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TI-652-260-001

# COVER SHEET

DOCUMENT TITLE

5 Kwe Reactor Power Distributions

AUTHOR

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ABSTRACT

Power distributions representative of beginning and end of life are presented for the 5 Kwe reactor. These distributions were obtained from DOT calculations in R, Z geometry. Reactivity worth of the sliding control segments was also obtained.

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I. Introduction

The reflector design of the 5 Kwe Reactor represents a significant departure from previously operated SNAP reflector designs. Previous designs used rotating drums to regulate neutron leakage and thereby control the reactivity of the reactor. Instead of rotating drums, the 5 Kwe Reactor uses sliding segments to regulate neutron leakage. The reflector in previous designs was also uniform in thickness over the height of the core. The 5 Kwe Reactor reflector is tapered.

The purpose of the change in reflector design is to reduce the weight of the shadow shield through a reduction in shieldable envelope. Sliding segments rather than rotating drums permit a reduction in envelope diameter approximately equal to twice the reflector thickness. Tapering the reflector permits an additional reduction in shield diameter because there is no beryllium outside the shadow cone from which radiation can scatter.

The current 5 Kwe Reactor design has two sliding segments each of which subtend  $120^{\circ}$  of the reactor diameter. The beryllium between each segment is stationary. The sliding segments are split two inches above the axial centerline. The thinner top pieces remain stationary and the thicker bottom pieces slide a maximum of four inches. Rather than sliding parallel to the core vessel the segments move along the taper angle.

The effect the sliding segments have on the power distributions could not be very well determined using the customary one-dimensional technique. A two-dimensional model of the reactor was set up in R, Z geometry for calculations using the DOT code. Calculations were made with the segments closed, two inches open, and four inches open.

In R, Z geometry there is no azimuthal variation; therefore, it was necessary to make some approximations concerning the fixed reflector when the sliding segments were not fully closed. In one calculation with the segments two inches open, the fixed beryllium was homogenized in the gap. From this a power distribution was obtained for use in the preliminary thermal, hydraulic analysis. This distribution did not contain the desired azimuthal variation. To obtain it, a power distribution was synthesized from two DOT calculations. A calculation with the segments fully closed and a calculation with a  $360^\circ$  two inch opening were used. The power distributions from the two calculations were combined using a sinusoidal azimuthal variation. A timeshare computer program was written for this purpose.

An additional calculation was made with an untapered reflector to determine the reactivity effect of tapering. A calculation with the segments four inches open was used to determine the total control worth.

## II. Summary of Results

The synthesized power distribution at beginning-of-life is shown in Figure 1- Figure 5. Figure 1 shows for one quadrant the peak power per element relative to the core average power. Due to symmetry the power distribution in the remaining quadrants is the same. The axial power distribution for each element in the quadrant is shown in Figure 2- Figure 5.

Power contours are shown in Figure 6 and Figure 7. Figure 6 shows the power contours with the segments fully closed which would be representative of end-of-life conditions. Figure 7 shows the power contours with the segments two inches open with homogenized beryllium in the opening.

The calculated reactivity change between segments fully closed and  $360^\circ$  segments four inches open was \$12.63. It is assumed that the worth is directly proportional to the angle subtended by the movable segments. The total control worth would therefore be \$8.42. The reactivity lost between a straight reflector and one with an  $8^\circ$  taper angle was 24¢.

## III. Method of Calculation

The calculations were made using the two-dimensional transport theory code, DOT II. The  $S_4$  approximation and the standard 28 group SNAP library was used. The two-dimensional model which was set up to describe the reactor is shown in Figure 8. The composition

of the various zones is given in Table I. Macroscopic cross sections are available in the SNAP library for most materials, therefore, volume fractions may be used instead of atom densities. The only exceptions were fuel and tungsten. The atom densities of these materials without volume weights is given in Table II.



Figure 1  
 Peak Power Per Element  
 Relative to Core Average  
 5 Kwe Reactor at BOL

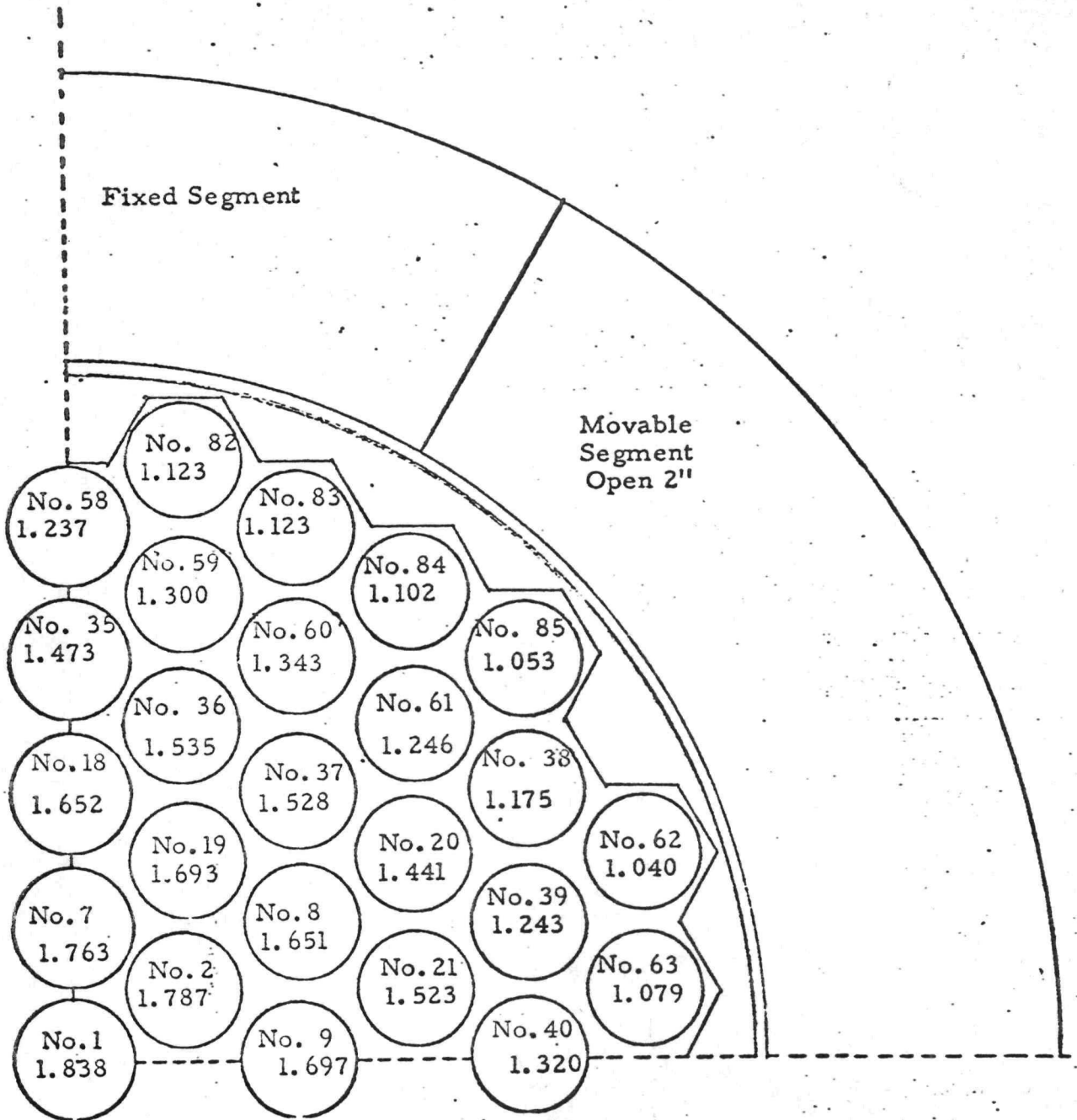
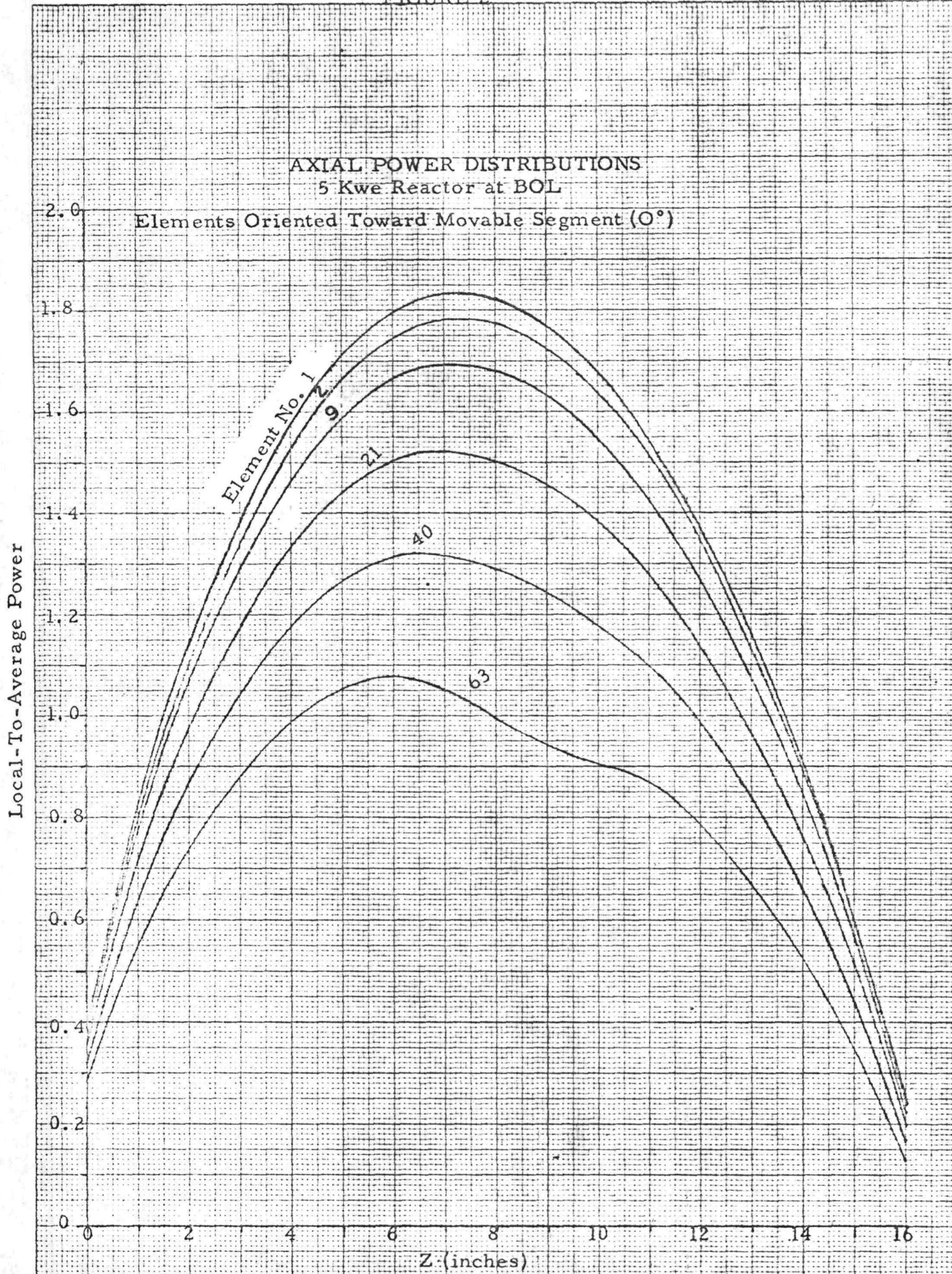


FIGURE 2



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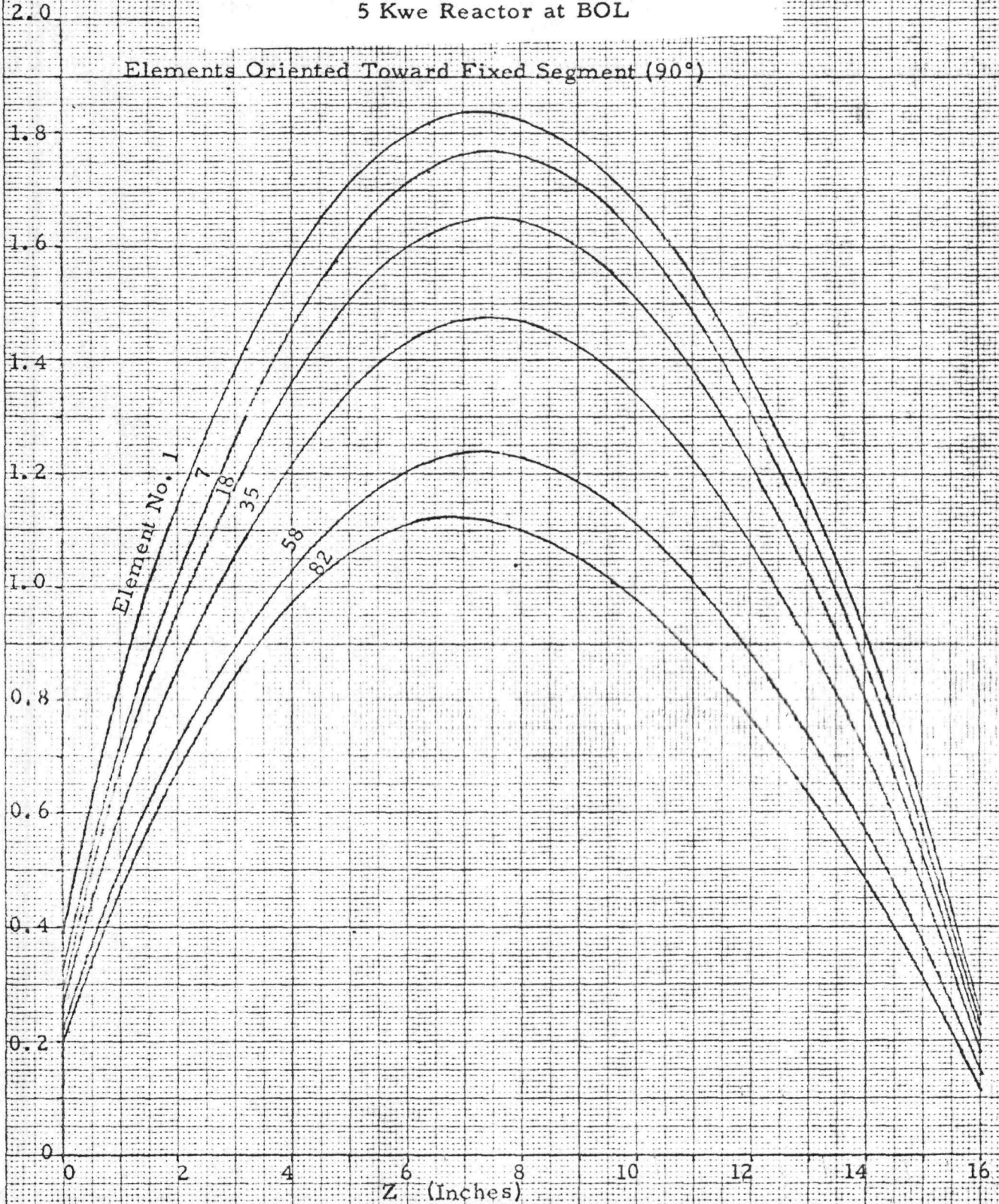
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### AXIAL POWER DISTRIBUTIONS

5 Kwe Reactor at BOL

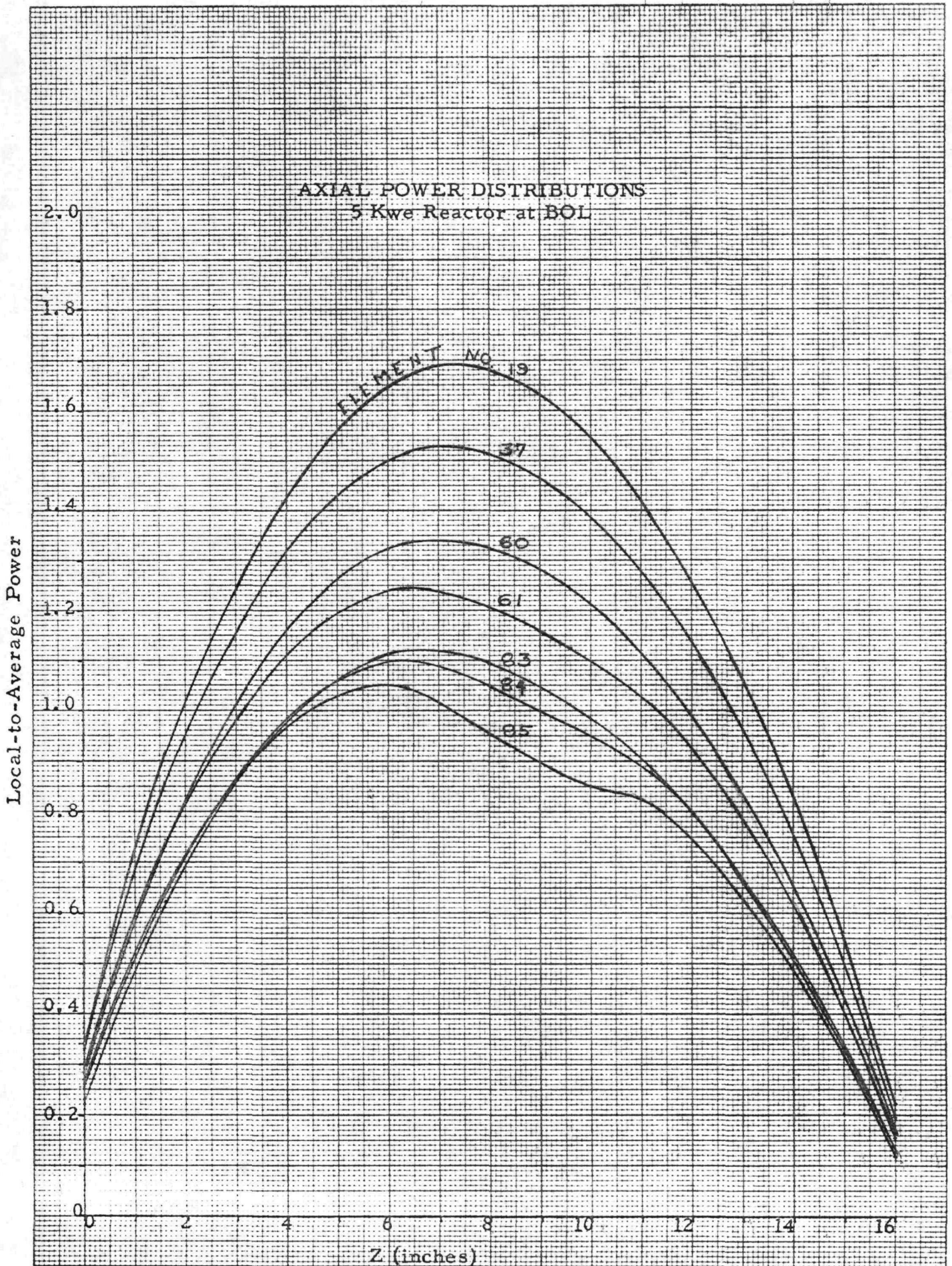
Elements Oriented Toward Fixed Segment (90°)

Local-To-Average Power



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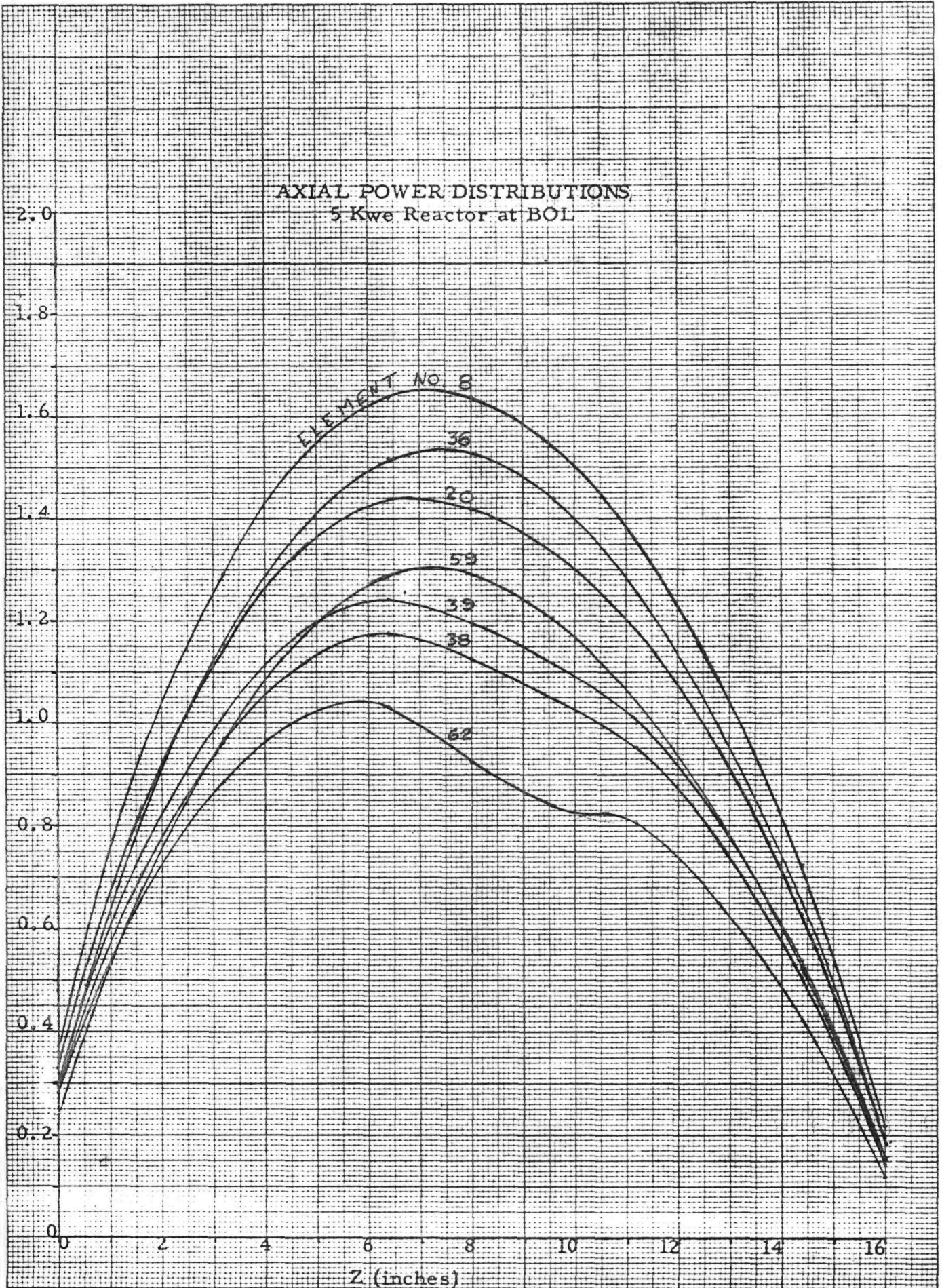


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5 Kwe Reactor at BOL

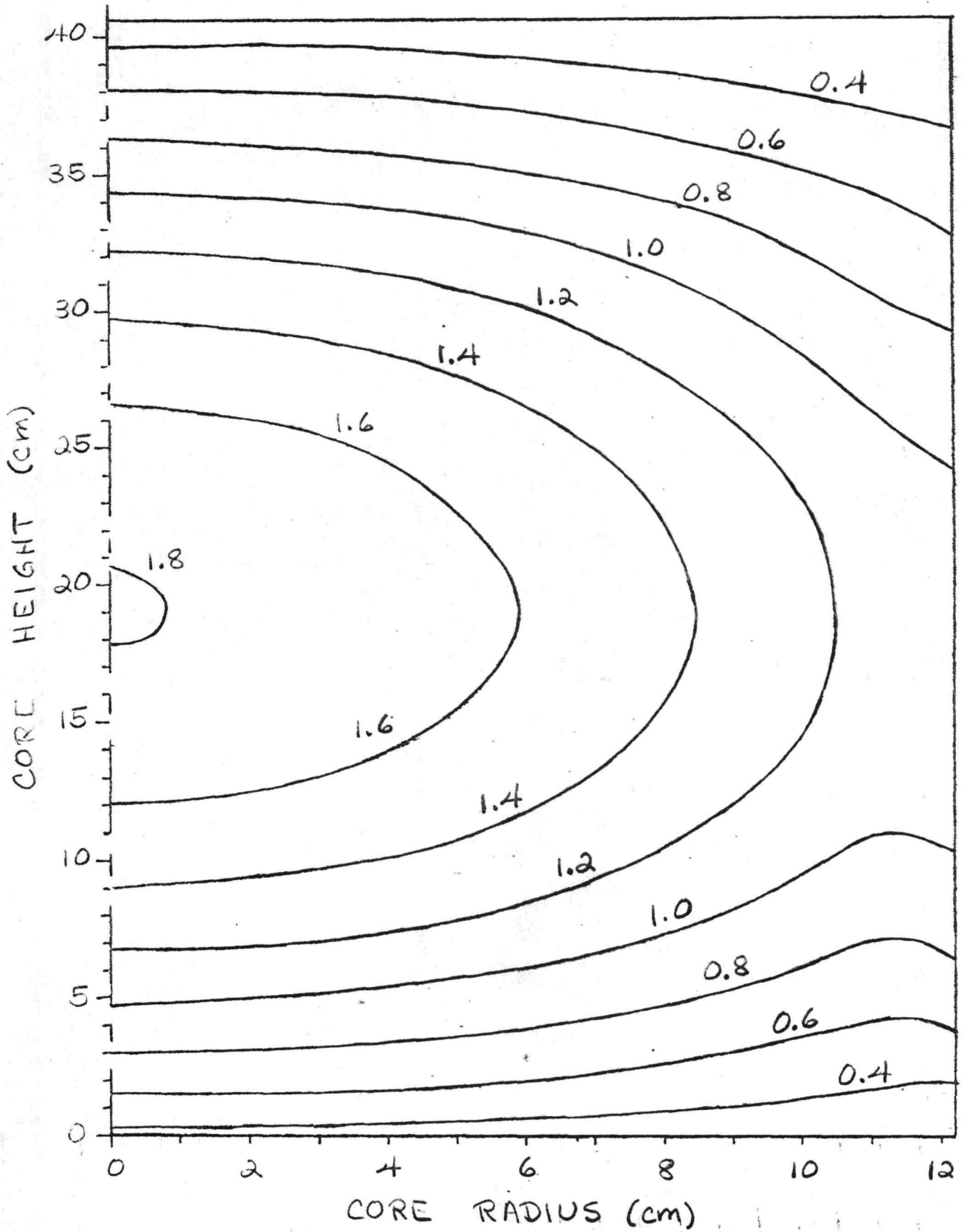
Local-to-Average Power



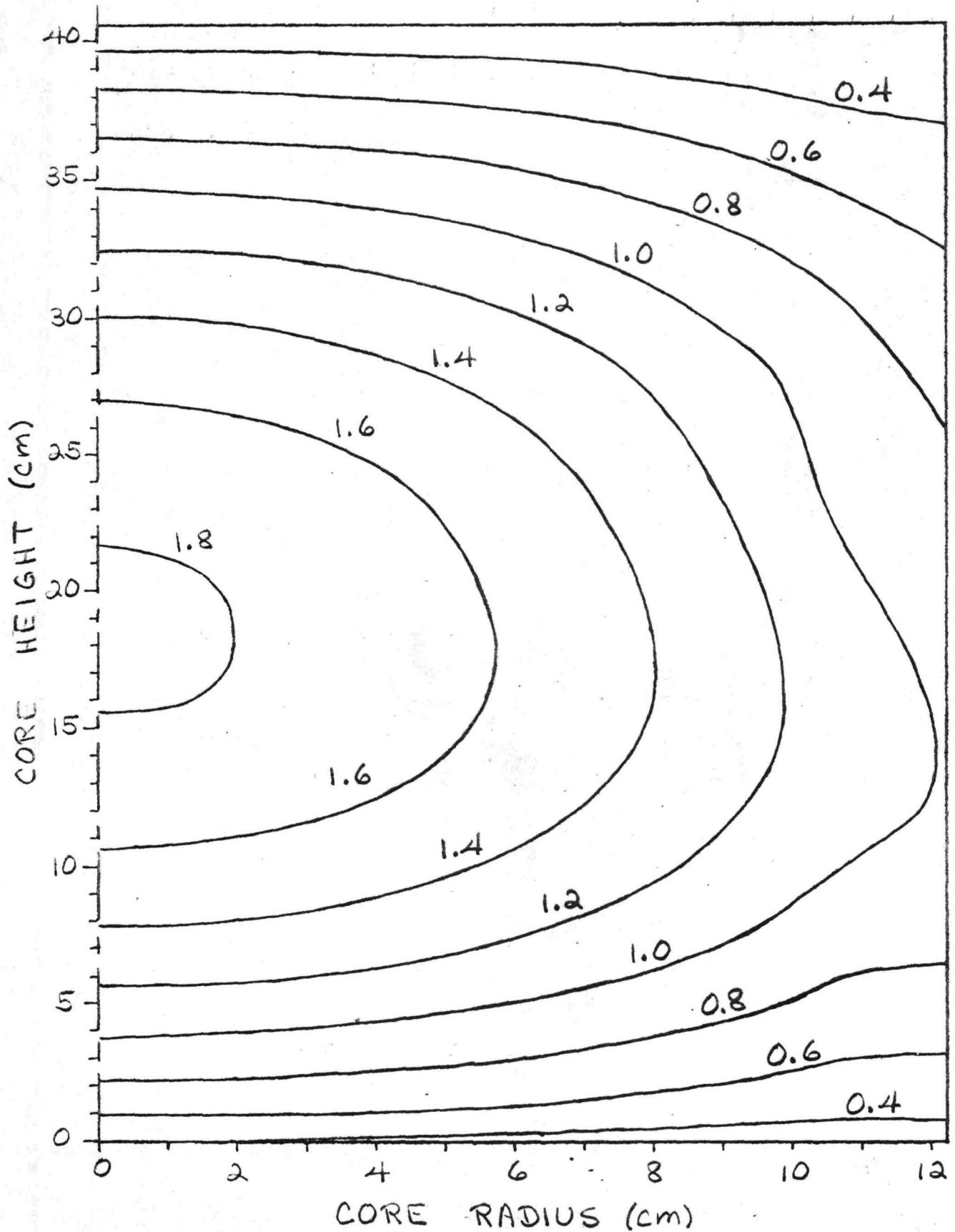
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FIGURE 6  
LOCAL-TO-AVERAGE POWER  
- SEGMENTS CLOSED -

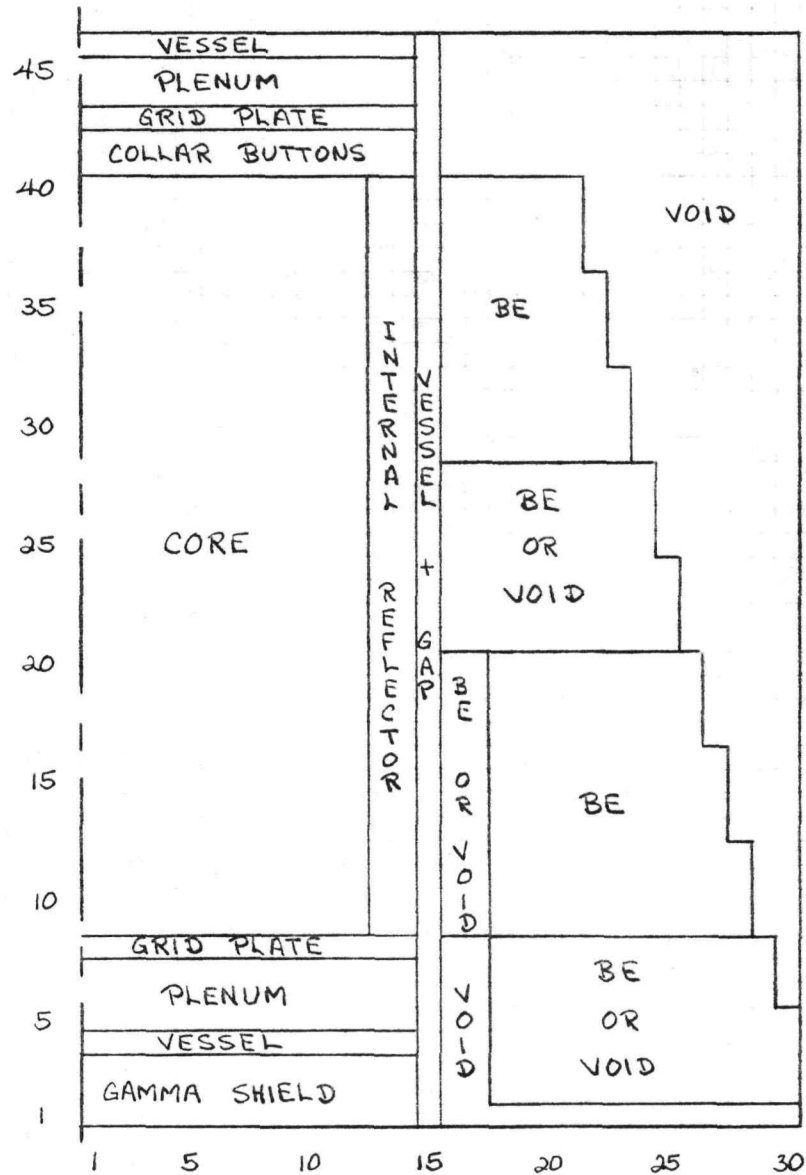


LOCAL-TO-AVERAGE POWER  
- SEGMENTS 2" OPEN - HOMOGENIZED BERYLLIUM -



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CHECKED BY:		PAGE NO. OF
DATE:	FIGURE 8	Page 13
		REPORT NO.
		MODEL NO.

### DOT R,Z MODEL



<u>INTERVALS</u>	<u>RADIUS</u>	<u>INTERVALS</u>	<u>HEIGHT</u>	<u>INTERVALS</u>	<u>HEIGHT</u>
1 - 12	12.236 cm	1	1.588 cm	8	11.748 cm
13 - 14	12.879	2	4.128	9 - 40	52.388
15	13.139	3	5.08	41 - 42	54.928
16 - 17	14.567	4	5.398	43	56.198
18 - 21	16.990	5	6.668	44 - 45	61.278
22 - 30	23.416	6	9.208	46	61.595
		7	10.488		



Table I  
DOT Model Volume Fractions

<u>Zone</u>	<u>Material</u>	<u>Volume Fraction</u>
Core	Fuel	0.7546
	Glass	0.0068
	Clad + Fins	0.0919
	NaK	0.1243
	Gap	0.0224
Internal	BeO	0.6
Reflector	316 SS	0.4
Vessel + Gap	316 SS	0.39
	Void	0.61
Grid Plate	NaK	0.32
	316 SS	0.57
	Hast X-280	0.11
Collar Buttons	NaK	0.1243
	Hast X-280	0.1
	Void	0.7757
Plenum	NaK	1.0
Vessel	316 SS	1.0
Gamma Shield	Tungsten	1.0
Reflector	Be	1.0

Table II  
Atom Densities

<u>Isotope</u>	<u>Atom Density (atoms/barn-cm)</u>
U-235	0.001503
U238	0.000109
Zr	0.03534
H	0.05725
W	0.0619