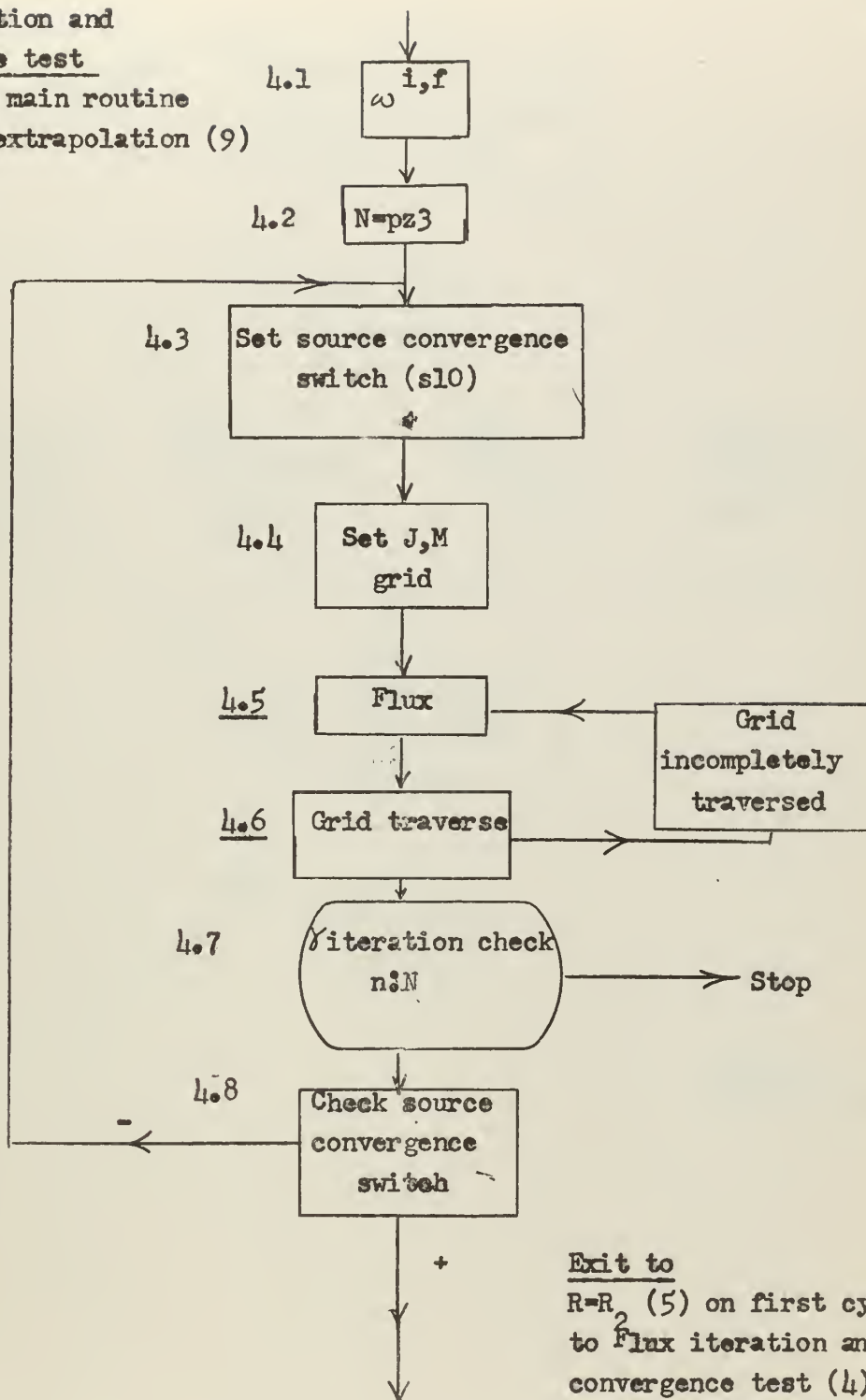
Main routine



Flux iteration and convergence test

Enter from main routine or radius extrapolation (9)

4

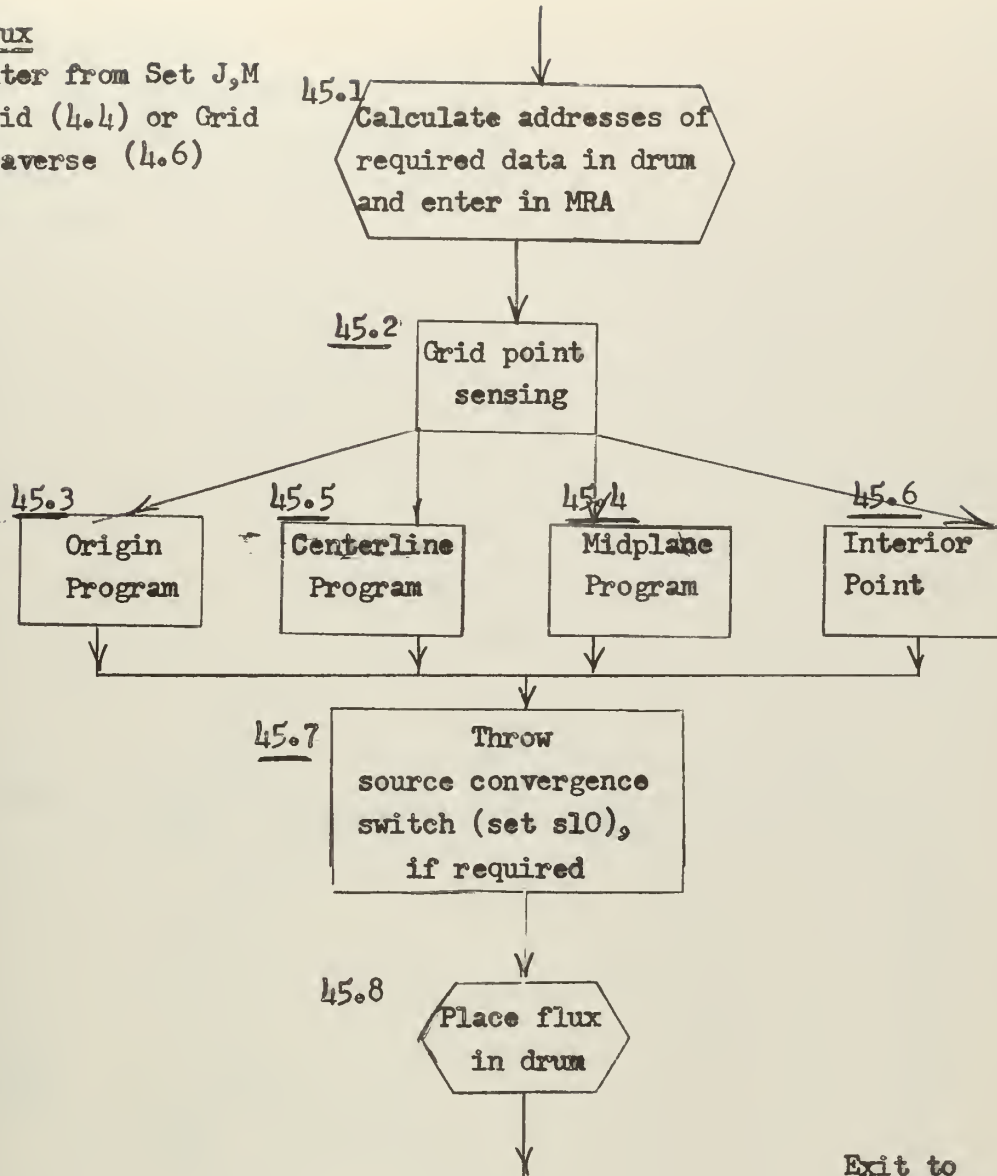


Exit to $R=R_2$ (5) on first cycle, to Flux iteration and convergence test (4) on subsequent cycles

Flux

Enter from Set J,M
grid (4.4) or Grid
Traverse (4.6)

4.5

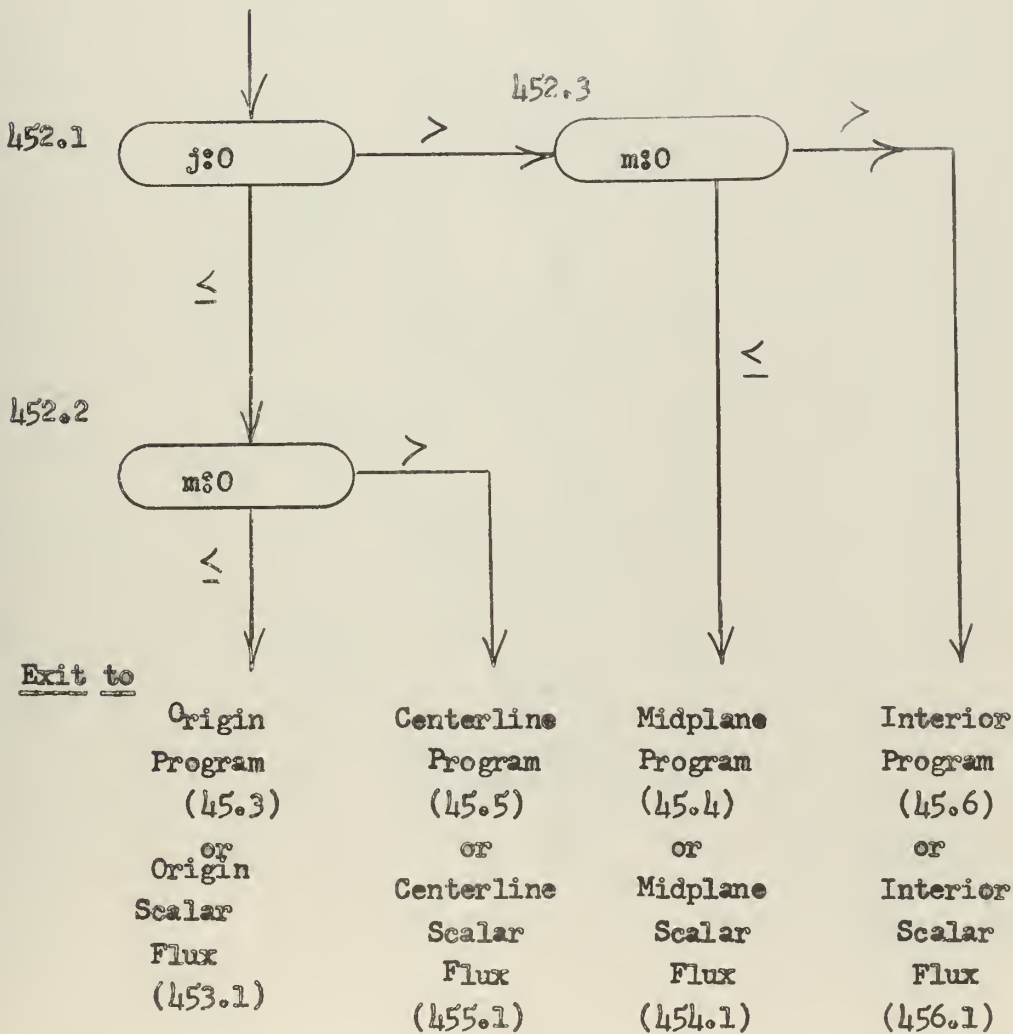


Exit to
Grid traverse (4.6)

Grid Point Sensing

45.2

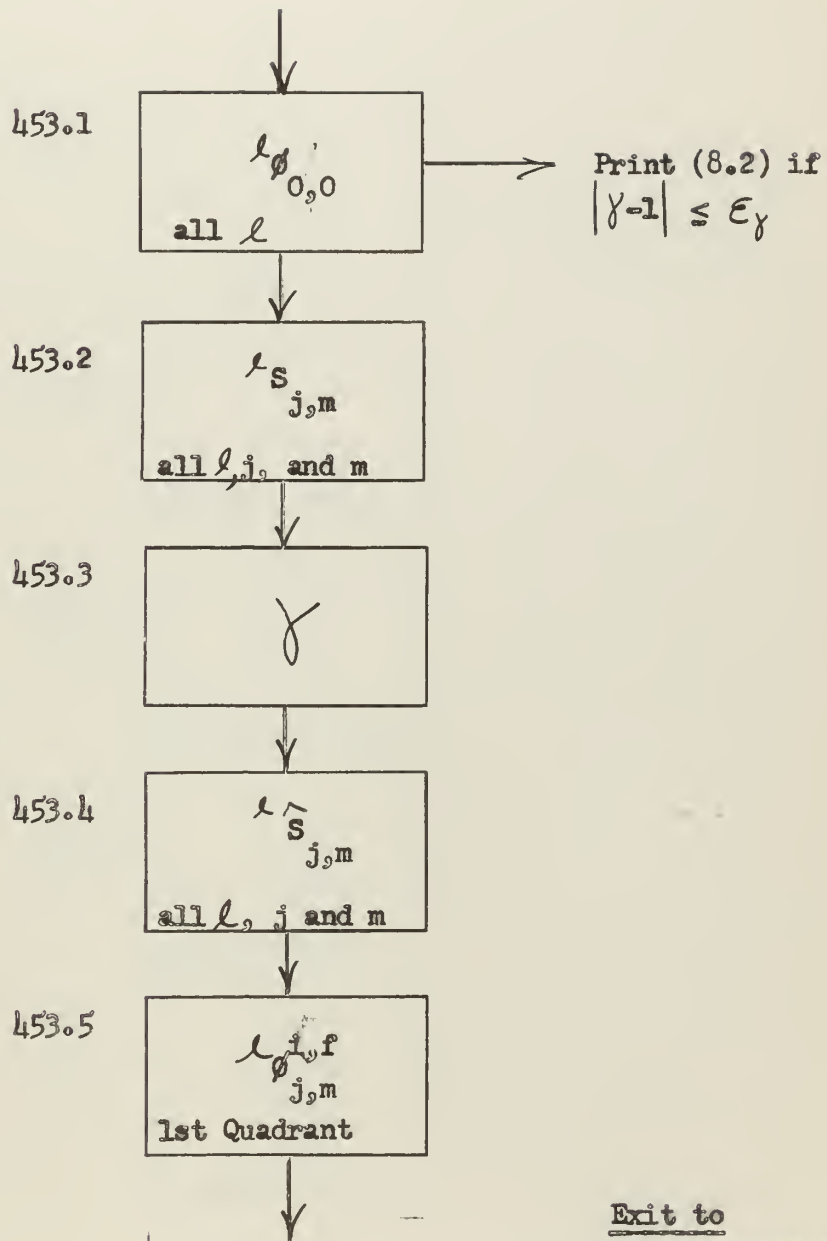
Enter from Address calculation and data entry (45.1) or calculate drum address of required data and enter data (8.1)



Origin program

45.3

Enter from Grid Point
Sensing (45.2), origin
exit

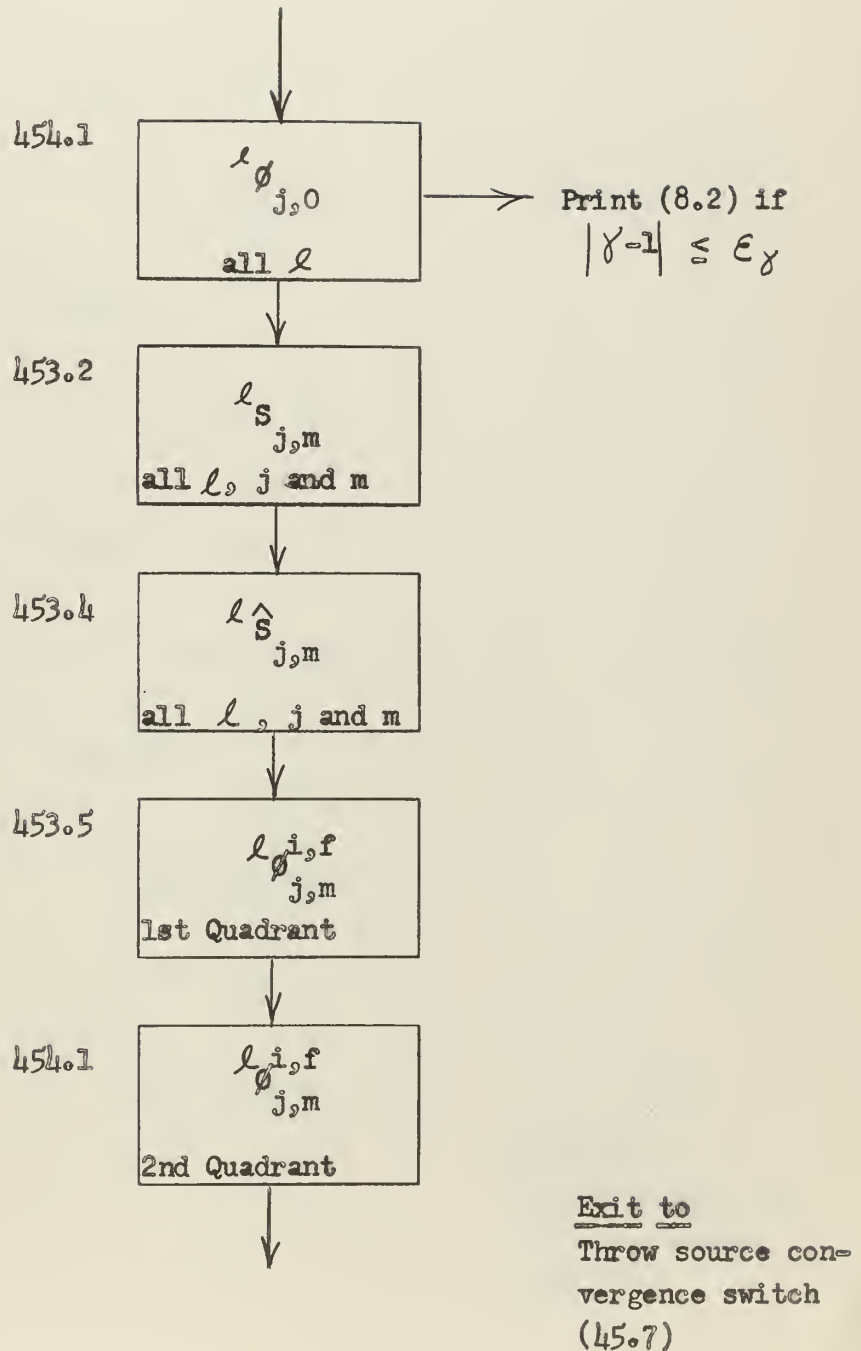


Exit to
Throw source con-
vergence switch
(45.7)

Midplane program

45.4

Enter from Grid Point
Sensing (45.2), midplane
exit



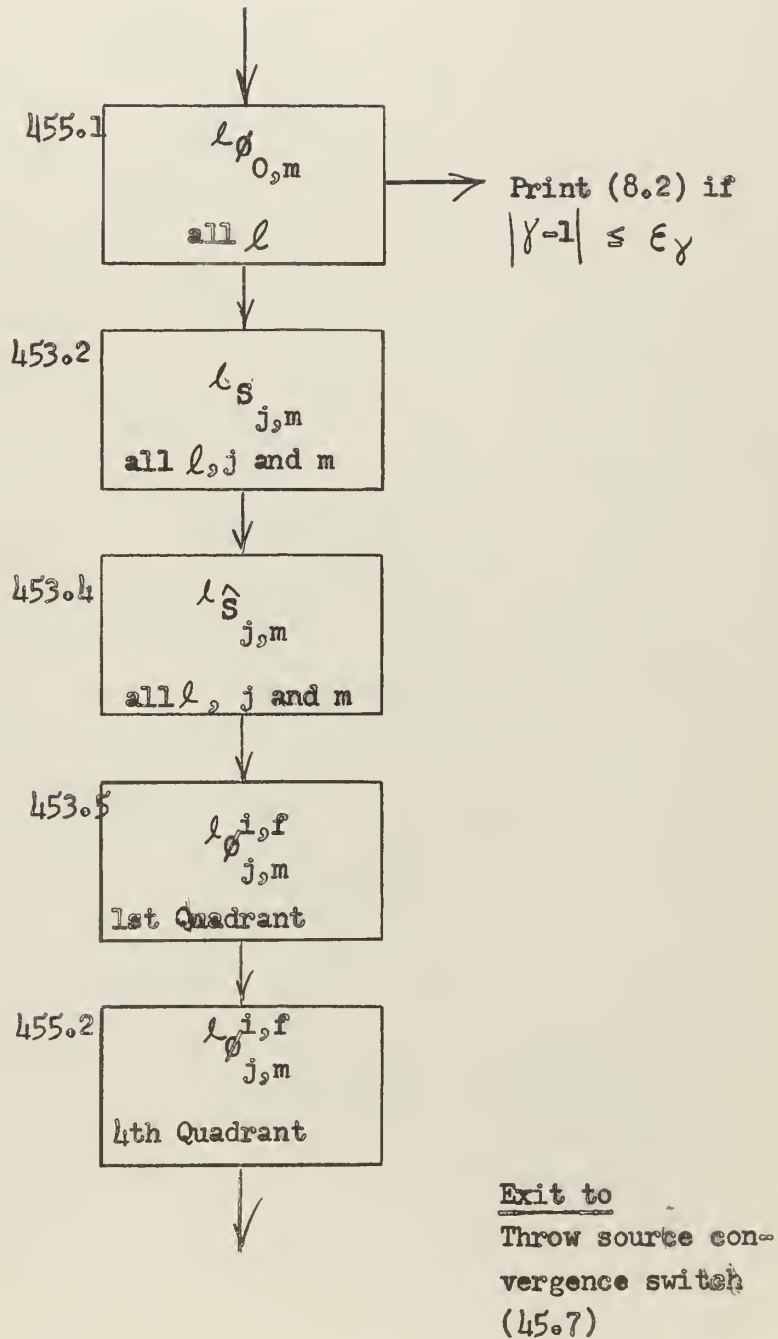
Centerline program

45.5

Enter from Grid Point

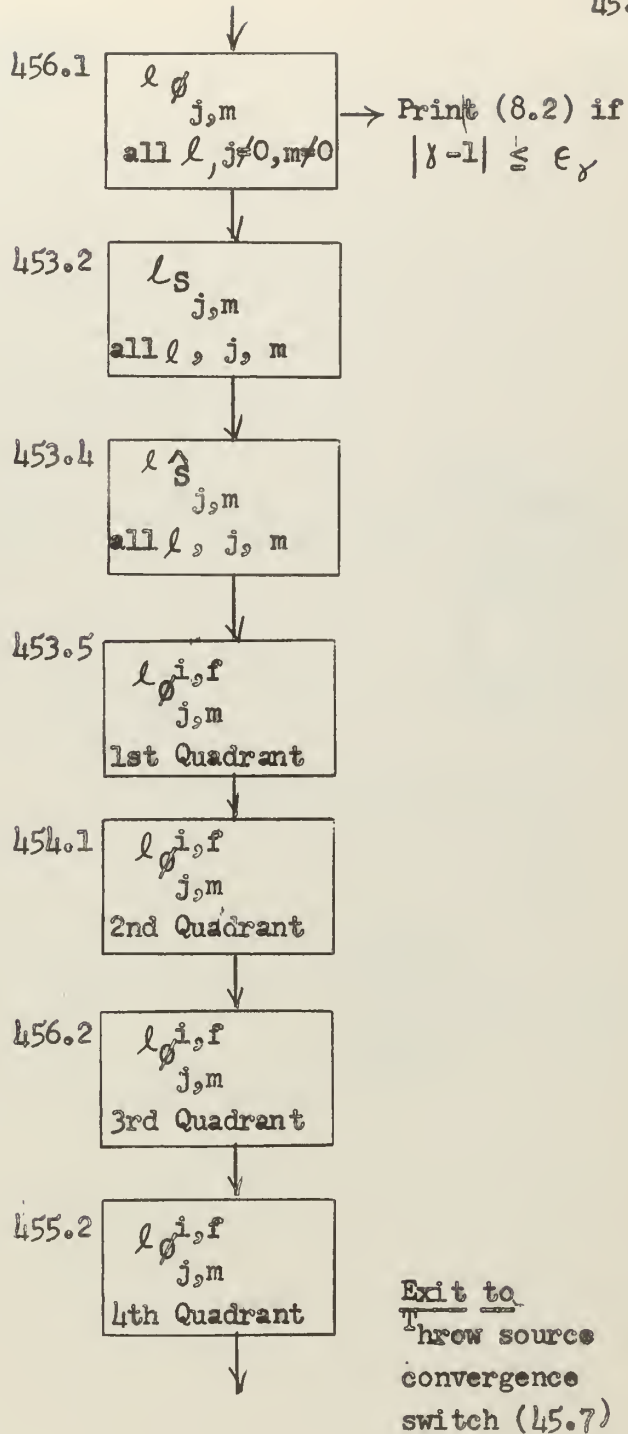
Sensing (45.2), centerline

exit



Interior point program
 Enter from Grid Point
 sensing (45.2), general
 point exit

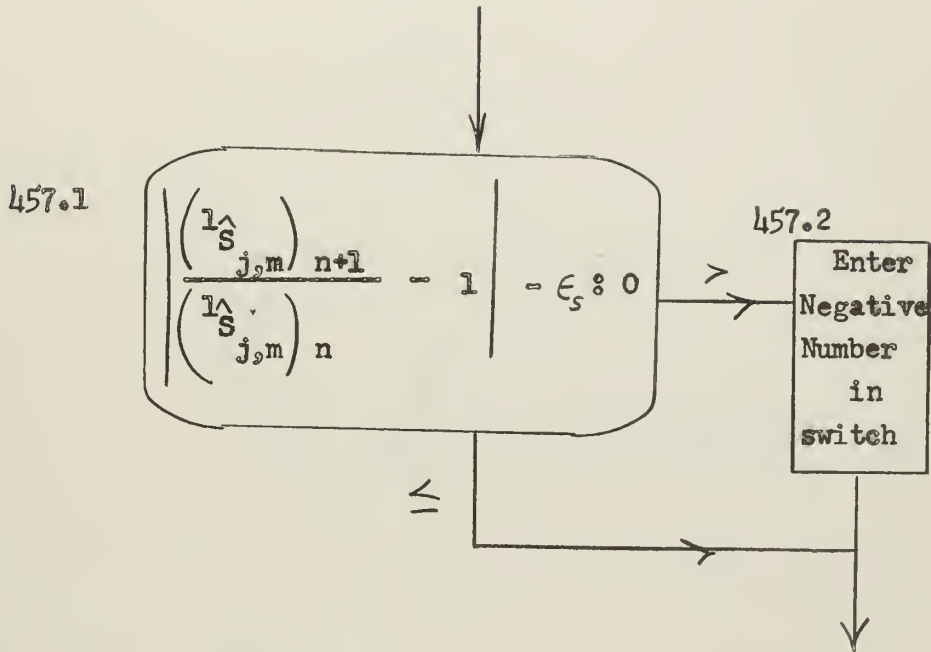
45.6



Throw source convergence switch

Enter from origin, midplane,
centerline or interior point
programs (45.3 to 45.6)

45.7



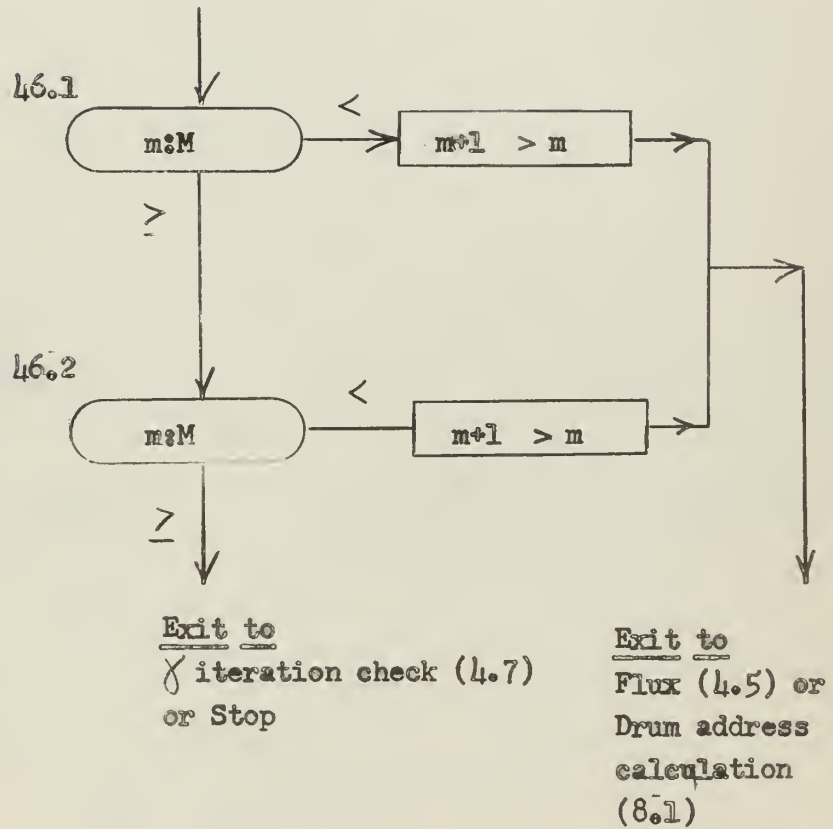
Exit to

Place flux in drum
(45.8)

Grid Traverse

4.6

Enter from Flux (4.5)
or Print (8.2)

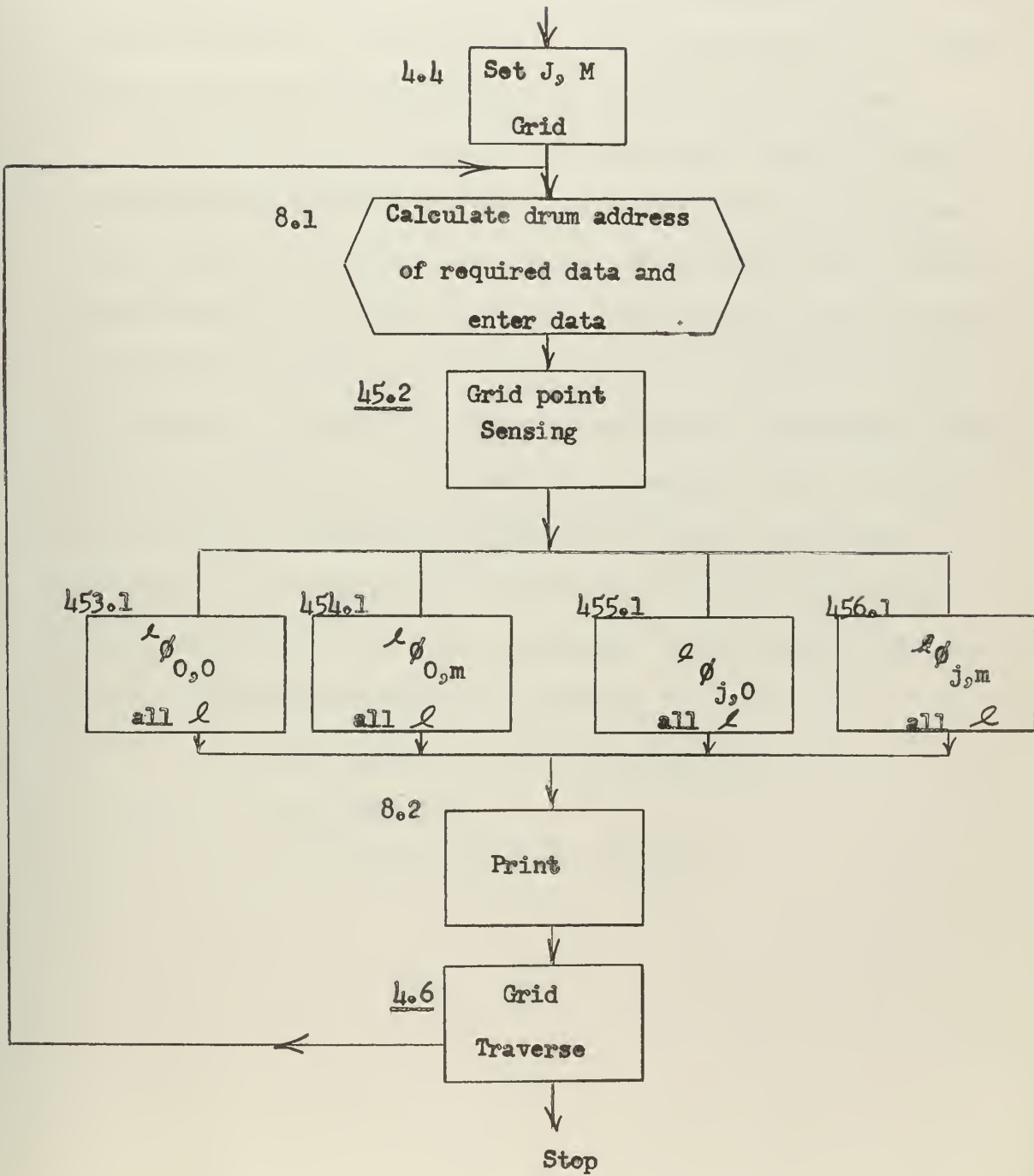


Print out routine

8

Enter from place print

out routine in MRA (7)



RESULTS

Numerous runs of subroutines and of the entire program were made on Whirlwind with the program contained in Appendix B. This program was adapted for a reactor radius of 75 and 95 centimeters and a height of 168 centimeters. The grid placed on the reactor was a 6 by 6 grid giving for the first trial radius a grid size 15 cm by 34 cm.

The final run had programmed a first trial radius of 75 cm; the maximum allowable number of iterations with a constant radius was twenty and the tolerance of convergence of the fast source input was three percent. The computer stopped, as programmed, upon completion of twenty iterations at a set radius when the fast neutron input at all points was not within the tolerance specified. The computer running time per iteration is two minutes. The computed values of scalar flux at each point and energy group were printed out for each iteration. The normalizing constant decreased from 0.27 for the first iteration to 0.01 for the twentieth. Some values of the scalar flux at the following point were negative

- a) interface of core and reflector,
- b) reflector,
- c) outside boundary of reflector.

DISCUSSION OF RESULTS

The negative values computed for scalar fluxes are known to be unrealistic. The possible causes of this error are

- 1) improper programming of the theoretically derived flux equation,
- 2) grid size being excessive.

The theoretical background for the flux equations is sound, and these equations have been programmed for other geometrical arrangements on other computers and give correct flux distributions [11]. The need for a complete hand calculation is well realized, but due to lack of time these hand calculations were not completed fully. Therefore, there is no reason to assume or expect that these equations have been programmed correctly. To accomplish the required hand calculations, computer verification of the hand calculations, and program trouble shooting it is estimated that another three to four weeks full time would be required.

For accuracy in evaluation of neutron transport equations, the grid spacing should be less than or approximately equal to the mean free path of the neutron. If this condition does not prevail, the accuracy and reliability of the calculations decrease rapidly.

A two fold increase in the number of grid points will increase the time of computation by a factor of eight. Any increase in the time required for one iteration thru the reactor is not allowable since now two minutes are required.

The code developed is in the 'interpretative' or 'floating point' mode for Whirlwind, for the reason that this mode is simple to learn and to trouble shoot. Whirlwind has also available a 'Whirlwind' or 'fixed point' mode in which the speed is increased by a factor of roughly 35. Use of the fixed point mode would make the two dimensional S_2 calculation on the Whirlwind computer practical.

CONCLUSIONS

- 1) The code as presently programmed is incorrect.
- 2) The grid spacing as used is much too coarse. This difficulty is not inherent in the coded program.
- 3) The mode of computer operation is unsuitable for an iterative process such as this because it results in a code that progresses too slowly.

RECOMMENDATIONS

- 1) Prior to the use of any computer a formal or semi-formal course in programming should be completed.
- 2) A grid spacing comparable to the mean free path of the neutron should be selected.
- 3) The fastest possible mode of calculation on the computer available should be used.
- 4) For a coded program of this difficulty a faster machine with a much larger fast memory is very desirable.

Appendix A

Development of Equations Using S_2 Approximation to the
Boltzman Equation for Cylindrical Geometry

Equations will be developed for the time independent case using cylindrical geometry.

The vector flux at a point in space is a function of two variables; one describes distance along the vertical axis of the cylinder, and the other describes radial distances. The axial variable will be designated as μ and the radial variable as η . From Figure I it can be seen that:

$$\eta = \cos \Psi$$

$$\mu = \cos \Theta$$

$$\sin \Theta = \sqrt{1 - \mu^2}$$

$$\sin \Psi = \sqrt{1 - \eta^2}$$

and where $\hat{\Omega}$ represents the unit vector in which direction neutrons

travel at a point. Therefore,

$$\nabla \phi = \hat{i}_r \frac{\partial \phi}{\partial r} + \hat{i}_\theta \frac{1}{r} \frac{\partial \phi}{\partial \theta} + \hat{i}_z \frac{\partial \phi}{\partial z}$$

$$\text{and } \hat{\Omega} = \hat{i}_r \sin \Theta \cos \Psi -$$

$$\hat{i}_\theta \sin \Theta \sin \Psi \frac{\partial \phi}{\partial \psi} + \hat{i}_z \cos \Theta.$$

Therefore,

$$\hat{\Omega} \cdot \nabla \phi = \sin \Theta \cos \Psi \frac{\partial \phi}{\partial r} - \frac{\sin \Theta \sin \Psi}{r} \frac{\partial \phi}{\partial \psi} + \cos \Theta \frac{\partial \phi}{\partial z} \quad (1)$$

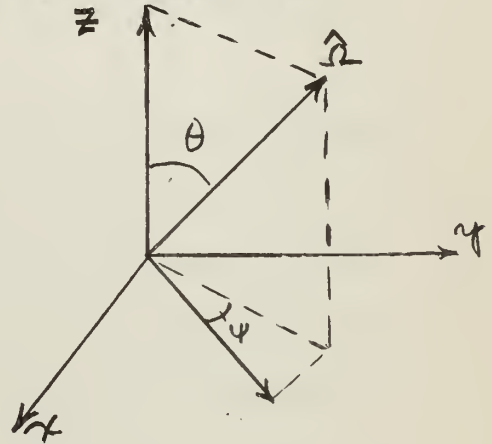
If we substitute the expression for η and μ which were derived from Figure I in equation 1, we find

$$\hat{\Omega} \cdot \nabla \phi = \eta \sqrt{1 - \mu^2} \frac{\partial \phi}{\partial r} + \frac{(1 - \eta^2)}{r} \sqrt{1 - \mu^2} \frac{\partial \phi}{\partial \eta} + \mu \frac{\partial \phi}{\partial z} \quad (2)$$

The quantity $\hat{\Omega} \cdot \phi$ represents the leakage of neutrons out of an element of volume per unit time in the direction $\hat{\Omega}$ per unit solid angle.

Figure I

Components of neutron flux at a space point.



In the time independent case this leakage plus the rate of loss of neutrons, $\Sigma \phi$, per unit time must equal the source strength, S. In problems in which cross sections vary with energy and neutrons alter their speeds, the multigroup approximation of the Boltzmann equation is introduced: [6]

$$\frac{\partial^l \phi}{\partial r} + \frac{1-\eta^2}{r} \frac{\partial^l \phi}{\partial \eta} + \frac{\mu}{\sqrt{1-\mu^2}} \frac{\partial^l \phi}{\partial z} = \frac{S}{\sqrt{1-\mu^2}} - \frac{l \Sigma^l \phi}{\sqrt{1-\mu^2}}, \quad (3)$$

where $^l \Sigma$ is the total cross section within the energy group $^l Q$ and $^l S$ is the neutron source term which results from fission and energy degradation of the total vector flux at the point. Slowing down is incorporated into the source term.

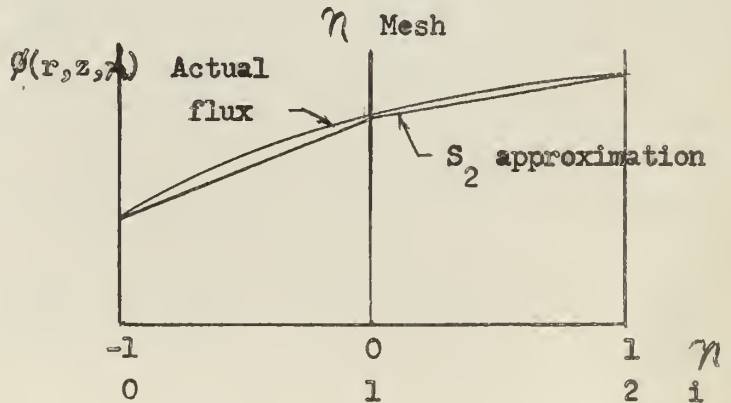
Equation 3 is identical in form to the time dependent multigroup equation for spherical geometry in which the z term is replaced by a term representing the rate of change of flux with respect to time.

The S_n approximation is a method of evaluating integrals over a variable in which the variable is quantized into n intervals and intermediate values of a function of the variable are approximated by a straight line between points at the boundaries of the intervals.

We shall consider η space as the first variable which will be integrated. The flux at any point

(r, z, η) in Figure II can be represented by the following equation where i represents the interval number.

Figure II



$$\phi(r, z, \eta) = \frac{\eta - \eta_{i-1}}{\eta_i - \eta_{i-1}} \phi(r, z, \eta_i) + \frac{\eta_i - \eta}{\eta_i - \eta_{i-1}} \phi(r, z, \eta_{i-1}) \quad (4)$$

For the S_2 approximation $i = +1, 0,$ and -1 . If we denote $\eta_i - \eta_{i-1}$ as Δ_i and $\phi(r, z, \eta_i)$ as ϕ^i we obtain

$$\phi = \frac{(\eta - \eta_{i-1})}{\Delta_i} \phi^i + \frac{(\eta_i - \eta)}{\Delta_i} \phi^{i-1}$$

If we substitute this in equation 3, we find

$$\begin{aligned} & \frac{\eta(\eta - \eta_{i-1})}{\Delta_i} \frac{\partial^2 \phi^i}{\partial r^2} + \frac{\eta(\eta_i - \eta)}{\Delta_i} \frac{\partial^2 \phi^{i-1}}{\partial r^2} + \frac{(1-\eta^2)(\eta - \eta_{i-1})}{r \Delta_i} \frac{\partial^2 \phi^i}{\partial \eta^2} \\ & + \frac{(1-\eta^2)(\eta_i - \eta)}{\Delta_i} \frac{\partial^2 \phi^{i-1}}{\partial \eta^2} + \frac{\mu(\eta - \eta_{i-1})}{\sqrt{1-\mu^2} \Delta_i} \frac{\partial^2 \phi^i}{\partial z^2} + \frac{\mu(\eta_i - \eta)}{\Delta_i} \frac{\partial^2 \phi^{i-1}}{\partial z^2} \\ & = \frac{S}{\sqrt{1-\mu^2}} - \frac{L_{\Sigma}(\eta - \eta_{i-1}) \phi^i}{\Delta_i \sqrt{1-\mu^2}} - \frac{L_{\Sigma}(\eta_i - \eta) \phi^{i-1}}{\Delta_i \sqrt{1-\mu^2}} \end{aligned} \quad (5)$$

Since the source, S , is isotropic, integrating this expression over η from η_{i-1} to η_i , we obtain n equations in the form

$$\begin{aligned} & b_i \frac{\partial^2 \phi^i}{\partial r^2} + c_i \frac{L_{\Sigma} \phi^i}{r} + h \frac{\partial^2 \phi^i}{\partial z^2} + g_{\Sigma} L_{\Sigma} \phi^i + d_i \frac{\partial^2 \phi^{i-1}}{\partial r^2} - c_i \frac{L_{\Sigma} \phi^{i-1}}{r} + \\ & + h \frac{\partial^2 \phi^{i-1}}{\partial z^2} + g_{\Sigma} L_{\Sigma} \phi^{i-1} = e_i g S \end{aligned} \quad (6)$$

where

$$b_i = \frac{2\eta_i + \eta_{i-1}}{3}, \quad (7a)$$

$$c_i = \frac{2}{3\Delta_i} (3 - \eta_i^2 - \eta_i - \eta_{i-1} - \eta_{i-1}^2), \quad (7b)$$

$$d_i = \frac{\eta_i + 2\eta_{i-1}}{3}, \quad (7c)$$

$$h = \frac{\mu}{\sqrt{1-\mu^2}}, \quad (7e)$$

$$e_i = 2 \quad (7d)$$

$$g_f = \frac{1}{\sqrt{1-\mu^2}}. \quad (7f)$$

A different method of finite differencing is used for the other directional variable, μ , since use of an expression analogous to (4) for μ in the integration of the equation (5) with respect to μ would result in an extremely complex expression. Instead, the equation (5) is directly quantized with respect to μ giving the results

$$b_i \frac{\partial^2 \phi^{i,f}}{\partial r^2} + c_i \frac{\partial \phi^{i,f}}{\partial r} + h_f \frac{\partial^2 \phi^{i,f}}{\partial z^2} + g_f \Sigma \phi^{i,f} \quad (8)$$

$$d_i \frac{\partial^2 \phi^{i-1,f}}{\partial r^2} - d_i \frac{\partial \phi^{i-1,f}}{\partial r} + h_f \frac{\partial^2 \phi^{i-1,f}}{\partial z^2} + g_f \Sigma \phi^{i-1,f} = e_i g_f S$$

We may now proceed to evaluate the source integral at a point. For isotropic scattering and three energy groups of neutrons

$$S(r,z) = \tau^p \sum_{k=1}^3 \sum_f k \phi(r,z) + \sum_{k=1}^3 \tau^k \phi(r,z), \quad (9)$$

where

$$\phi^k(r,z) = \frac{1}{2\pi} \int_{-1}^{+1} d\mu \int_{-1}^{+1} \phi(r,z,\mu,\eta) \frac{d\eta}{\sqrt{1-\eta^2}}$$

As in equation (4) $\phi(r,z,\mu,\eta)$ is approximated by

$$\phi(r,z,\mu) = \frac{\eta - \eta_{i-1}}{\Delta_i} \phi^i(r,z,\mu) + \frac{\eta_i - \eta}{\Delta_i} \phi^{i-1}(r,z,\mu),$$

and

$$\begin{aligned}
 {}^k\phi(r, z, \mu) = & \frac{1}{2\pi} \sum_{i=0}^2 \left\{ \phi^i \left[\int_{\eta_{i-1}}^{\eta_i} \frac{\eta d\eta}{\sqrt{1-\eta^2}} - \eta_{i-1} \int_{\eta_{i-1}}^{\eta_i} \frac{d\eta}{\sqrt{1-\eta^2}} \right] \right. \\
 & \left. + \phi^{i-1} \left[\eta_i \int_{\eta_{i-1}}^{\eta_i} \frac{d\eta}{\sqrt{1-\eta^2}} - \eta_{i-1} \int_{\eta_{i-1}}^{\eta_i} \frac{\eta d\eta}{\sqrt{1-\eta^2}} \right] \right\} . \quad (10)
 \end{aligned}$$

Upon integration

$$\begin{aligned}
 {}^k\phi(r, z, \mu) = & \frac{1}{2} \sum_{i=0}^2 \left[\frac{\phi^i}{\pi \Delta_i} \left\{ \sqrt{1-\eta_{i-1}^2} - \sqrt{1-\eta_i^2} + \eta_{i-1} \sin^{-1} \eta_{i-1} \right. \right. \\
 & \left. \left. - \eta_{i-1} \sin^{-1} \eta_i \right\} + \frac{1}{2} \left[\frac{\phi^{i-1}}{\pi \Delta_i} \left\{ \eta_i \sin^{-1} \eta_i \right. \right. \right. \\
 & \left. \left. - \eta_i \sin^{-1} \eta_{i-1} + \sqrt{1-\eta_i^2} - \sqrt{1-\eta_{i-1}^2} \right\} \right] . \quad (11)
 \end{aligned}$$

If values of $\eta_i = +1, 0$ and -1 (S_2 approximation) are substituted in the above expression, we find that

$${}^k\phi(r, z, \mu) = \frac{1}{2\pi} \sum_{i=0}^2 P_i \phi^i(r, z, \mu) , \quad (12)$$

where $P_0 = 0.3183$, $P_1 = 0.3634$, and $P_2 = 0.3183$.

By use of the Gauss quadrature method [8] the remaining integral can be evaluated in the form of a sum of the products of coefficient times the vector flux where μ_f values are the $n+2$ positive roots of

$$P_{n+2}(\mu_f) = 0 .$$

The table below lists the values of μ_f and the coefficient H_f used

f	μ_f	H_f
0	-0.8611363	0.3478548
1	-0.3399810	0.6521452
2	+0.3399810	0.6521452
3	+0.8611363	0.3478548

Therefore, since $H_0 = H_3$ and $H_1 = H_2$ the scalar flux at a point is

$${}^k\phi(r,z) = \sum_{f=0}^3 H_f \sum_{i=0}^2 P_i {}^k\phi_i{}^f(r,z), \quad (13)$$

and the source term equals

$$S = \sum_p \sigma_p^l \sum_{k=1}^3 k \sum_f {}^k\phi(r,z) + \sum_{k=1}^3 \sigma_{\gamma}^k {}^k\phi(r,z), \quad (14)$$

where

σ_p^l is the probability that a fission neutron will have an energy l and σ_{γ}^k is the probability that a neutron having an energy k will be scattered into the energy group l , $k \geq l$.

The scattering probability requires further explanation. We define μ_{lk} as the probability that a neutron which is scattered in the k^{th} group will land in the l^{th} group, k as the energy group in which the scattering occurs (1 for fast, 2 for epithermal and 3 for thermal) and l as the energy group into which the neutron is scattered. The energy range covered by scattered neutrons is $E - \alpha E$, and the probability of landing in the interval ΔE is $\frac{\Delta E}{E - \alpha E}$. Let E be the energy of the neutron before scattering.

$$E_{k+1} < E < E_k$$

The neutron flux per unit energy interval is $\frac{d\phi}{dE}$.

Then $\frac{d\phi}{dE} dE \Sigma_s(E)$ is the number of neutrons scattered per unit volume and per unit time, [9] but

$$\frac{d\phi}{dE} = \frac{q}{\Sigma_s(E)}$$

Therefore, $\frac{d\phi}{dE} dE \Sigma_s(E) = \frac{q}{\Sigma_s(E)} \frac{dE}{E}$ (15)

$\frac{E_{l+1} - E_l}{E_l(1-\alpha)}$ is the probability a neutron will land in the l th group.

${}^{\ell k} \gamma$ is the number of neutrons scattered in the k th group that land in the l th group. Since the absorption, sources and leakage in a large thermal reactor are small, the slowing down density q , is rather constant over an energy group. This fact is used to compute the number of neutrons scattered from the group k to group l . [10]

The equations for each case are listed below.

Case	Equation
1. $E_{l+1} > \alpha E_k$	${}^{\ell k} \gamma = \frac{\Sigma_s^k}{(1-\alpha) \ln(E_k/E_{k+1})} (E_{l+1} - E_l) \left(\frac{1}{E_{k+1}} - \frac{1}{E_k} \right)$ (16a)
2. $E_l > \alpha E_k$ $\alpha E_k > E_{l+1} > \alpha E_{k+1}$	${}^{\ell k} \gamma = \frac{\Sigma_s^k}{(1-\alpha) \ln(E_k/E_{k+1})} \left[\frac{E_{l+1} - E_l}{E_{k+1}} + \alpha - \frac{E_l}{E_k} - \alpha \ln \left(\frac{\alpha E_k}{E_{l+1}} \right) \right]$ (16b)

$$\begin{aligned}
3. \quad E_l > \alpha E_k & \quad \ell^k \gamma = \frac{\sum_s^k}{(1-\alpha) \ln(E_k/E_{k+1})} \left\{ E_l \left[\frac{1}{E_{k+1}} - \frac{\alpha}{E_k} \right] \right. \\
E_{l+1} < \alpha E_{k+1} & \quad \left. + \alpha \ln \frac{E_k}{E_{k+1}} \right\} \quad (16c)
\end{aligned}$$

$$\begin{aligned}
4. \quad \alpha E_{k+1} < E_l < \alpha E_k & \quad (16d) \\
E_{l+1} < \alpha E_{k+1} & \quad \ell^k \gamma = \frac{\sum_s^k}{(1-\alpha) \ln(E_k/E_{k+1})} \left\{ E_l \left[\frac{1}{E_{k+1}} - \frac{\alpha}{E_l} - \alpha \ln \frac{E_l}{\alpha E_{k+1}} \right] \right\}
\end{aligned}$$

$$5. \quad E_l < \alpha E_{k+1} \quad \ell^k \gamma = 0 \quad (16e)$$

For the cases where the scattering occurs within the energy group we find

$$\begin{aligned}
E_{l+1} > \alpha E_l & \quad \ell^l \gamma = \frac{\sum_s^k}{1-\alpha} \left[1 + \frac{E_{l+1} - E_l}{E_l \ln E_l/E_{l+1}} \right] \quad (16f) \\
E_{l+1} < \alpha E_l & \quad \ell^l \gamma = \left[1 - \frac{\xi}{\ln E_l/E_{l+1}} \right] \sum_s^k
\end{aligned}$$

We may now proceed to evaluate equation 4 by integrating over the r and z variables. However, before doing so we must consider the direction in which integration should be carried out. It is always best to integrate in the direction of neutron flow since any errors produced in the process of integration will be diminished as we proceed further in this direction.

If integration is carried out in a direction opposite to that of neutron flow, any errors produced will be magnified. [7]

We must now difference our equations with respect to r and z.

We can arrange terms in equation (8) in order to facilitate the integration. Thus equation (8) becomes

$$\left[h_f \frac{\partial}{\partial z} + b_i \frac{\partial}{\partial r} \right] \left[l_{\phi}^{i,f} + l_{\phi}^{i-1,f} \right] = - \left[g_f l_{\Sigma} + \frac{c_i}{r} \right] l_{\phi}^{i,f} - \left[g_f l_{\Sigma} - \frac{c_i}{r} \right] l_{\phi}^{i-1,f} + \left[b_i - d_i \right] \frac{\partial l_{\phi}^{i-1,f}}{\partial r} + e_i g_f l_s \quad (17)$$

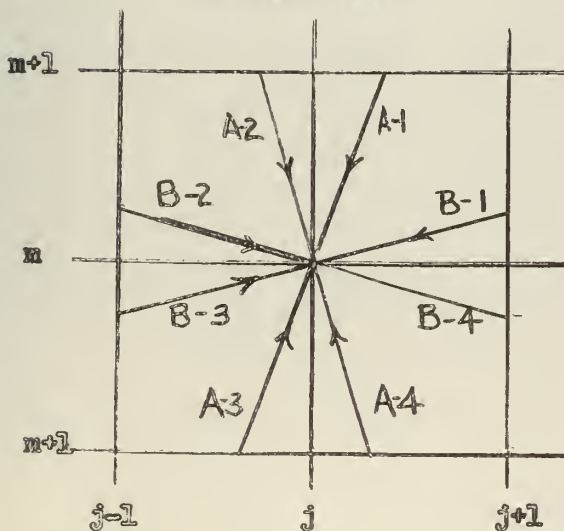
We must now further divide the results of the integration into two separate cases:

Case A - The component of velocity in the r direction is less than the component of velocity in the z direction.

Case B - The component of velocity in the r direction is greater than the component of velocity in the z direction.

If we consider Case A, and difference r and z as illustrated in the figure below, we obtain four distinct subcases. These will be classified in

Figure III
Position Mesh



accordance with the direction of travel being determined by the signs of b_i and h_f . A positive sign indicates a direction of travel away from the center of the core. The various cases are listed in Table I.

Table I

In all A cases $|b_i| < |h_f|$

In all B cases $|b_i| > |h_f|$

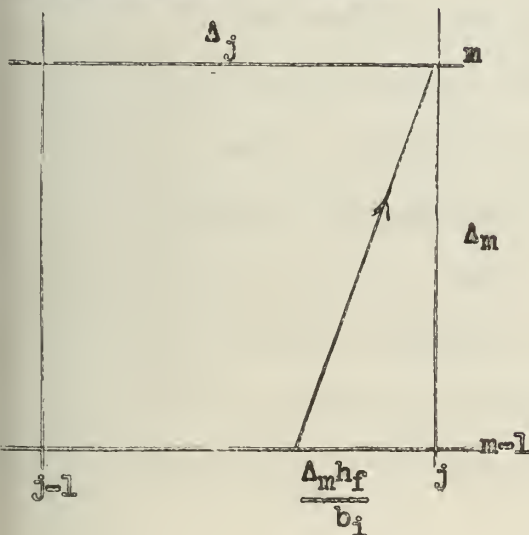
<u>Subcase</u>	$\frac{b_i}{h_f}$	$\frac{h_f}{b_i}$
A1, B1	-	-
A2, B2	+	-
A3, B3	+	+
A4, B4	-	+

Considering subcase A-3 and integrating equation 5 and knowing that

$$\int_{r_{j-1}}^{r_j} r \phi^{i,f}(r,z) dr = \frac{\Delta_j}{2} \left[\phi_{j,m}^{i,f} + \phi_{j-1,m}^{i,f} \right]$$

Figure IV

Position Mesh for Case A



and

$$\int_{z_{m-1}}^{z_m} \phi^{i,f}(r,z) dz = \frac{\Delta_m}{2} \left[\phi_{j,m}^{i,f} + \phi_{j,m-1}^{i,f} \right]$$

where $\Delta_j = r_j - r_{j-1}$ and $\Delta_m = z_m - z_{m-1}$

we obtain the following equation

$$\begin{aligned}
& \left| \frac{h_f}{\Delta_m} \right| \left[l_{\phi_{j,m}^{i,f}} + l_{\phi_{j,m}^{i-1,f}} - l_{\phi_{j,m-1}^{i,f}} - l_{\phi_{j,m-1}^{i-1,f}} \right] + \left| \frac{b_i}{\Delta_j} \right| \left[l_{\phi_{j,m-1}^{i,f}} \right. \\
& \quad \left. + l_{\phi_{j,m-1}^{i-1,f}} - l_{\phi_{j-1,m-1}^{i,f}} - l_{\phi_{j-1,m-1}^{i-1,f}} \right] = - \left[g_f \ l_{\Sigma_{j-1}^{m-1}} \right. \\
& \quad \left. + \frac{c_i}{\bar{r}_{j-1}} \right] \left[\frac{l_{\phi_{j,m}^{i,f}} + l_{\phi_{j-1,m-1}^{i,f}}}{2} \right] - \left[g_f \ l_{\Sigma_{j-1}^{m-1}} - \frac{c_i}{\bar{r}_{j-1}} \right] \left[\frac{l_{\phi_{j,m}^{i-1,f}} + l_{\phi_{j-1,m-1}^{i-1,f}}}{2} \right] \\
& \quad + \frac{1}{\Delta_j} \left[\frac{b_i - d_i}{2} \right] \left[l_{\phi_{j,m}^{i-1,f}} + l_{\phi_{j,m-1}^{i-1,f}} - l_{\phi_{j-1,m}^{i-1,f}} - l_{\phi_{j-1,m-1}^{i-1,f}} \right] \\
& \quad + \frac{e_{i g_f}}{2} \left[l_{S_{j,m}} + l_{S_{j-1,m-1}} \right] \tag{18}
\end{aligned}$$

Where Σ_{j-1}^{m-1} is the total cross section of the material at the center of the grid in figure 4; i.e., the material to the left and below the j,m point and \bar{r}_{j-1} is $\frac{r_j + r_{j-1}}{2}$.

The radial direction of integration is determined by the sign of η , i.e., by the sign of the b_i or d_i . Consequently, the signs of η , b_i , d_i , and Δ_j all change together. Likewise, the direction of integration axially is determined by the sign of μ , i.e., by the sign of h_f . Consequently the signs of μ , h_f and Δ_m all change together. If we use absolute magnitude signs for these quantities all Case A equations will be identical insofar as the signs of various terms are concerned. Similarly, all Case B equations will be identical.

If we now add the following expression to both sides of equation (18)

$$\frac{|b_i|}{|\Delta_j|} \left[l_{\phi_{j-1,m-1}}^{i,f} + l_{\phi_{j-1,m-1}}^{i-1,f} - l_{\phi_{j,m}}^{i,f} - l_{\phi_{j,m}}^{i-1,f} \right],$$

we obtain

$$\begin{aligned} & \left[\frac{|h_f|}{|\Delta_m|} - \frac{|b_i|}{|\Delta_j|} \right] \left[l_{\phi_{j,m}}^{i,f} + l_{\phi_{j-1,m-1}}^{i-1,f} - l_{\phi_{j,m-1}}^{i,f} - l_{\phi_{j,m-1}}^{i-1,f} \right] = \\ & - \left[g_f l_{\Sigma_{j-1}}^{m-1} + \frac{c_i}{r_{j-1}} \right] \left[\frac{l_{\phi_{j,m}}^{i,f} + l_{\phi_{j-1,m-1}}^{i,f}}{2} \right] - \left[g_f l_{\Sigma_{j-1}}^{m-1} \right. \\ & + \frac{c_i}{r_{j-1}} \left. \left[\frac{l_{\phi_{j,m}}^{i-1,f} + l_{\phi_{j-1,m-1}}^{i-1,f}}{2} \right] + \frac{|b_i| - |d_i|}{2|\Delta_j|} \left[l_{\phi_{j,m}}^{i-1,f} + l_{\phi_{j,m-1}}^{i-1,f} \right. \right. \\ & \left. \left. - l_{\phi_{j-1,m}}^{i-1,f} - l_{\phi_{j-1,m-1}}^{i-1,f} \right] + \frac{c_i g_f}{2} \left[l_{\bar{S}_{j,m}} + l_{\bar{S}_{j-1,m-1}} \right] \right. \\ & \left. + \frac{|b_i|}{|\Delta_j|} \left[l_{\phi_{j-1,m-1}}^{i,f} + l_{\phi_{j-1,m-1}}^{i-1,f} - l_{\phi_{j,m}}^{i,f} - l_{\phi_{j,m}}^{i-1,f} \right] \right]. \end{aligned} \quad (18a)$$

If this expression is multiplied by $|\Delta_j| \omega^{i,f}$ where $|\Delta_j| \omega^{i,f} = \frac{|b_i|}{|h_f|} |\Delta_m|$

we obtain

$$\begin{aligned} & |b_i| (1 - \omega^{i,f}) \left[l_{\phi_{j,m}}^{i,f} + l_{\phi_{j,m}}^{i-1,f} - l_{\phi_{j,m-1}}^{i,f} - l_{\phi_{j,m-1}}^{i-1,f} \right] = \\ & \omega^{i,f} \left[|b_i| + \frac{|\Delta_j|}{2} \left(g_f l_{\Sigma_{j-1}}^{m-1} + \frac{c_i}{r_{j-1}} \right) \right] l_{\phi_{j,m}}^{i,f} + A_3 \omega^{i,f}. \end{aligned} \quad (18b)$$

Where

$$\begin{aligned}
 A_3 = & \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_f \ell_{\Sigma_{j-1}}^{m-1} + \frac{c_i}{r_{j-1}} \right) \right] \phi_{j-1,m-1}^{i,f} \\
 & - \left[|b_i| + \frac{|\Delta_j|}{2} \left(g_f \ell_{\Sigma_{j-1}}^{m-1} - \frac{c_i}{r_{j-1}} \right) \right] \phi_{j,m}^{i-1,f} \\
 & + \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_f \ell_{\Sigma_{j-1}}^{m-1} - \frac{c_i}{r_{j-1}} \right) \right] \phi_{j-1,m-1}^{i-1,f} + \frac{i g_f}{2} |\Delta_j| \left[\bar{S}_{j,m} + \bar{S}_{j-1,m-1} \right] \\
 & + \left(\frac{|b_i| - |d_i|}{2} \right) \left[\phi_{j,m}^{i-1,f} + \phi_{j,m-1}^{i-1,f} - \phi_{j-1,m}^{i-1,f} - \phi_{j-1,m-1}^{i-1,f} \right] \quad (18c)
 \end{aligned}$$

And finally transposing terms in equation (18b) we obtain an expression for the vector flux at the point j,m .

$$\begin{aligned}
 \phi_{j,m}^{i,f} = & \frac{b_i (1-\omega^{i,f}) \left(\phi_{j,m-1}^{i,f} + \phi_{j,m-1}^{i-1,f} - \phi_{j,m}^{i-1,f} \right) + \omega^{i,f} A_3}{|b_i| + \omega^{i,f} \frac{|\Delta_j|}{2} \left(g_f \ell_{\Sigma_{j-1}}^{m-1} + \frac{c_i}{r_{j-1}} \right)} \quad (18d)
 \end{aligned}$$

By use of similar procedures for developing the equations for the three remaining A subcases we obtain the three equations on the following pages.

For subcase A-4

$$l_{\phi, j, m}^{i, f} = \frac{|b_i| (1 - \omega^{i, f}) \left(l_{\phi, j, m-1}^{i, f} + l_{\phi, j, m-1}^{i-1, f} - l_{\phi, j, m}^{i-1, f} \right) + \omega^{i, f} A_4}{|b_i| + \omega^{i, f} \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j+1}}^{m-1} + \frac{c_i}{\bar{r}_{j+1}} \right)} \quad (19a)$$

where

$$\begin{aligned} A_4 = & \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j+1}}^{m-1} + \frac{c_i}{\bar{r}_{j+1}} \right) \right] l_{\phi, j+1, m-1}^{i, f} \\ & - \left[|b_i| + \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j+1}}^{m-1} - \frac{c_i}{\bar{r}_{j+1}} \right) \right] l_{\phi, j, m}^{i-1, f} \\ & + \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j+1}}^{m-1} - \frac{c_i}{\bar{r}_{m+1}} \right) \right] l_{\phi, j+1, m-1}^{i-1, f} \\ & + \frac{g_i}{2} |\Delta_j| \left[l_{\bar{S}_{j, m}} + l_{\bar{S}_{j+1, m-1}} \right] + \frac{|b_i| - |d_i|}{2} \left[l_{\phi, j, m}^{i-1, f} + l_{\phi, j, m-1}^{i-1, f} \right. \\ & \left. - l_{\phi, j+1, m}^{i-1, f} - l_{\phi, j+1, m-1}^{i-1, f} \right]. \end{aligned} \quad (19b)$$

For subcase A-1

$$l_{\phi_{j,m}^{i,f}} = \frac{|b_i| (1-\omega)^{i,f} \left(l_{\phi_{j,m+1}^{i,f}} + l_{\phi_{j,m+1}^{i-1,f}} - l_{\phi_{j,m}^{i-1,f}} \right) + \omega^{i,f} A_1}{|b_i| + \omega^{i,f} \frac{|\Delta_j|}{2} \left(\frac{l_{\Sigma_{j+1}^{m+1}}}{g_{f,j+1}} + \frac{c_i}{\bar{r}_{j+1}} \right)} \quad (20a)$$

where

$$\begin{aligned} A_1 &= \left[|b_i| - \frac{|\Delta_j|}{2} \left(\frac{l_{\Sigma_{j+1}^{m+1}}}{g_{f,j+1}} + \frac{c_i}{\bar{r}_{j+1}} \right) \right] l_{\phi_{j+1,m+1}^{i,f}} \\ &- \left[|b_i| + \frac{|\Delta_j|}{2} \left(\frac{l_{\Sigma_{j+1}^{m+1}}}{g_{f,j+1}} - \frac{c_i}{\bar{r}_{j+1}} \right) \right] l_{\phi_{j,m}^{i-1,f}} \\ &+ \left[|b_i| - \frac{|\Delta_j|}{2} \left(\frac{l_{\Sigma_{j+1}^{m+1}}}{g_{f,j+1}} - \frac{c_i}{\bar{r}_{j+1}} \right) \right] l_{\phi_{j+1,m+1}^{i-1,f}} + \frac{e_i g_f}{2} \left[l_{s_{j,m}} + l_{s_{j+1,m+1}} \right] \\ &+ \left[\frac{|b_i| - |d_i|}{2} \right] \left(l_{\phi_{j,m}^{i-1,f}} + l_{\phi_{j,m+1}^{i-1,f}} - l_{\phi_{j+1,m}^{i-1,f}} - l_{\phi_{j+1,m+1}^{i-1,f}} \right) \end{aligned} \quad (20b)$$

For subcase A-2

$$l_{\phi_{j,m}^{i,f}} = \frac{|b_i| (1-\omega)^{i,f} \left(l_{\phi_{j,m+1}^{i,f}} + l_{\phi_{j,m+1}^{i-1,f}} - l_{\phi_{j,m}^{i-1,f}} \right) + \omega^{i,f} A_2}{|b_i| + \omega^{i,f} \frac{|\Delta_j|}{2} \left(\frac{l_{\Sigma_{j-1}^{m+1}}}{g_{f,j-1}} + \frac{c_i}{\bar{r}_{j-1}} \right)} \quad (21a)$$

where

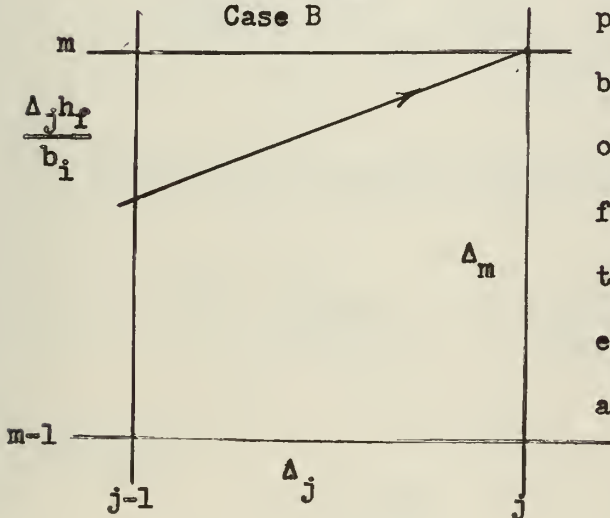
$$\begin{aligned}
 A_2 = & \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j-1}}^{l_{m+1}} + \frac{c_i}{r_{j-1}} \right) \right] \phi_{j-1, m+1}^{i, f} \\
 & - \left[|b_i| + \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j-1}}^{l_{m+1}} - \frac{c_i}{r_{j-1}} \right) \right] \phi_{j, m}^{i-1, f} \\
 & + \left[|b_i| - \frac{|\Delta_j|}{2} \left(g_{f \Sigma_{j-1}}^{l_{m+1}} - \frac{c_i}{r_{j-1}} \right) \right] \phi_{j-1, m+1}^{i-1, f} + \frac{g_{i f}}{2} |\Delta_j| \left[\psi_{j, m} + \psi_{j-1, m+1} \right] \\
 & + \left[\frac{|b_i| - |d_i|}{2} \right] \left[\phi_{j, m}^{i-1, f} + \phi_{j, m+1}^{i-1, f} - \phi_{j-1, m}^{i-1, f} - \phi_{j-1, m+1}^{i-1, f} \right] \quad (21b)
 \end{aligned}$$

A similar procedure is used in deriving the equations for the

Figure V

Position Mesh

Case B



B cases; except that as shown in Figure 5 the r and z components of the flux are different because $b_i \neq h_f$. The method of deriving the flux equation for case B-3 is outlined on the following page. The equations for the other B cases are listed also.

$$\begin{aligned}
& \left| \frac{h_f}{\Delta_m} \right| \left(l_{\phi_{j-1,m}}^{i,f} + l_{\phi_{j-1,m}}^{i-1,f} - l_{\phi_{j-1,m-1}}^{i,f} - l_{\phi_{j-1,m-1}}^{i-1,f} \right) \\
& + \left| \frac{b_i}{\Delta_j} \right| \left(l_{\phi_{j,m}}^{i,f} + l_{\phi_{j,m}}^{i-1,f} - l_{\phi_{j-1,m}}^{i,f} - l_{\phi_{j-1,m}}^{i-1,f} \right) = \\
& - \left[\frac{l_{\Sigma_{j-1}}^{m-1}}{g_f} + \frac{c_i}{\bar{r}_{j-1}} \right] \left[\frac{l_{\phi_{j,m}}^{i,f} + l_{\phi_{j-1,m-1}}^{i,f}}{2} \right] \\
& - \left[\frac{l_{\Sigma_{j-1}}^{m-1}}{g_f} - \frac{c_i}{\bar{r}_{j-1}} \right] \left[\frac{l_{\phi_{j,m}}^{i-1,f} + l_{\phi_{j-1,m-1}}^{i-1,f}}{2} \right] \\
& + \left[\frac{|b_i| - |d_i|}{2|\Delta_j|} \right] \left[l_{\phi_{j,m}}^{i-1,f} + l_{\phi_{j,m-1}}^{i-1,f} - l_{\phi_{j-1,m}}^{i-1,f} - l_{\phi_{j-1,m-1}}^{i-1,f} \right] \\
& + \frac{e_i g_f}{2} \left[l_{S_{j,m}} + l_{S_{j-1,m-1}} \right] \tag{22a}
\end{aligned}$$

To both sides of equation (22a) we add

$$\left| \frac{b_i}{\Delta_j} \right| \left(l_{\phi_{j-1,m-1}}^{i,f} + l_{\phi_{j-1,m-1}}^{i-1,f} - l_{\phi_{j,m}}^{i,f} - l_{\phi_{j,m}}^{i-1,f} \right)$$

and define

$$\omega^{if} = \frac{|\Delta_j| |h_f|}{|\Delta_m| |b_i|}$$

we obtain

$$|b_i| (1 - \omega^{i,f}) \left[l_{\phi_{j-1,m}}^{i,f} - l_{\phi_{j-1,m}}^{i-1,f} + l_{\phi_{j-1,m-1}}^{i,f} + l_{\phi_{j-1,m-1}}^{i-1,f} \right] =$$

$$\left[|b_i| + \frac{\Delta_j}{2} \left(g_f \sum_{j-1}^{m-1} + \frac{c_i}{\bar{r}_{j-1}} \right) \right] l_{\phi_{j,m}}^{i,f} + A_3. \quad (22b)$$

From this expression

$$l_{\phi_{j,m}}^{i,f} = \frac{|b_i| (1 - \omega^{i,f}) \left(l_{\phi_{j-1,m}}^{i,f} + l_{\phi_{j-1,m}}^{i-1,f} - l_{\phi_{j-1,m-1}}^{i,f} - l_{\phi_{j-1,m-1}}^{i-1,f} \right) + A_3}{|b_i| + \frac{\Delta_j}{2} \left(g_f \sum_{j-1}^{m-1} + \frac{c_i}{\bar{r}_{j-1}} \right)} \quad (22c)$$

where A_3 is defined by equation (18c)

For case B-4

$$l_{\phi_{j,m}}^{i,f} = \frac{|b_i| (1 - \omega^{i,f}) \left[l_{\phi_{j+1,m}}^{i,f} + l_{\phi_{j+1,m}}^{i-1,f} - l_{\phi_{j+1,m-1}}^{i,f} - l_{\phi_{j+1,m-1}}^{i-1,f} \right] + A_4}{|b_i| + \frac{\Delta_j}{2} \left(g_f \sum_{j+1}^{m-1} + \frac{c_i}{\bar{r}_{j+1}} \right)} \quad (23)$$

where A_4 is defined by equation (19b)

For case B-1

$$l_{\phi_{j,m}}^{i,f} = \frac{|b_i| (1 - \omega^{i,f}) \left[l_{\phi_{j+1,m}}^{i,f} + l_{\phi_{j+1,m}}^{i-1,f} - l_{\phi_{j+1,m+1}}^{i,f} - l_{\phi_{j+1,m+1}}^{i-1,f} \right] + A_1}{|b_i| + \frac{\Delta_j}{2} \left(g_f \sum_{j-1}^{m+1} + \frac{c_i}{\bar{r}_{j-1}} \right)} \quad (24)$$

where A_2 is defined by equation (20b)

For case B-2

$$l_{\phi_{j,m}^{i,f}} = \frac{|b_i| (1 - \omega^{i,f}) \left[l_{\phi_{j-1,m}^{i,f}} + l_{\phi_{j-1,m}^{i-1,f}} - l_{\phi_{j-1,m+1}^{i,f}} - l_{\phi_{j-1,m-1}^{i-1,f}} \right] + A_2}{|b_i| + \frac{|\Delta_j|}{2} \left(\epsilon_f \lambda_{\Sigma_{j-1}}^{m+1} + \frac{c_i}{r_{j-1}} \right)} \quad (25)$$

where A_2 is defined by equation (21b).

The vector fluxes defined by equations (18d), (19a), (20a), (21a), (22c), (23), (24), (25), are valid for all values of μ with the exceptions of the end points of our μ space where $\mu = \pm 1$. At these points trouble is experienced with singularities.^[11] However, a modified equation may be derived by letting the terms containing $\frac{1}{r}$ in all equations following equation (1) be replaced by zero, since for $\mu = \pm 1$ the term in $\frac{1}{r}$ in equation (1) vanishes. This special equation was not needed since values of μ were chosen which avoided these singular points.

The procedures described above result in n equations. One more equation is needed and is provided by setting $n = -1$ in equation (3). The development above may be used when the values of b , c , ϕ_0 , ϕ_{-1} , and ϵ used above are replaced by -1 , 0 , $\phi(\eta = -1)$, 0 , and 1 , respectively. These values reduce equation (6) to that found for the special case. Thus equation (6c) becomes with $i = 0$

$$l_{\phi_{j,m}^{0,f}} = \frac{(1 - \omega^{0,f}) l_{\phi_{j,m-1}^{0,f}} + \omega^{0,f} A_3}{1 + \frac{\omega^{0,f} |\Delta_j|}{2} \left(\epsilon_f \lambda_{\Sigma_{j-1}}^{m-1} \right)}$$

where

$$A_3' = \left[+1 - \frac{|\Delta_j|}{2} \binom{l_{\Sigma, m-1}}{g_f, j-1} \right] \left[l_{j-1, m-1}^{0, f} + \frac{g_f |\Delta_j|}{2} \left[l_{j, m}^- + l_{m-1, m-1}^- \right] \right].$$

Nomenclature

b_i c_i d_i e_i	}	Values obtained by differencing and integration of the transport equation over η , and defined by Equations 7a to 7d of Appendix A.
g_f h_f	}	Values depending upon the quantized values of μ , and defined by Equations 7e and 7f of Appendix A.
\hat{i}		a unit vector in the direction indicated by a subscript.
l_p		Fraction of fission neutrons which are produced in a particular energy group l .
r		The radial distance to a point.
\bar{r}_h		The mean radius between two adjacent grid points.
z		The axial distance.
E		Neutron energy from a point. See Figure 1, Appendix A.
H_f		Values obtained by the use of the Gauss quadrature approximation to sum the flux over μ space, defined by Equation 13, Appendix A.
P_i		Values obtained by the use of the S_2 approximation to sum the flux over μ space, defined by Equation 12, Appendix A.
α		Ratio of the minimum neutron energy after scattering to its energy before scattering.

- Δ Grid spacing of a mesh.
- η Cosine of the angle ψ
- θ Angle between the axis and the direction in which neutrons travel. See Figure I, Appendix A.
- μ Cosine of the angle θ .
- μ_{lk} The probability that a neutron which is scattered in the k^{th} group will land in the l^{th} group.
- ν Number of neutrons produced per fission.
- Σ Total macroscopic cross section.
- Σ_f Fission macroscopic cross section.
- ρ_{lk} The probability per unit length of neutron travel that a neutron is scattered in group k .
- $\phi(r, z, \mu, \eta, v)$ Neutron flux which is classified according to speed, v , colatitude and azimuthal direction cosines μ and η , and position r and z .
- ψ Colatitude angle denoting neutron travel with respect to the radius. See Figure I, Appendix A.
- ω See defining equation Page A-12, Appendix A.
- ω' See defining equation Page A-17, Appendix A.
- \hat{n} Direction of unit vector in which neutrons travel.

Superscripts and Subscripts

- i Index labeling quantized values of η .
- f Index labeling quantized values of μ .
- j grid position index in the radial direction.
- m grid position index in the axial direction.
- l neutron speed index.
- k neutron speed index of group in which neutron is scattered.

Appendix B

Whirlwind Program

Prior to the use of this program the following preset parameters must be defined and/or calculated:

pz1 = number of grid points in a radial direction less one

pz2 = number of grid points in an axial direction less one

pz3 = number of iterations, N , allowed at a set radius, in attempting to achieve a convergence of source pattern.

pz4 = number of adjustments, (N') , of radius allowed in attempting to achieve criticality.

pf2 \cong drum address of cross section

$$pf1 + 92(pz1 + 2) (pz1 + 2)$$

pf3 = $92(pz2 + 2)$

pf4 = $6(pz2 + 1)$

The preset parameters listed below were used in the development of the program and have the values as indicated

pf1 = 5000; drum address of first flux at origin point.

pf5 = 704; address in fast memory at which the print-out routine will be placed when needed

pf10 = 2650; address of the print-out routine in the drum

pf15 = 300; length of the print-out routine

pf13 = 1470; address in the fast memory of the vector flux subroutine of the quadrant being calculated.

It is necessary also to furnish the program at

- 1) a9, the value of $+(2pz1).0$. This value is used in the calculation of $\frac{\Delta j}{2}$

2) r_4 , the value of the first trial radius of core

3) r_3 , the value of the second trial radius of core

For a three group solution, values of each of the thirty-six fluxes and the three sources at each point are assumed and placed on tape in the order indicated in Table I.

Table I

Sequence of fluxes at a point

<u>f</u>	<u>i</u>	<u>l</u>
0	0	0
		1
		2
	1	0
		1
		2
	2	0
		1
		2
1	0	0
.	.	.
$0 \leq f < 3$	$0 \leq i < 2$	$0 \leq l < 2$

Table III

Sequence of additional data for a point

γ^{11}

γ^{21}

γ^{31}

γ^{22}

γ^{32}

γ^{33}

Σ_f

The total scattering cross section is associated with the region between points and are ordered as in Table IV.

Table IV

Sequence of Total cross sections

<u>i</u>		<u>m</u>
0		0
		1
		•
		•
		•
		pz2
2	•	
	•	
	•	
pz1	•	0
	•	
	•	
	•	pz2

fc TAPE 410-316-1032 S² AND GAUSS CAMPBELL AND PANCIERA
(24,6)

pz1=5 pz2=5 pz3=20 pz4=20 pfl=5000 pf2=10000 pf3=644
pfl4=36 pf5=704 pfl0=2650 pfl3=1470 pfl5=300

/Constants

a1,+1.0 =0.333333 +0.6666667 a2,+0.0 +1.333333 +2.0
a3,+2.46 a4,-1.0 a5,+1.9677054 +1.1306938
+1.306938 +1.9677054 a6,+0.03 +0.03 +10.0 a7,+0.3183
+0.3634 +0.3183 a8,+0.6521452 +0.3478548 +0.3478548
+0.6521452 a9,+22.0 a10,+4.0 a11,+1.6944625 +0.3843341

/Cross Sections

b1,+1.0 DITTO 12r/b2,+1.0 DITTO 12r/m6,+0 m7,+0
m8,pf3 m9,pf4 m10,+92 m11,+6 m12,pf1 m13,pf2 m14,+0
m15,+0

/RADII

r3,+0.0 r4,+0.0 r5,+0.0

/NORMALIZING CONSTANTS

d1,+0.0 d2,+0.0 d3,+0.0 r1,+0.0 r2,+0.0 s10,+0.0

/VARIABLES

c1,+4.8 c2,=5.6 c3,+0.6885 +0.2295 +0.4590 +3.0355
+1.0118 +2.0236

/j=1,m-1 numbers

n1,+1.0 DITTO 72r/s1,+1.0 +1.0 +1.0 t1,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j=1,m numbers

n2,+1.0 DITTO 72r/s2,+1.0 +1.0 +1.0 t2,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j=1,m+1, numbers

n3,+1.0 DITTO 72r/s3,+1.0 +1.0 +1.0 t3,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j, m=1, numbers

n4,+1.0 DITTO 72r/s4,+1.0 +1.0 +1.0 t4,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j,m numbers

n5,+1.0 DITTO 72r/s5,+1.0 +1.0 +1.0 t5,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j,m+1, numbers

n6,+1.0 DITTO 72r/s6,+1.0 +1.0 +1.0 t6,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j+1,m-1, numbers
n7,+1.0 DITTO 72r/s7,+1.0 +1.0 +1.0 t7,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j+1,m, numbers
n8,+1.0 DITTO 72r/s8,+1.0 +1.0 +1.0 t8,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/j+1,m+1, numbers
n9,+1.0 DITTO 72r/s9,+1.0 +1.0 +1.0 t9,+1.0
+1.0 +1.0 +1.0 +1.0 +1.0 +1.0

/CONVERGED GAMMA

m1, itam18 itr5 idva9 itscl isc0 icr3 iscl
icr6 isc2 icr2 m2,icac2 idvcl isc2 idvall+c
isc0 imral+c icl1 itsc3+c isc0 ictm2 isc2
ictm2 isc0 icrpz3 m3,icaal itssl0 icac1 itsrl
iscl icrpz1+1 isc2 icrpz2+1 m4,ispil isc2 ictm4
icar1 itr2 iadcl iadcl itsrl iscl ictm4
isc0 ictm5 iSTOP m5,icas10 icpm3 m18,ispo

/j,m SENSE AND CONTROL

il,itai7 iscl itim6 isc2 itim7 OUT cam7
mhm10 slhl5 adm12 tsm15 cam8 mhm6 slhl5
tsm14 adm15 tsi6 cam15 sum10 tsm15 adm14
tsm29 cam14 sum8 adnk5 tsm26 cam14 adm8
adm15 tsm30 cam7 sul mhm11 slhl5 adm13
tsm15 cam9 mhm6 slhl5 tsm14 adm15 tsm32
cam14 sum9 adm15 tsm31 cam6 dm0 cpi2
cam7 dm0 cpi3 IN ispm23 ispm27 ispi5
i2,cam7 dm0 cpi4 IN ispm23 ispm21 ispi5
i3,IN ispm23 ispm24 ispi5 i4,IN ispm23 ispm16
i5,iDOB n5 i6,+0 +78 i7,ispo

/DATA READ IN PROGRAMME

m23,itam33 iDIB n1 m26,+0 +276 iDIB n4
m29,+0 +276 iDIB n7 m30,+0 +276 iDIB
b1 m31,+0 +6 iDIB b2 m32,+0 +6

m33,ispo

/ORIGIN PROGRAM

m16,itam17 ispp10 ispp13 icaal idvt7 itsdl ispe3
iDIB pfl3 +2050 +175 ispf1 m17,isp0

/CENTERLINE PROGRAM

m21,itam22 ispp4 ispp13 ispe3 iDIB pfl3 +2050
 +175 ispf1 iDIB pfl3 +2475 +175 ispf18
 m22,isp0

/MIDPLANE PROGRAM

m24,itam25 ispp7 ispp13 ispe3 iDIB pfl3 +2050
 +175 ispf1 iDIB pfl3 +2225 +125 ispf10
 m25,isp0

/GENERAL POINT PROGRAM

m27,itam28 ispp1 ispp13 ispe3 iDIB pfl3 +2050
 +175 ispf1 iDIB pfl3 +2225 +125 ispf10
 iDIB pfl3 +2350 +125 ispf27 iDIB pfl3
 +2475 +175 ispf18 m28,isp0

/Source S^0, S^1, S^2 at all Points

pl3, itapl4 icat6 imrt5 itst1 icaa3 imrt5+12 imrt6+4
 iadtl itst7 icat6 imrt5+2 itst1 icat6+2 imrt5+6
 iadtl itst7+2 icat6 imrt5+4 itst1 icat6+2 imrt5+8
 iadtl itst1 icat6+4 imrt5+10 iadtl itst7+4 pl4,isp0

/Source Normalization+Tolerance Program

e3, itae7 icat7+2 imrd1 itss5+2 icat7+4 imrd1 itss5+4
 icat7 imrd1 itst7 isus5 icpe4 idvt7 isua6
 icpe5 e6, icsa6 itss10 ispe5 e4, idvt7 iada6 icpe6
 e5, icat7 itss5 e7, isp0

/SCALAR Flux-General Point

pl, itap3 isc3 icr3 p2, icaa7 imrn5+c itst1 icaa7+2
 imrn5+6+c iact1 itst1 icaa7+4 imrn5+12+c iadtl
 imra8 itst1 icaa7 imrn5+18+c itst2 icaa7+2
 imrn5+24+c iadt2 itst2 icaa7+4 imrn5+30+c iadt2
 imra8+2 iadtl itst1 icaa7 imrn5+36+c itst2
 icaa7+2 imrn5+42+c iadt2 itst2 icaa7+4 imrn5+48+c
 iadt2 imra8+4 iadtl itst1 icaa7 imrn5+54+c
 itst2 icaa7+2 imrn5+60+c iact2 itst2 icaa7+4
 imrn5+66+c iadt2 imra6+6 iadtl itst6+c ictp2 p3, isp0

/Scalar Flux-Centerline Point

pl, itap6 isc3 icr3 p5, icaa8 imrn5+c itst1 icaa8+2
 imrn5+18+c iadtl itst1 icaa8+4 imrn5+36+c iadtl
 itst1 icaa8+6 imrn5+54+c iadtl imra7 itst1
 icaa8 imrn5+6+c itst2 icaa8+2 imrn5+24+c iadt2

itst2 icaa8+4 imrn5+42+c iadt2 itst2 icaa8+6
imrn5+60+c iadt2 imra7+2 iadt1 imra2+4 itst6+c ictp5
p6,isp0

/Scalar Flux-Midplane Point

p7,itap9 isc3 icr3 p8,icaa7 imrn5+c itst1 icaa7+2
imrn5+6+c iadt1 itst1 icaa7+4 imrn5+12+c iadt1
imra8 itst1 icaa7 imrn5+18+c itst2 icaa7+2
imrn5+24+c iadt2 itst2 icaa7+4 imrn5+30+c iadt2
imra8+2 iadt1 imra2+4 itst6+c ictp8 p9,isp0

Scalar Flux - Origin

pl0,itapl2 isc3 icr3 pl1,icaa7 imrn5+c itst1 icaa7+2
imrn5+6+c iadt1 imra8 itst1 icaa7 imrn5+18+c
itst2 icaa7+2 imrn5+24+c iadt2 imra8+2 iadt1
imra2+4 imra2+4 itst6+c ictpl1 pl2,isp0

/RADIUS EXTRAPOLATION ROUTINE

qi,itaq9 icad2 isual itst1 icpq2 q3,itst3 icad3
isual itst2 icpq4 q5,itst4 isut3 icpq6 icad2
iexd3 itsd2 icar3 iexr4 itsr3 icat3 iext4
itst3 icat1 iext2 itst1 q6,icat3 isua6+2 icpq7
q8,icad2 isud3 itst6 icar4 isur3 imrt1 idvt6
iadr3 itsr3 q9,isp0 q2,imra4 ispq3 q4,imra4 ispq5

/Control

x1,icar4 isch4 icrpz4 ispm1 icad1 itsd3 x2,icar3
ispm1 icad1 itsd2 isch4 ictx3 istop x3,ispql
ispx2 q7,ldib pf5 pf10 pf15 ql0,ispff5
istop

/WORKING FLUX

1470/y1,+0 DITTO 175r/y2,+0

/First Quadrant Flux

/First Flux

DA2050/ 1470/ f1,itaf3 isc3 icr3 f2,icsb2+6+c
imra5 imrc1 itst1 iadal imrn9+c itst2 icas5+c
iads9+c imra5 imrc1 iadt2 imrc3 itst2 icst1
imrc3 iadal itst1 icaal isuc3 imrn6+c iadt2
idvt1 itsn5+c

/Second Flux

f4,icsa2+2 idvr1 itst1 icsb2+6+c imra5 itst2 iadt1

imrc1	itst3	iadal+2	imrn9+6+c	itst4	icst1	iadt2
imrc1	itst6	iadal+2	imrn9+c	iadt4	itst4	icat6
isual+2	f5,imrn5+c		iadt4	itst4	icas5+c	iads9+c
imra5	imrc1	imra2+4	iadt4	itst4	icsal+4	iadal+2
idva2+4	itst1	ican6+c	iadn5+c	isun9+c	isun8+c	imrt1
iadt4	imrc3+2	itst4	f6,icst3	imrc3+2	iadal+2	itst2
icaal	isuc3+2	imra1+2	itst3	ican6+6+c	iadn6+c	isun5+c
imrt3	iadt4	idvt2	itsn5+6+c			

/Third Flux

f7,icsb2+6+c		imra5+2	imrc1	itst1	iadal	imrn9+18+c
itst2	icas5+c	iads9+c	imra5+2	imrc1	iadt2	imrc3+6
itst2	icst1	iadal	imrc3+6	itst1	icac3+6	isual
itst3	ican8+18+c		isun9+18+c		imrt3	iadt2
idvt1	itsn5+18+c					

/Fourth Flux

f8,icsa2+2	idvr1	itst1	icsb2+6+c	imra5+2	itst2	iadt1
imrc1	itst3	iadal+2	imrn9+24+c		itst4	icst1
iadt2	imrc1	itst6	iadal+2	imrn9+18+c		iadt4
itst4	icat6	isual+2	imrn5+18+c		iadt4	itst4
f9,icas5+c	iads9+c	imra5+2	imrc1	imra2+4	iadt4	itst4
icsal+4	iadal+2	idva2+4	itst1	ican6+18+c		iadn5+18+c
isun9+18+c	isun8+18+c		imrt1	iadt4	imrc3+8	itst4
icst3	imrc3+8	iadal+2	itst2	icaal	isuc3+8	imra1+2
itst3	ican6+24+c		iadn6+18+c		isun5+18+c	
imrt3	iadt4	idvt2	itsn5+24+c		ictf2	f3,isp0

/Second quadrant flux

/First flux

DA2225/ 1470/ f10, itaf17			isc3	icr3	f11,icsa2+2	
idvr2	itst1	icsb1+6+c	imra5	itst2	iadt1	imrc1
itst3	iadal+4	imrn3+12+c		itst4	icst1	iadt2
imrc1	itst6	iadal+4	imrn3+6+c	iadt4	itst4	icat6
isual+4	imrn5+6+c	f12, iadt4	itst4	icas5+c	iads3+c	imra5
imrc1	imra2+4	iadt4	itst4	icaal+4	isual+2	idva2+4
itst1	ican6+6+c	iadn5+6+c	isun3+6+c	isun3+6+c	itst3	ican6+12+c
iadn6+6+c	isun5+6+c	imrt3	iadt4	idvt2	itsn5+12+c	

/Second Flux

f14,icsa2+2	idvr2	itst1	icsb1+6+c	imra5+2	itst2	iadt1
imrc1	itst3	iadal+4	imrn3+30+c		itst4	icst1
iadt2	imrc1	itst6	iadal+4	imrn3+24+c		iadt4
itst4	icat6	isual+4	imrn5+24+c		f15, iadt4	itst4

icas5+c	iads3+c	imra5+2	imrc1	imra2+4	iadt4	itst4
icaal+4	isual+2	idva2+4	itst1	ican6+24+c		iadn5+24+c
isun3+24+c	isun2+24+c		imrt1	iadt4	imrc3+10	itst1
icst3	iadal+4	imrc3+10	itst2	f16, icac3+10		isual
imral+4	itst3	ican2+30+c		iadn2+24+c		isun3+30+c
isun3+24+c	imrt3	iadt1	idvt2	itsn5+30+c		ictf11

f17,isp0

/Third Quadrant Flux

/First Flux

DA2350/ 1470/ f27, itaf33		isc3	icr3	f28, icsa2+2		
idvr2	itst1	icsb1+c	imra5+4	itst2	iadt1	imrc1
itst3	iadal+4	imrn1+48+c		itst4	icst1	iadt2
imrc1	itst1	iadal+4	imrn1+42+c		iadt4	itst4
icat1	isual+4	imrn5+42+c		iadt4	itst4	f29, icas5+c
iadsl+c	imra5+4	imrc1	imra2+4	iadt4	itst4	icaal+4
isual+2	idva2+4	itst1	ican5+42+c		iadn4+42+c	
isun2+42+c	isun1+42+c		imrt1	iadt4	imrc3+10	itst1
icst2	iadal+4	imrc3+10	itst2	icac3+10	isual	imral+4
itst3	ican2+48+c		iadn2+42+c		isun1+48+c	
isun1+42+c	imrt3	iadt1	idvt2	f30, itsn5+48+c		

/Second Flux

f31, icsa2+2	idvr2	itst1	icsb1+c	imra5+6	itst2	iadt1
imrc1	itst3	iadal+4	imrn1+66+c		itst4	icst1
iadt2	imrc1	itst1	iadal+4	imrn1+60+c		iadt4
itst4	icat1	isual+4	imrn5+60+c		iadt4	itst4
f32, icas5+c	iadsl+c	imra5+6	imrc1	imra2+4	iadt4	itst4
icaal+4	isual+2	idva2+4	itst2	ican5+60+c		iadn4+60+c
isun2+60+c	isun1+60+c		imrt2	iadt4	imrc3+4	itst1
icst3	imrc3+4	iadal+4	itst2	icaal	isuc3+4	imral+4
itst3	ican4+66+c		iadn4+60+c		isun5+60+c	
imrt3	iadt1	idvt2	itsn5+66+c		ictf28	f33, isp0

/Fourth Quadrant Flux

/First Flux

DA2475/ 1470/ f18, itaf26		isc3	icr3	f19, icsb2+c		
imra5+4	imrc1	itst1	iadal	imrn7+36+c		itst2
icas5+c	iads7+c	imra5+4	imrc1	iadt2	imrc3+6	itst2
icst1	iadal	imrc3+6	itst1	icac3+6	isual	itst3
ican8+36+c	isun7+36+c		imrt3	iadt2	idvt1	itsn5+36+c

/Second Flux

f2o,icsa2+2	idvr1	itst1	icsb2+c	imra5+4	itst2	iadt1
imrc1	itst3	iadal+2	imrn7+42+c		itst4	icst1
iadt2	imrc1	itst6	iadal+2	imrn7+36+c		iadt4
itst4	icat6	isual+2	imrn5+36+c		iadt4	itst4
f21,icas5+c	iads7+c	imra5+4	imrc1	imra2+4	iadt4	itst4
icaal+4	idva2+4	itst1	ican5+36+c		iadn4+36+c	
isun8+36+c	isun7+36+c		imrt1	iadt4	imrc3+8	itst1
icst3	imrc3+8	iadal+2	itst2	f22,icaal	isun3+8	imral+2
itst3	ican4+42+c		iadn4+36+c		isun5+36+c	
imrt3	iadt1	idvt2	itsn5+42+c			

/Third Flux

f23,icsb2+c	imra5+6	imrc1	itst1	iadal	imrn7+54+c	
itst2	icas5+c	iads7+c	imra5+6	imrc1	iadt2	imrc3
itst2	icst1	imrc3	iadal	itst1	icaal	isuc3
imrn4+54+c	iadt2	idvt1	itsn5+54+c			

/Fourth Flux

f24,icsa2+2	idvr1	itst1	icsb2+c	imra5+6	itst2	iadt1
imrc1	itst3	iadal+2	imrn7+60+c		itst4	icst1
iadt2	imrc1	itst1	iadal+2	imrn7+54+c		iadt4
itst4	icat1	isual+2	imrn5+54+c		iadt4	itst4
f25,icas5+c	iads7+c	imra5+6	imrc1	imra2+4	iadt4	icaal+2
isual+4	idva2+4	itst1	ican5+54+c		iadn4+54+c	
isun8+54+c	isun7+54+c		imrt1	iadt4	imrc3+2	itst1
icst3	imrc3+2	iadal+2	itst2	icaal	isuc3+2	imral+2
itst3	ican4+60+c		iadn4+54+c		isun5+54+c	
imrt3	iadt1	idvt2	itsn5+60+c		ictf19	f26,isp0
DA2650/704/	z1,itazl4		isc1	i crpz1+2	isc2	icrpz2+2
icad1	ispl36z13	isc1	z12,itim6	isc2	z11,itim7	OUT
cam7	mhm10	slh15	adml2	tsml5	cam8	mhm6
slh15	adml5	tsz2	cam6	dm0	cpz3	cam7
dm0	cpz4	IN	ispz5	isppl	ispz6	z3,cam7
dm0	cpz7	IN	ispz5	isppl4	ispz6	z4,IN
ispz5	isppl7	ispz6	z7,IN	ispz5	isppl0	ispz6
isc2	ictz11	isc1	ictz12	z14,isp0	z5,itaz8	IDIB
n5	z2,+0	+72	z8,isp0	z6,ispz6+1		z14,ispo
z10,icat6	ispl39z13	icat6+2	ispl39z13	icat6+4	ispl36z13	z9,ispz7+4
z13.						

Appendix C

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