

**BORON CARBIDE ABSORBER RODS IN URANIUM
(10% ^{235}U) NITRATE SOLUTION**

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IDENTIFICATION NUMBER: LEU-SOL-THERM-006

SPECTRA

KEY WORDS: absorber, absorber rods, acceptable, boron, boron carbide, critical experiment, cylinder, homogeneous, low-enriched, moderated, nitric acid, solution, thermal, uranium, uranyl nitrate, water-reflected

1.0 DETAILED DESCRIPTION

1.1 Overview of Experiment

A large number of critical experiments with absorber elements of different types in uranium nitrate solution of different enrichments and concentrations were performed in 1961 - 1963 at the Solution Physical Facility of the Institute of Physics and Power Engineering (IPPE), Obninsk, Russia. The purpose of these experiments was to determine the effects of enrichment, concentration, geometry, neutron reflection, and type, diameter, number, and arrangement of absorber rods on the critical mass of light-water-moderated homogeneous uranyl nitrate solutions. The experiments included ones with a central boron carbide or cadmium rod, clusters of boron carbide rods, and triangular lattices of boron carbide rods in cylindrical tanks of different dimensions filled with solutions of uranyl nitrate.^a

The five experiments included in this evaluation were performed with uranium enriched to 10 wt.% ^{235}U . Uranium nitrate solution with uranium concentration of 420.5 g/l was pumped into the core or inner tank, a stainless steel cylindrical tank with inner diameter 110 cm. One experiment was performed without absorber rods. In each of four experiments a different number of boron carbide absorber rods was inserted in the core tank. The absorber rods were arranged in a hexagonal lattice with different pitches. There was a thick side and bottom water reflector in these experiments. All five configurations are considered to be acceptable for use as criticality safety benchmark experiments.

Experiments performed using the same tank and absorber rods with uranium enrichments of 5.64% and 89% are reported in evaluations LEU-SOL-THERM-005 and HEU-SOL-THERM-035.

1.2 Description of Experimental Configuration

A diagram of the critical assembly is shown in Figure 1 (vertical cut). The experiment was performed in the same room as the other uranium-solution experiments at IPPE (see Appendix B of

^a See LEU-SOL-THERM-005, HEU-SOL-THERM-027, HEU-SOL-THERM-028, HEU-SOL-THERM-029, HEU-SOL-THERM-030, HEU-SOL-THERM-031, HEU-SOL-THERM-035, and HEU-SOL-THERM-037.

HEU-SOL-THERM-014). The room had dimensions $7.5 \times 5.5 \times 8.8$ m. The core central axis was 2 m from the north wall, 5.5 m from the south wall, and 2.75 m from the west and east walls. All the walls, the ceiling, and the floor were concrete. The thickness of the walls was approximately 100 cm, the thickness of the ceiling was 75 cm, and the thickness of the floor was 20 cm.

The critical assembly consisted of two open-topped coaxial cylindrical tanks. The core tank was 110.0 cm in inner diameter, 250.0 cm tall, with wall thickness 0.6 cm, and bottom thickness 1.5 cm. The reflector tank had 198.4-cm inner diameter, was 300 cm tall, with wall thickness 0.8 cm, and bottom thickness 1.0 cm. The core tank stood on a pedestal inside the reflector tank. The height of the pedestal was 36.0 cm. No other characteristics of the pedestal are known. The reflector tank stood on the floor.

The core tank was partially filled with the aqueous solution of uranyl nitrate $\text{UO}_2(\text{NO}_3)_2$ with some excess of nitric acid (HNO_3). The reflector tank was filled with distilled water in all cases. The height of the water reflector measured from the bottom (inner surface) of the core tank was 108.0 cm.

One experiment was performed without absorber rods. The arrangement of the absorber rods for the different cases is shown in Figures 2 - 5. The experimental results obtained for the different experimental conditions are presented in Table 1.

Table 1. Critical Dimensions (see Figure 1).

Case Number	Number of Absorber Rods	Number of Holes in the Lattice Plate	Solution Volume, liters
1	none	85	210.0 ± 0.3
2	7	85	231.5 ± 0.4
3	18	163	251.3 ± 0.4
4	19	163	255.5 ± 0.4
5	31	85	306.3 ± 0.5

The average inside diameter of the core tank was measured by filling the tank with a known volume of water and then measuring the water level. Several volumes were used to establish the inside diameter of the tank. The uncertainty of the core-tank average inside diameter is ± 0.2 cm.

The core tank was filled with solution through a polyethylene feed pipe, with inner diameter 1.2 cm. The pipe wall thickness was 0.2 cm. The end of this pipe was positioned under the solution surface. Pumping the solution out of the tank was performed through a stainless steel pipe, with outer diameter 1.2 cm and wall thickness 0.15 cm. The pipes entered the tank from the top. These pipes are not shown in Figure 1.

Technical drawing of a core tank assembly, showing dimensions and components. The drawing is a cross-section view of a cylindrical tank with a total height of 300.0 cm and a total diameter of D203.4 cm. The tank is divided into three main sections: a top section (height 1.7 cm), a middle section (height 250.0 cm), and a bottom section (height 1.5 cm). The middle section contains three vertical absorber rods, each with a diameter of D5.5 cm, spaced at a center-to-center distance of D110.0 cm. The rods are supported by an upper lattice plate at the top and a lower lattice plate at the bottom. The bottom section is filled with a yellow material, likely a reflector or moderator. The tank is surrounded by a reflector tank, which has a thickness of 0.6 cm. The total diameter of the reflector tank is D200.0 cm. The drawing includes labels for the following components: Upper lattice plate, Absorber rods, Core tank, Reflector tank, and Lower lattice plate. Dimensions are given in cm.

Dimensions in cm

Revision: 0
Date: September 30, 1998

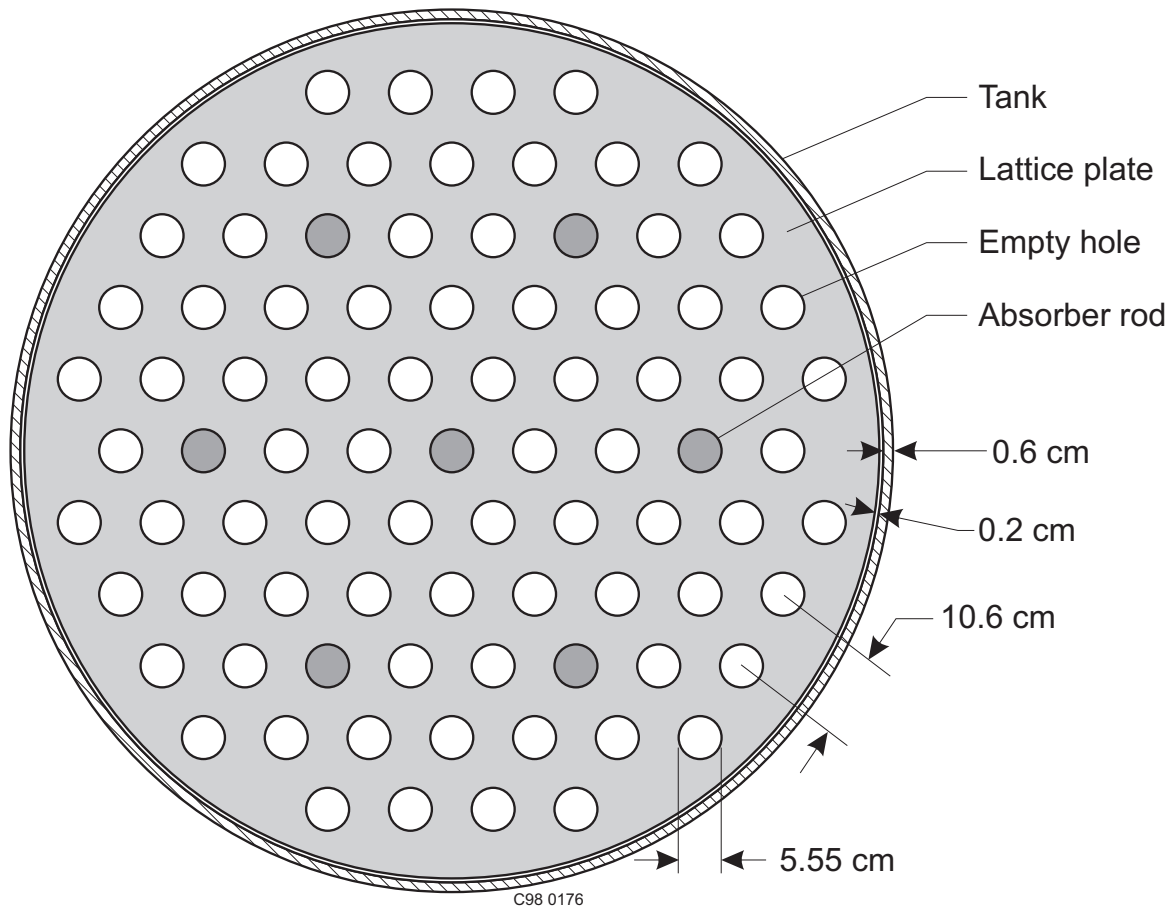


Figure 2. Core Tank of the Critical Assembly.
(Case 2 - horizontal cut through the lower lattice plate.)

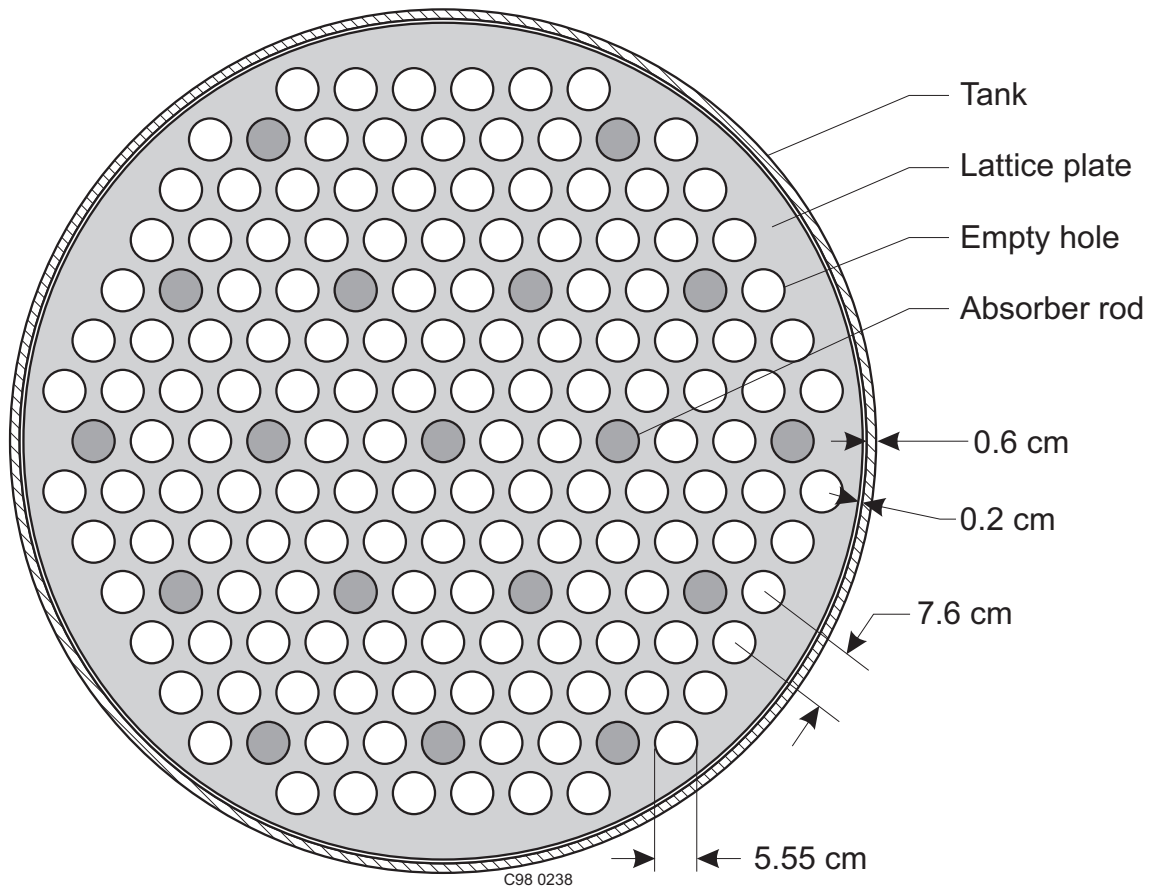


Figure 3. Core Tank of the Critical Assembly.
(Case 3 - horizontal cut through the lower lattice plate.)

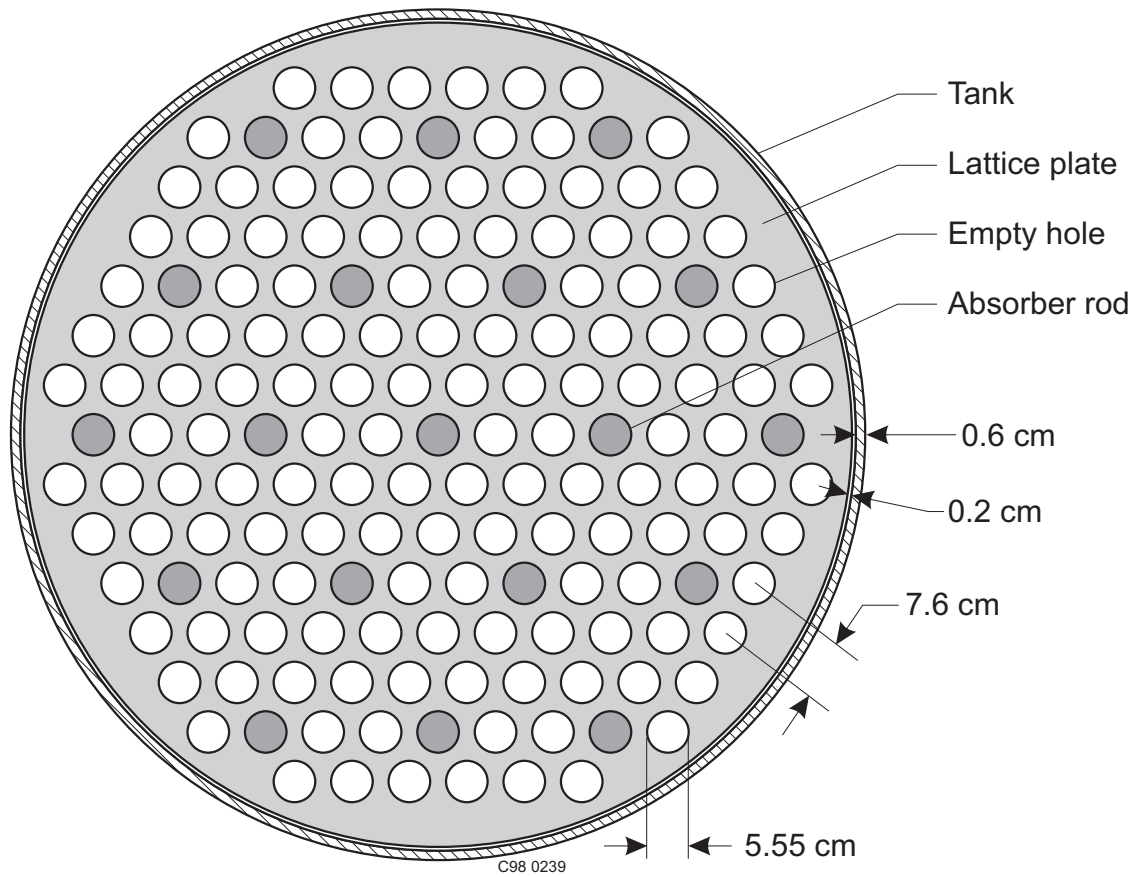


Figure 4. Core Tank of the Critical Assembly.
(Case 4 - horizontal cut through the lower lattice plate.)

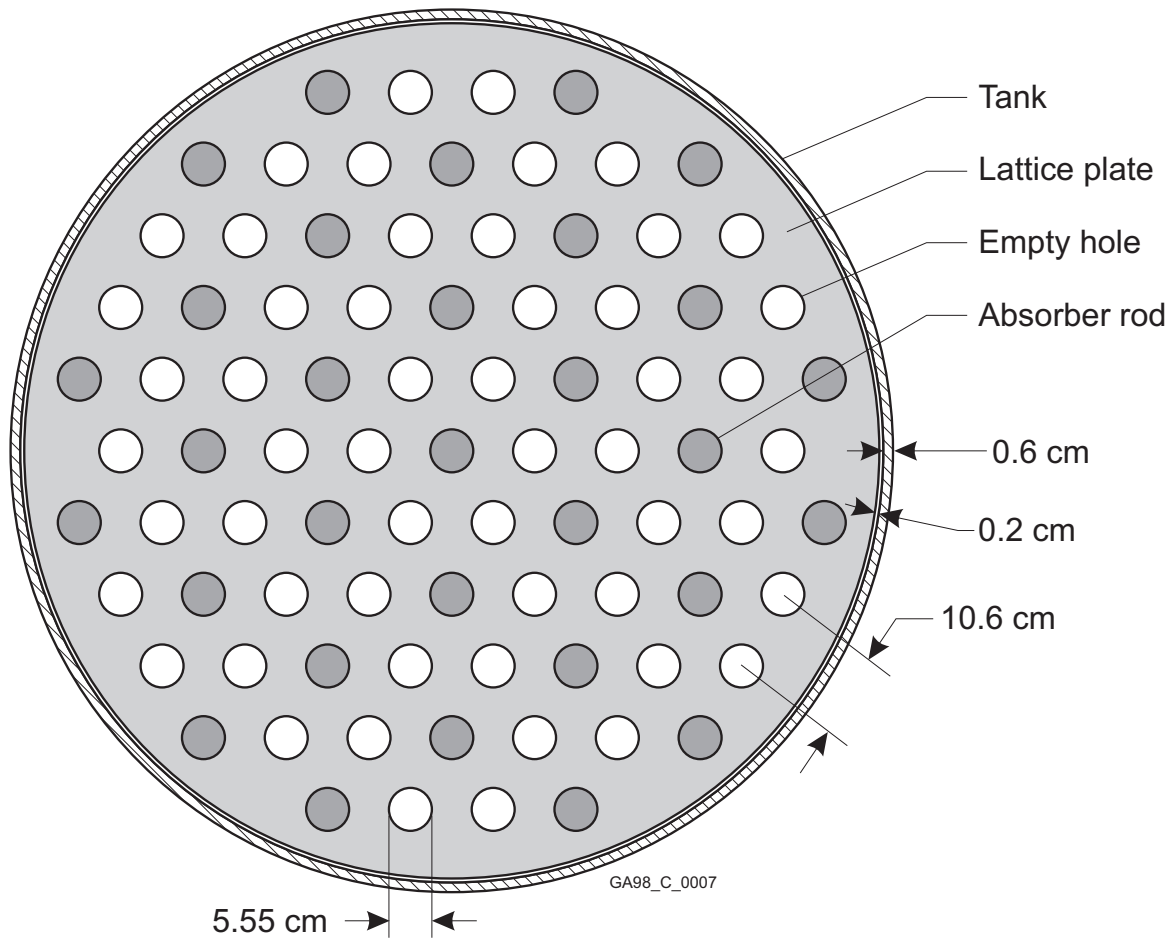


Figure 5. Core Tank of the Critical Assembly.
(Case 5 - horizontal cut through the lower lattice plate.)

The space between the absorber rods was ensured by two stainless steel lattice plates. These plates were 1.7 cm thick. The upper plate was placed on top of the core tank, as shown in Figure 1. The diameter of the upper plate was 115 cm. The lower plate lay on the bottom of the core tank. The diameter of this plate was 109.6 cm.

Two sets of lattice plates were used in these experiments. There were 85 holes 5.55 cm in diameter arranged in a triangular lattice with pitch 10.6 cm for the absorber rods in the first set of the plates, as shown in Figure 2 and Figure 5. There were 163 holes 5.55 cm in diameter arranged in a triangular lattice with pitch 7.6 cm for the absorber rods in the second set of the plates, as shown in Figure 3 and Figure 4.

There was only the lower lattice plate and no upper lattice plate in Case 1, the experiment without absorber rods.

Three boron carbide control rods were suspended above the surface of the solution. The control rods were stainless steel tubes with outer diameter 3 cm and wall and bottom thickness 0.2 cm filled with boron carbide powder. The distance between the critical solution surface and the bottom of the control rods varied from 10 to 20 cm above the solution. Empty holes in the upper lattice plate was used for the boron carbide control rods. Where these rods were located exactly is not known.

Six ionization chambers of the control system were located in six "dry" vertical stainless steel tubes, 6.3 cm in diameter with wall thickness 1.5 mm, symmetrically arranged outside the core tank at a surface-to-surface distance of 20 cm from it. The elements of the control system are not shown in Figure 1.

There was no other apparatus or equipment in the room at the time of the experiment that might affect criticality.

More detailed description of the experimental room and safety/control-rod support system is given in Appendix B in HEU-SOL-THERM-014.

The typical procedure for determining the critical condition was as follows: The core tank, with absorber rods in place, was filled with uranyl nitrate solution portion by portion. The volume of a successive portion was determined from the reciprocal multiplication curve. Portions were small (the smallest sized portion was 0.05 liter) near criticality. The experiments were reported as "critical" but further information about determining the critical condition was not retained. Other solution experiments performed at the facility were taken to a slightly supercritical state.

The experiments were performed at room temperature (approximately 20°C).

1.3 Description of Material Data

Low-enriched uranium was used in the experiments. The enrichment of uranium by ^{235}U according to certificate data was 10.0 wt.% (10.12 at.%) with no shown uncertainty. Therefore, the adopted uncertainty of this value was ± 0.05 wt.%. According to results of mass spectrometric analysis, the

^{235}U content is 10.26 ± 0.07 at.%. An average value was adopted (see Table 2). The ^{234}U at.% was estimated from the mass spectrometric analysis. The uranium isotopic composition is given in Table 2.

Table 2. Uranium Isotopic Composition.

Isotope	at.%
^{234}U	0.09 ± 0.04
^{235}U	10.19 ± 0.10
^{238}U	89.72 ± 0.10
Total:	100.00

The uranium solution was uranyl nitrate, $\text{UO}_2(\text{NO}_3)_2$, from UO_2 dissolved in nitric acid (HNO_3) and diluted with distilled water. The solution density was measured by a float-type hydrometer with the value of one scale division equal to 0.001 g/cm^3 . This value was adopted as the measure of the inaccuracy of the density determination.

The uranium concentration in the solution was measured by weighing U_3O_8 which was prepared by precipitation of ammonium diuranate and its heating. Two or three analyses were done. As usual, two or three small portions of the solution were taken from the core tank (usually from the bottom, middle, and top of the core) after criticality was achieved and then a complete analysis of the solution properties was performed in the chemical laboratory, located in the same building in a neighboring room. The temperature in the chemical laboratory was the same as in the experimental room, therefore the density and uranium concentration of the solution remained constant. The range of the results was, as a rule, smaller than method inaccuracies. So the latter ($\pm 0.5\%$) was adopted as a measure of the uncertainty of the uranium concentration determination. Determinations of chemical-methodology inaccuracies were done previously for a wide range of measured concentrations. (See, for example, HEU-SOL-THERM-014, HEU-SOL-THERM-015, HEU-SOL-THERM-016, HEU-SOL-THERM-017, HEU-SOL-THERM-018, and HEU-SOL-THERM-019.)

The concentration of free nitric acid in the solution was determined by the alkalimetric titration method. The uncertainty of this method is $\pm 1\%$.

Solution properties are given in Table 3.

Table 3. Solution Properties.

Uranium Concentration, grams/liter	Solution Density, grams/cm^3	Concentration of HNO_3 , moles/liter
420.5 ± 2.0	1.581 ± 0.001	0.400 ± 0.004

All impurities in the solution were caused by the impurity of the uranyl nitrate used, not by water impurities. Concentrations of impurities in uranium dioxide determined by chemical analysis are given in Table 4. Uranium content in the uranium dioxide was 87.83 wt.%, according to the certificate on the uranium dioxide. Because the uranium concentration was measured by a chemical method that can detect only uranium,^a the measured uranium concentration did not include impurities.

Table 4. Concentrations of Impurities in Uranium Dioxide.

Element	Wt. %
Fe	0.0470
C	< 0.0200
Ni	0.0054
Si	0.0040
N	0.0010
Mn	0.0018
Cu	0.0007
Co	< 0.0001
B	0.000018
Total:	0.080018

Natural boron was used in the experiments. The density of the boron carbide powder was $1.25 \pm 0.13 \text{ g/cm}^3$.

The cylindrical tanks, absorber rod tubes, and auxiliary parts of the critical assemblies used in these experiments were made of 1X18H10T stainless steel. The composition and density of the stainless steel was determined by normative data on the steel 1X18H10T and was not analyzed specifically. Its composition and density, according to the USSR State Standard 5632-72, is given in Table 5. Cited inaccuracies are caused by possible deviations of the real composition from the normative one.

^a Personal communication, A. A. Krinitsin, analytical chemist IPPE, May, 1997.

Table 5. Stainless Steel^(a) Composition.

Element	Wt. %
Fe	69.1 ± 0.7
Cr	18.0 ± 0.5
Ni	10.0 ± 0.5
Mn	1.5 ± 0.2
Si	0.8 ± 0.1
Ti	0.6 ± 0.1
Total:	100.0

(a) Density is $7.93 \pm 0.02 \text{ g/cm}^3$

1.4 Supplemental Experimental Measurements

No additional experimental data were found.

2.0 EVALUATION OF EXPERIMENTAL DATA

2.1 General Notes

The results of the considered experiments were collected from unclassified internal reports that are published in Reference 1. Additional experimental data were extracted from the internal IPPE reports and retained working books. Many questions were resolved in talks with experimenters and with analytical chemists.

2.2 Influence of Uncertainties in Experimental Parameters on Criticality

The uncertainty of k_{eff} due to reported measurement uncertainties for this set of experiments was calculated. These calculations were performed using one-dimensional cylindrical models of assemblies with critical buckling by means of the CRAB-1 code and three-dimensional HEX-Z models by means of the TRIGEX code with the ABBN-90^a group cross sections set. Some results were additionally checked by the KENO-5A code.

Inaccuracy of the determination of criticality caused by uncertainties of the following parameters was estimated:

- isotopic composition of uranium;
- uranium concentration in solution;
- solution density;
- HNO₃ concentration in solution;
- impurity concentrations in solution;
- absorber material;
- absorber rods pitch;
- critical dimensions.

To determine the reactivity effect due to the uncertainty of the isotopic concentrations, data of Table 2 were used. The ²³⁵U concentration was varied by ± 0.10 at.% with a corresponding change of ²³⁸U atomic density. The corresponding changes in k_{eff} are shown in Table 6 as Δk_e .

To determine the reactivity effect due to the uncertainty of the uranium concentration in solution, data of Table 3 were used. The uranium concentration was varied by the uncertainty value of 2 g/liter. The corresponding changes in k_{eff} are shown in Table 6 as Δk_U .

To determine the reactivity effect due to the uncertainty of the solution density, data of Table 3 were used. The solution density was varied by ± 0.001 g/cm³. The corresponding changes in k_{eff} are shown in Table 6 as Δk_{sol} .

^a RSICC DLC-182 "ABBN-90: Multigroup Constant Set for Calculation of Neutron and Photon Radiation Fields and Functionals, Including the CONSYST2 Program."

To determine the reactivity effect due to the uncertainty of the HNO_3 concentration in solution, data of Table 3 were used. The concentration of HNO_3 was varied by the uncertainty value 0.004 moles/liter. The observed change in k_{eff} did not exceed 0.0003. The corresponding changes in k_{eff} are shown in Table 6 as Δk_{acid} .

As estimated by direct calculations, concentrations of impurities in the solution were so small that their influence on k_{eff} of the considered assemblies is $\sim 0.01\%$. Thus it was not necessary to include the influence of uncertainties of these concentrations in the k_{eff} uncertainty. (The calculations with impurities included the additional nitrate ions bound to the impurities.)

To estimate the reactivity effect due to the uncertainty of the absorber material dimensions, it was supposed that the uncertainty of the outer diameter of the stainless steel tube is $\pm 1\%$ (with constant wall thickness). The volume of the solution was conserved. This leads to change in the surface area of the absorber material. The calculated uncertainty was divided by the square root of the number of rods. The corresponding changes in k_{eff} are shown in Table 6 as Δk_s .

To estimate the reactivity effect due to the uncertainty of the wall thickness of the stainless steel tube of the absorber rod, it was supposed that the uncertainty of the wall thickness is ± 0.05 cm ($\pm 10\%$). The calculated uncertainty was divided by the square root of the number of rods. The corresponding changes in k_{eff} are shown in Table 6 as Δk_w .

To estimate the reactivity effect due to the uncertainty of the pitch of the absorber rods, the rod-separation value was changed by 0.1 cm. The calculated uncertainty was divided by the square root of the number of rods. The corresponding changes in k_{eff} are shown in Table 6 as Δk_p .

To determine the reactivity effect due to the uncertainty of the solution volume, the solution heights were changed by the value that corresponds to the uncertainty of the solution volume given in Table 1 at the constant tank inner radius. The corresponding changes in k_{eff} are shown in Table 6 as Δk_h .

To determine the reactivity effect due to the uncertainty of the core tank inner radius, the value of inner radius was changed by 0.1 cm, and the solution heights correspondingly were changed (thus the critical volumes were conserved). The tank thickness remained constant. The corresponding changes in k_{eff} are shown in Table 6 as Δk_r .

To estimate the reactivity effect due to the uncertainty of the core tank wall and bottom thickness, it was supposed that the uncertainty of the wall and bottom thickness of the tank is $\pm 10\%$. The corresponding changes in k_{eff} are shown in Table 6 as Δk_t .

The influence of the uncertainty in density of the absorber material was calculated by variation of the boron carbide density by ± 0.13 g/cm³. The effect is shown in Table 6 as Δk_d .

The effect of the uncertainty in the isotopic composition of natural boron (19.1% to 20.3% ^{10}B) was also calculated. The effect is shown in Table 6 as Δk_b .

Because the solution density was measured at the time of the experiments, and because the effect of water-reflector density variation due to a large (10°C) temperature change was calculated to be negligible, there is no additional k_{eff} uncertainty due to uncertain temperature.

All Δk 's were determined consistent with formulas provided in Section 3.0.

In Table 6 the constituents of the inaccuracy of k_{eff} are listed for all assemblies of the considered series. In the last column, the summary inaccuracy of k_{eff} is given

$$\Delta k = \sqrt{\Delta k_{\text{e}}^2 + \Delta k_{\text{U}}^2 + \Delta k_{\text{sol}}^2 + \Delta k_{\text{acid}}^2 + \Delta k_{\text{s}}^2 + \Delta k_{\text{w}}^2 + \Delta k_{\text{p}}^2 + \Delta k_{\text{h}}^2 + \Delta k_{\text{r}}^2 + \Delta k_{\text{t}}^2 + \Delta k_{\text{d}}^2 + \Delta k_{\text{b}}^2}$$

Table 6. Constituents of the Inaccuracy of k_{eff} (percents).

Case	Δk_{e}	Δk_{U}	Δk_{sol}	Δk_{acid}	Δk_{s}	Δk_{w}	Δk_{p}	Δk_{h}	Δk_{r}	Δk_{t}	Δk_{d}	Δk_{b}	Δk
1	0.30	0.06	0.03	0.02	--	--	--	0.08	0.15	0.08	--	--	0.36
2	0.30	0.04	0.02	0.03	0.03	0.08	0.05	0.09	0.13	0.06	0.03	0.05	0.37
3	0.31	0.05	0.02	0.03	0.08	0.14	0.05	0.08	0.11	0.05	0.06	0.06	0.40
4	0.28	0.05	0.02	0.03	0.09	0.19	0.05	0.08	0.11	0.05	0.06	0.06	0.40
5	0.29	0.05	0.02	0.03	0.17	0.25	0.05	0.07	0.10	0.04	0.09	0.08	0.46

Because sufficient data are known and uncertainties have been quantified as relatively small, the five configurations are acceptable benchmark experiments.

3.0 BENCHMARK SPECIFICATIONS

3.1 Description of Model

3.1.1 Description of Simplifications - The benchmark models were based on a series of simplifications to the actual experimental configurations. The following experimental details were omitted, after evaluating their influence on criticality during the benchmark specifications:

- impurities in fissile solution;
- small constructive details in and around the assembly, including the upper lattice plate, and reflection of neutrons from the room walls.

3.1.2 Influence of the Elimination of Impurities - As was pointed out in Section 2, the influence of solution impurities on k_{eff} was negligibly small. Thus impurities are not included in the benchmark models and no correction of k_{eff} was made.

3.1.3 Influence of the Elimination of the Experimental Surroundings - Because the thick water reflector was used, the experimental surroundings (i.e., room wall, ionization-chamber tubes) had no noticeable influence on the k_{eff} of the considered assemblies. The effect of removing the control rods located above the solution surface was estimated using the KENO-5A code. The effect was found equal to less than 0.1% with approximately the same uncertainty. Because the exact positions of the control rods in the experiments are not known, the additional 0.1% uncertainty is included in the benchmark-model k_{eff} 's. The influence of removing the fill and drain tubes was estimated earlier (see HEU-SOL-THERM-014, HEU-SOL-THERM-015, HEU-SOL-THERM-016, HEU-SOL-THERM-017, HEU-SOL-THERM-018, and HEU-SOL-THERM-019) using perturbation theory. It was assumed that in the currently considered series the fill and drain tubes also do not influence criticality. Making the top surfaces of the two tanks and absorber rods coplanar and thereby reducing the height of the reflector tank by 13 cm and of the rods by 5 cm to simplify the model was judged to have negligible effect.

3.2 Dimensions

The benchmark models of the experiments are shown in Figures 6 through 13. The model is two open-top coaxial stainless steel cylindrical tanks. The core tank has inner diameter 110.0 cm, wall thickness 0.6 cm, bottom thickness 1.5 cm, is 250.0 cm tall, filled with solution of uranyl nitrate to the height (measured from the inner surface of the tank bottom) shown in Table 7. The reflector tank, with inner diameter 198.4 cm, wall thickness 0.8 cm, bottom thickness 1.0 cm, 287.0 cm tall, is filled with water to the height (measured from the inner surface of the core tank bottom) 108.0 cm. The distance between the inner surface of the reflector tank bottom and the outer surface of the core tank bottom is 36.0 cm.

There is only solution in the core tank (no absorber rods) in Case 1. The boron absorber rods are inserted in the core tank in Cases 2 - 5. The number of the boron rods in the tank is shown in Table 7. The arrangement of the absorber rods in the tank is shown in Figures 9 - 13.

Table 7. Geometrical Sizes of Benchmark Models.

Case Number	Number of Absorber Rods	Number of Holes in the Lattice Plate	Solution Height, cm
1	none	85	23.4174
2	7	85	26.1371
3	18	163	28.7180
4	19	163	29.2573
5	31	85	36.3693

The boron absorber rods are stainless steel tubes with outer diameter of 5.5 cm, wall thickness 0.5 cm, and bottom thickness 0.7 cm, 248.5 cm long, filled with natural boron carbide. The absorber rods extend to the bottom of the core tank. The top surface of the absorber rods is coplanar with the top surfaces of the core and reflector tanks.

The lower stainless steel lattice plate is also included in the benchmark models. This plate is lying on the bottom of the core tank. It has diameter 109.6 cm and is 1.7 cm thick. Two kinds of lattice plates are used. There are 85 holes in the first type, arranged in a hexagonal lattice with pitch 10.6 cm. There are 163 holes in the second type, arranged in a hexagonal lattice with pitch 7.6 cm. The type of the lattice plate which is used in each particular case is shown in Table 7. The arrangement of the holes in the plates is shown in Figures 9 - 13. The holes have diameter 5.55 cm.

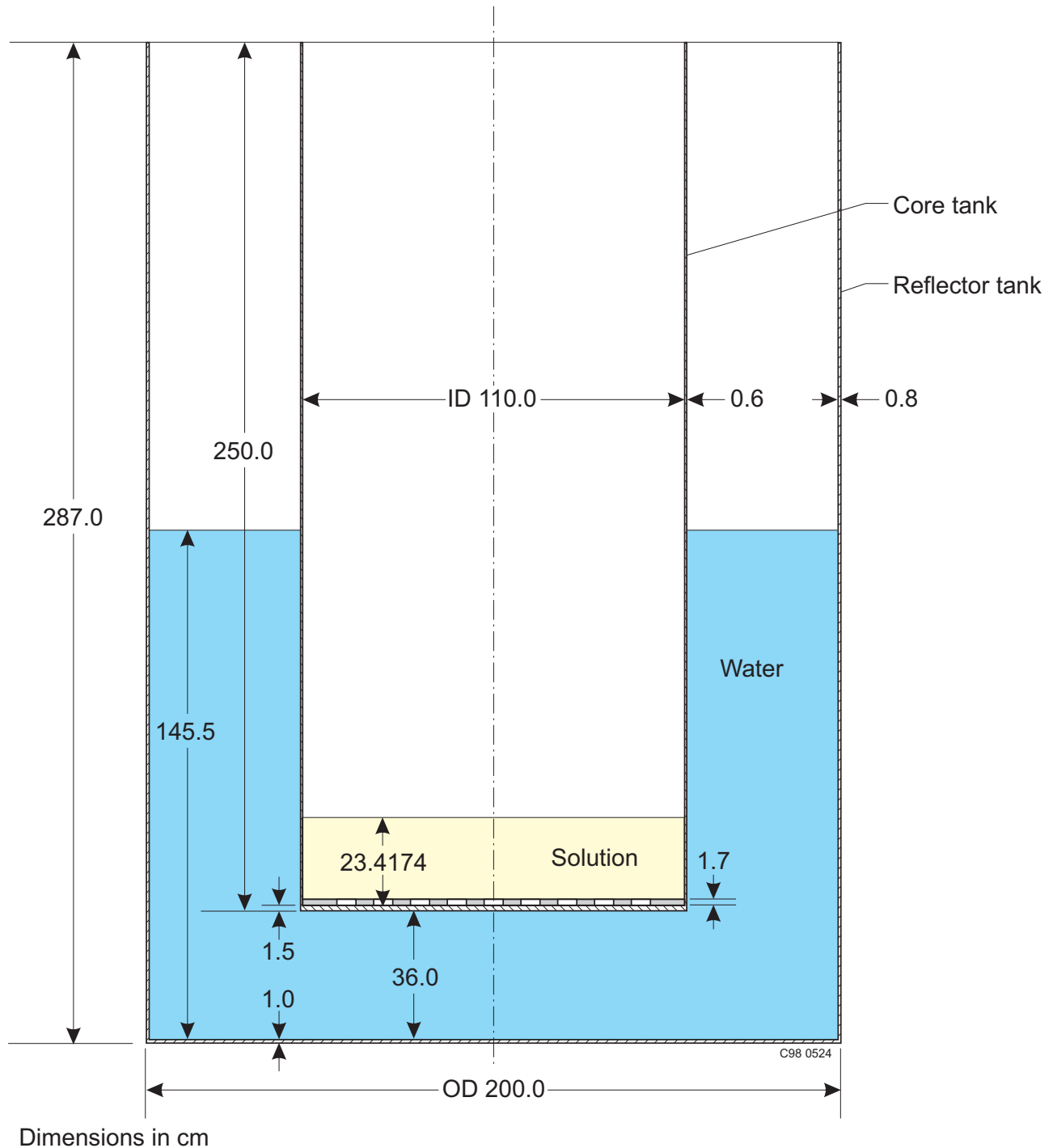


Figure 6. Benchmark Model of Case 1 (vertical cut).

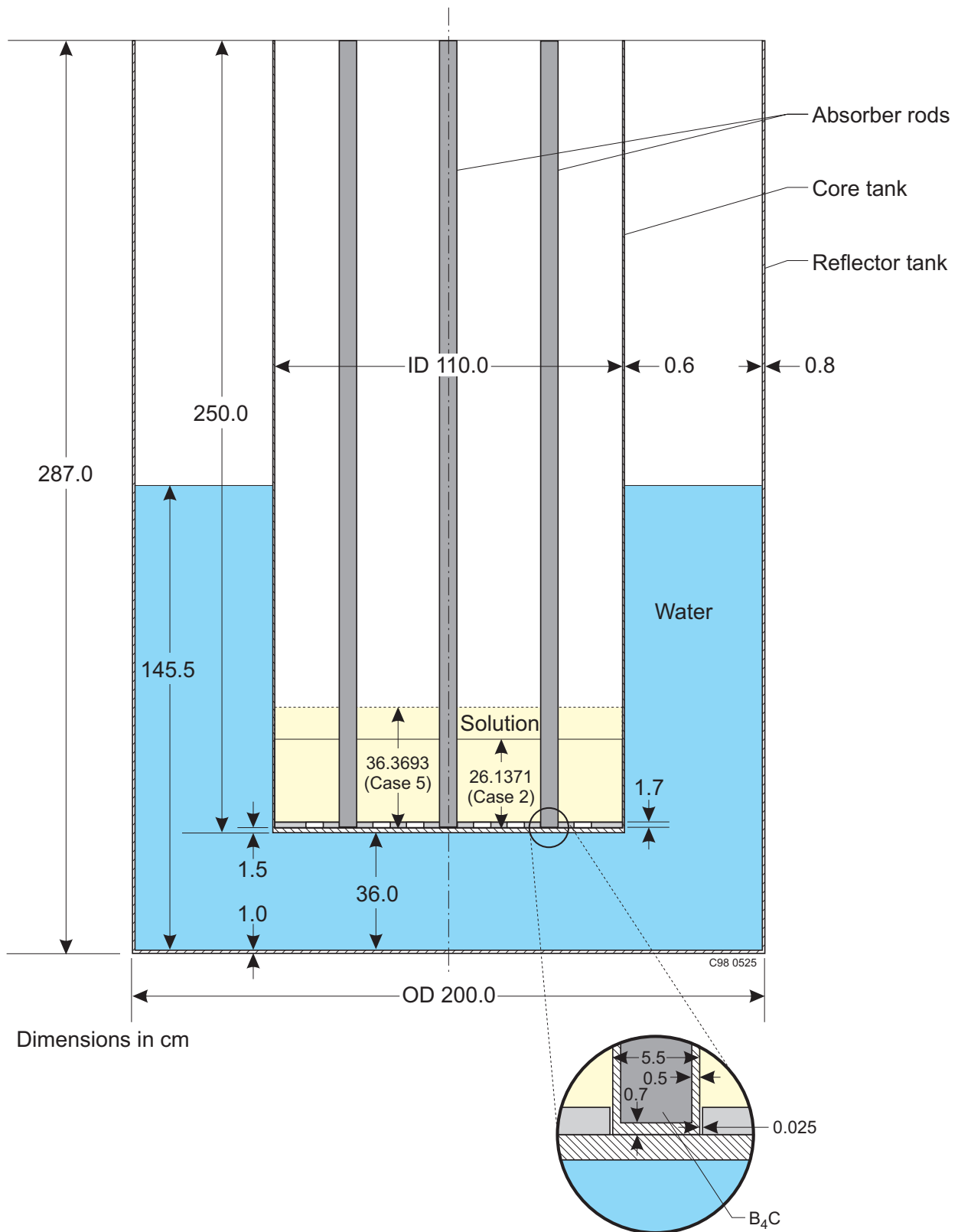


Figure 7. Benchmark Model of Cases 2 and 5 (vertical cut).

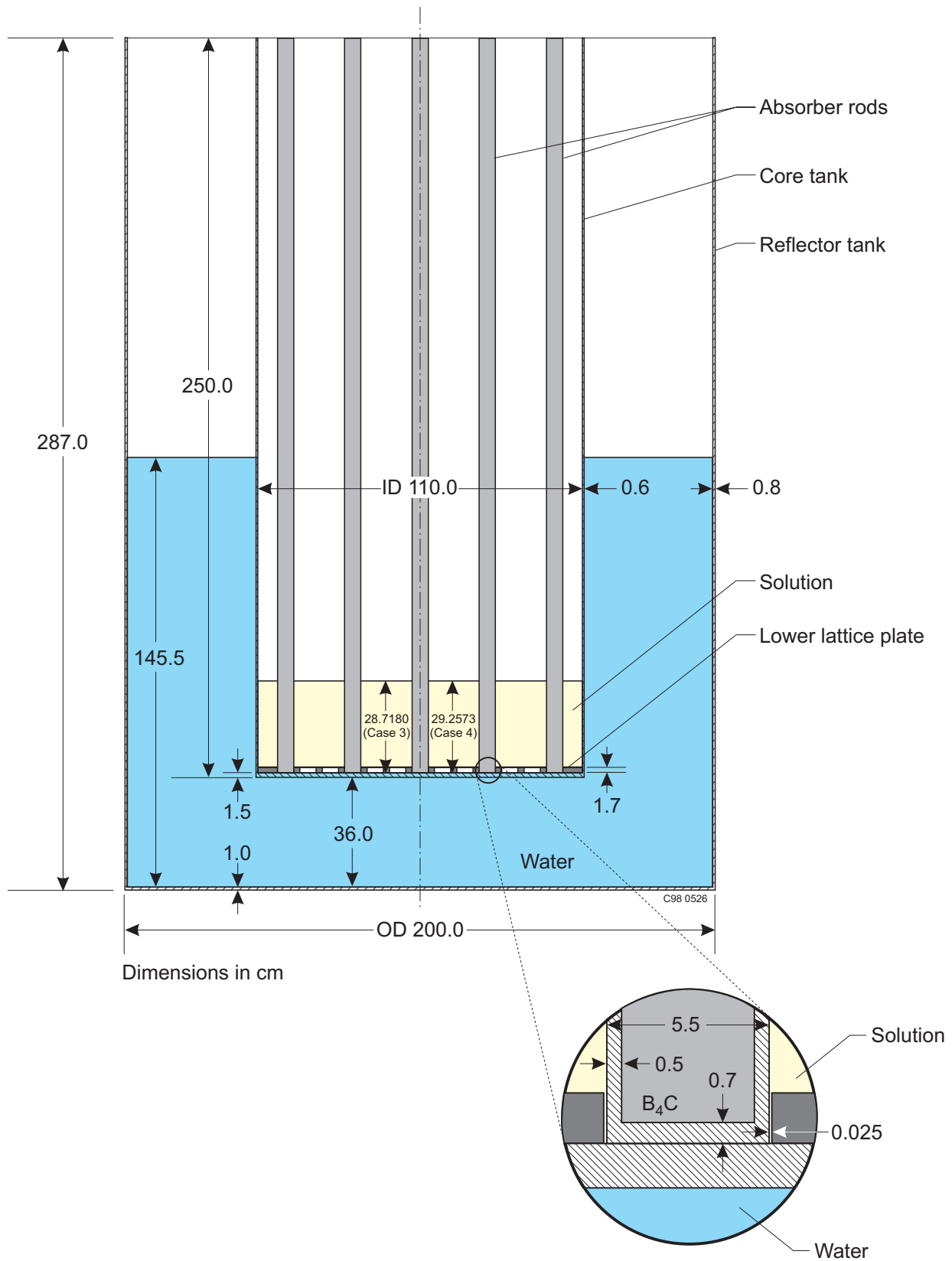


Figure 8. Benchmark Model of Cases 3 and 4 (vertical cut).

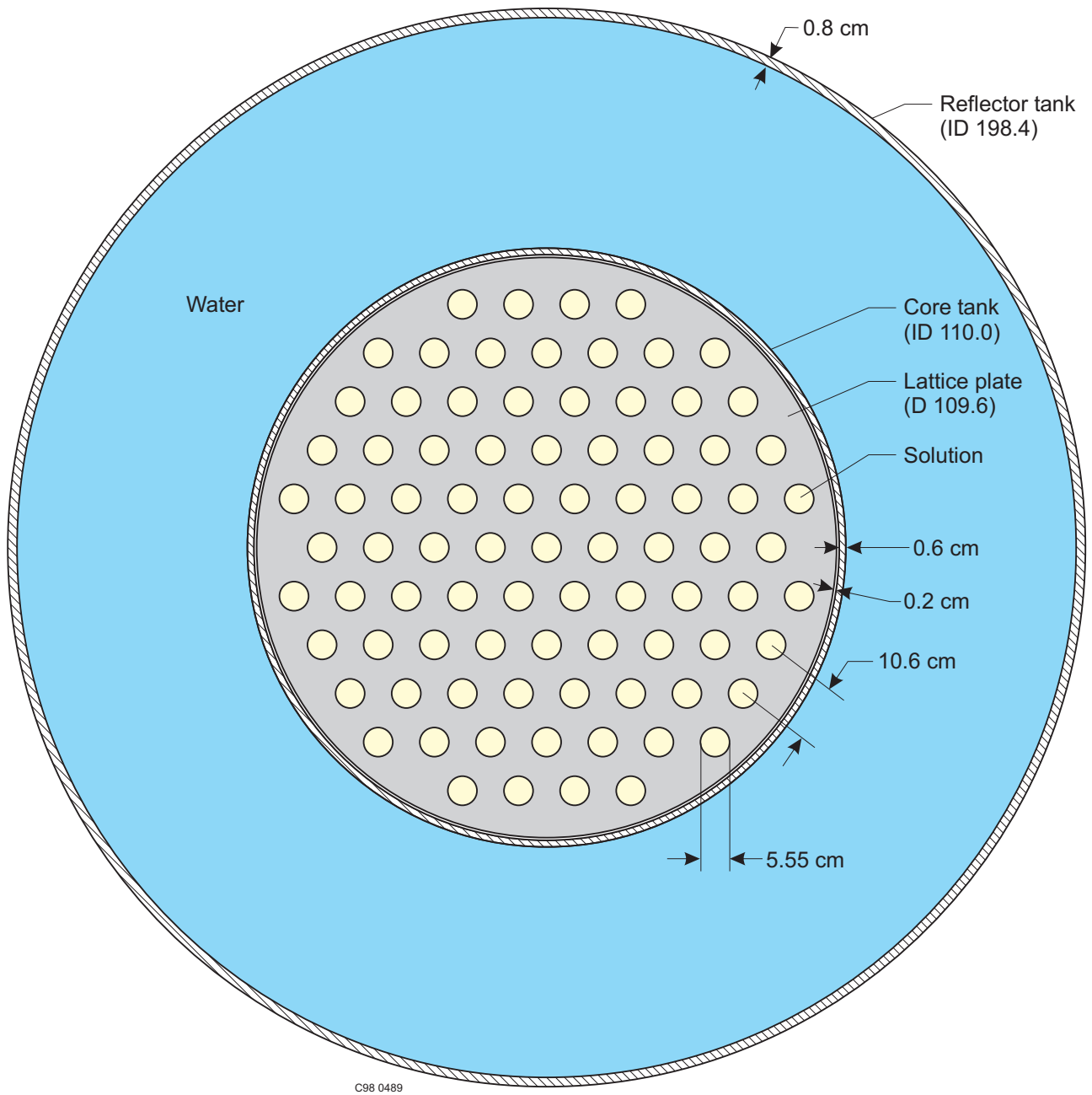


Figure 9. Benchmark Model of Case 1 (horizontal cut through lattice plate).

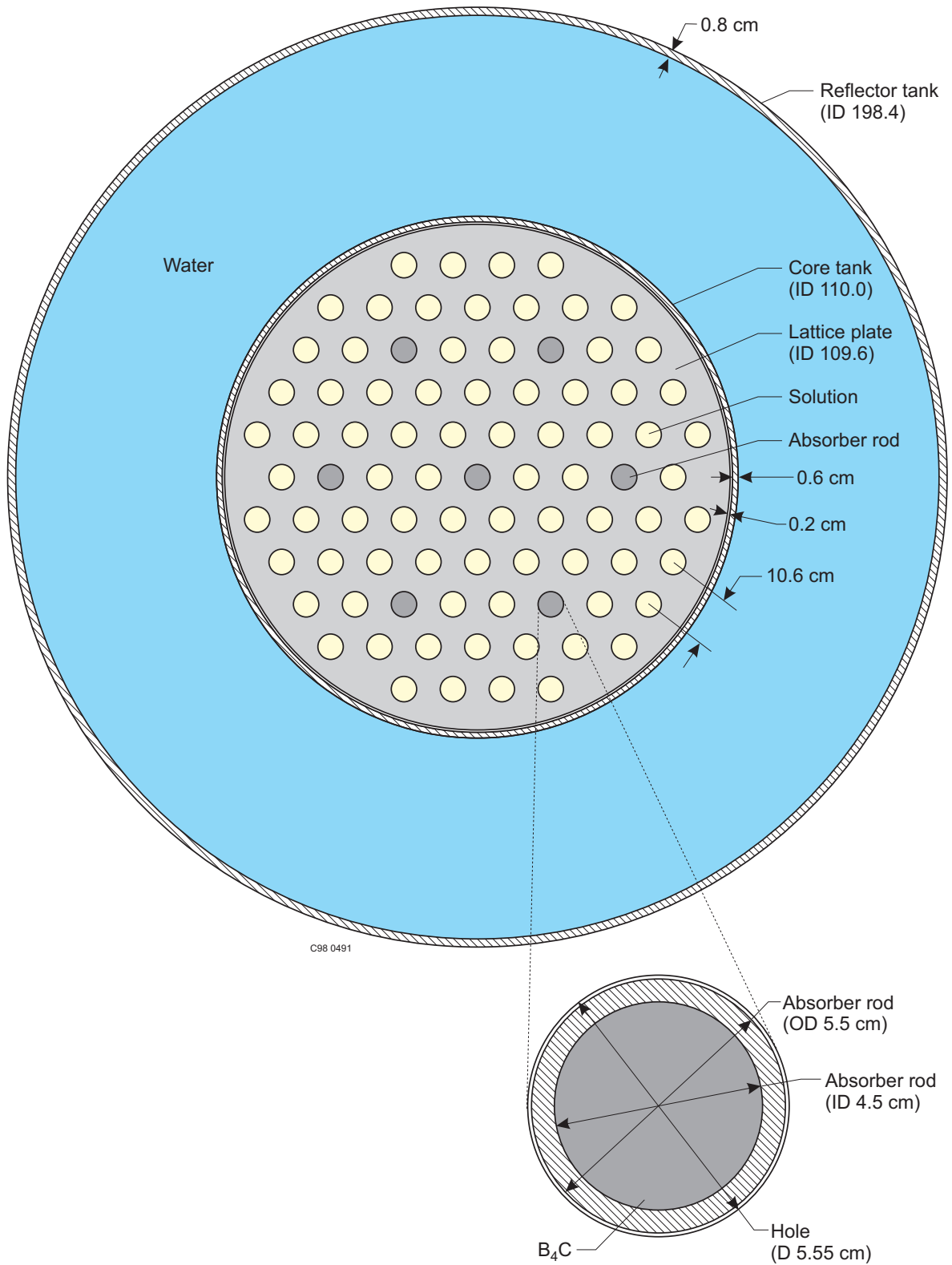


Figure 10. Benchmark Model of Case 2 (horizontal cut through lattice plate).

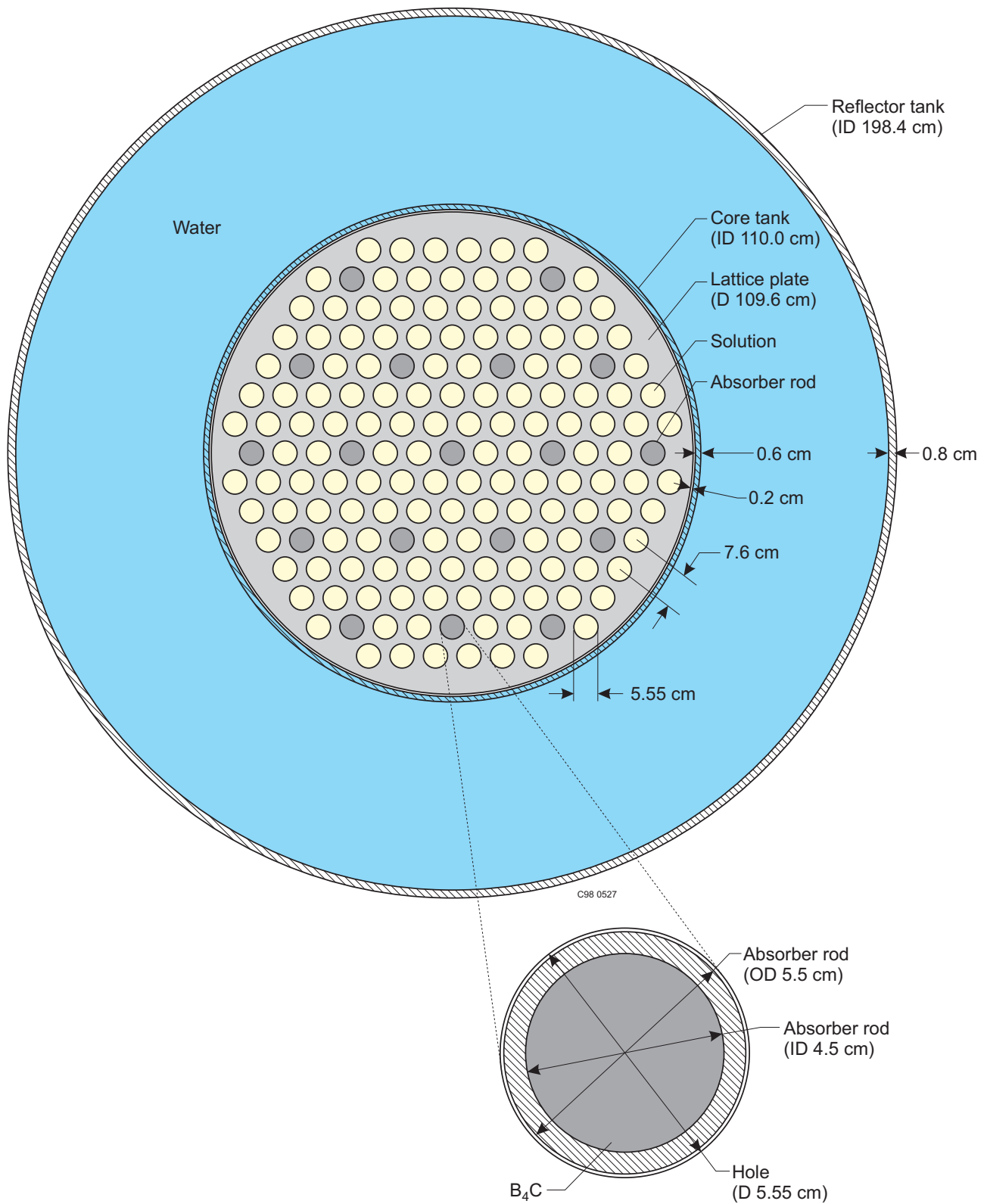


Figure 11. Benchmark Model of Case 3 (horizontal cut through lattice plate).

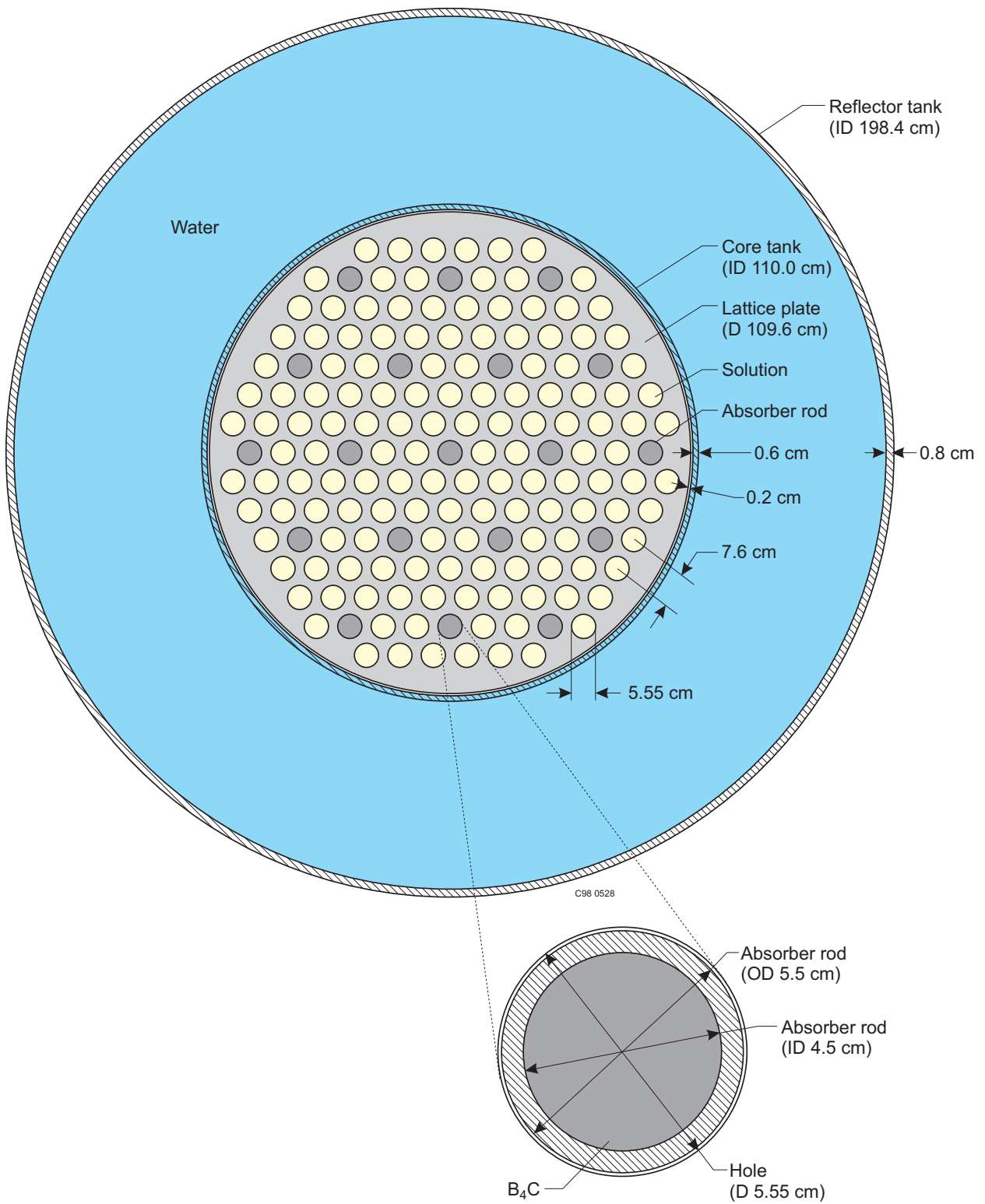


Figure 12. Benchmark Model of Case 4 (horizontal cut through lattice plate).

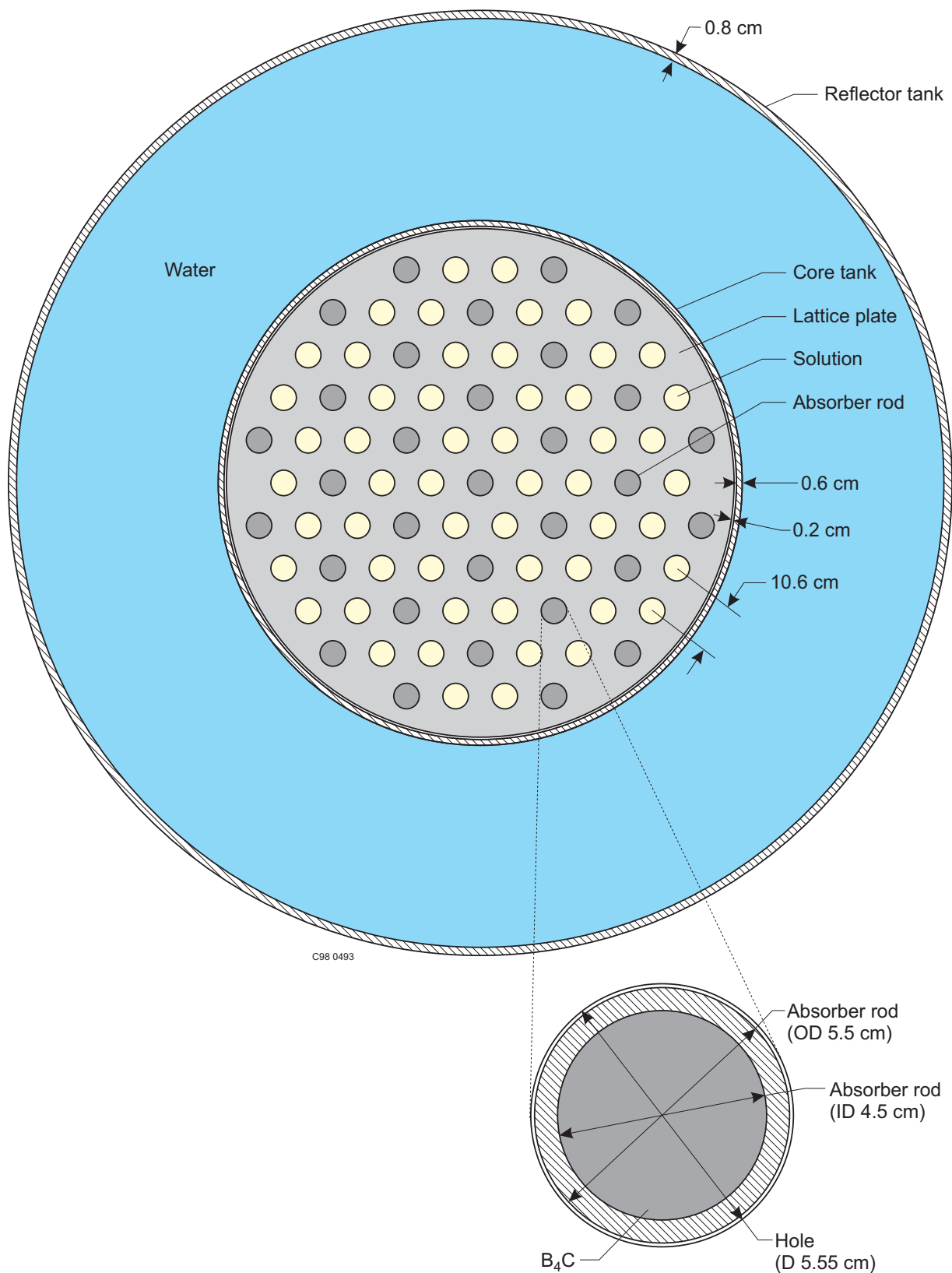


Figure 13. Benchmark Model of Case 5 (horizontal cut through lattice plate).

3.3 Material Data

3.3.1 Core Nuclear Densities - The nuclear densities of the individual uranium isotopes are given by

$$N_i = \frac{\rho_U N_A A_{f,i}}{A_{w,U}}, \text{ for } i = {}^{234}\text{U}, {}^{235}\text{U}, {}^{238}\text{U}.$$

where N_i is an atomic density of isotope "i";

$A_{f,i}$ is the atom fraction of isotope "i" (see at.%'s in Table 2);

ρ_U is the total density of uranium in solution in g/cm³ (see Table 3);

$A_{w,U}$ is the average atomic weight of the uranium

($A_{w,U} = A_{w,234} \times A_{f,234} + A_{w,235} \times A_{f,235} + A_{w,238} \times A_{f,238}$);

N_A is Avogadro's number.

The mass density of uranyl nitrate, $\text{UO}_2(\text{NO}_3)_2$, is given by

$$\rho_{\text{UO}_2(\text{NO}_3)_2} = \frac{\rho_U M_{w,\text{UO}_2(\text{NO}_3)_2}}{A_{w,U}}$$

where $A_{w,U}$ is the average atomic weight of the uranium, and

$M_{w,\text{UO}_2(\text{NO}_3)_2}$ is the molecular weight of the uranyl nitrate.

The mass density of nitric acid, HNO_3 , in grams per cubic centimeter is given by

$$\rho_{\text{HNO}_3} = 0.001 \cdot M_{w,\text{HNO}_3} \cdot M_{\text{HNO}_3}$$

where M_{w,HNO_3} is the molecular weight of nitric acid, and

M_{HNO_3} is the concentration of HNO_3 in moles/liter (see Table 3).

The mass density of water in solution can be then determined by difference

$$\rho_{\text{H}_2\text{O}} = \rho_{\text{solution}} - \rho_{\text{UO}_2(\text{NO}_3)_2} - \rho_{\text{HNO}_3}$$

The total atom densities of nitrogen (N_N), oxygen (N_O), and hydrogen (N_H) are then determined by the following formulas:

$$N_N = 2 (N_{\text{U}234} + N_{\text{U}235} + N_{\text{U}236} + N_{\text{U}238}) + 0.001 \cdot M_{\text{HNO}_3} \cdot N_A$$

$$N_O = 2 (N_{\text{U}235} + N_{\text{U}234} + N_{\text{U}236} + N_{\text{U}238}) + \frac{\rho_{\text{H}_2\text{O}} N_A}{M_{w,\text{H}_2\text{O}}} + 3 N_N$$

$$N_H = N_A \left(2 \frac{\rho_{H_2O}}{M_{w,H_2O}} + \frac{\rho_{HNO_3}}{M_{w,HNO_3}} \right)$$

Here M_{w,H_2O} is the molecular weight of water, and ρ_{H_2O} , ρ_{HNO_3} in grams per cubic centimeter are calculated by the above formulas.

Using the above equations, solution atom densities were calculated. Results of these calculations are given in Table 8.

Table 8. Calculated Atom Densities for 420.5 gU/l Solution.

Element	Atom Density, atoms/(barn \times cm)
^{234}U	9.5863×10^{-7}
^{235}U	1.0854×10^{-4}
^{238}U	9.5565×10^{-4}
N	2.3712×10^{-3}
O	3.7970×10^{-2}
H	5.7694×10^{-2}

3.3.2 Boron Carbide Atom Densities - Atom densities of boron carbide B_4C were calculated using the following isotopic composition^a of natural boron: ^{10}B - 19.9 at.%; ^{11}B - 80.1 at.%, and density of 1.25 g/cm^3 . Boron carbide atom densities are given in Table 9.

Table 9. Atom Densities of Boron Carbide.

Element	Atom Density, atoms/(barn \times cm)
^{10}B	1.0844×10^{-2}
^{11}B	4.3648×10^{-2}
C	1.3623×10^{-2}

3.3.3 Water Reflector Atom Density - The density of the water was selected for the temperature and pressure at which the experiment was conducted. The temperature of the water was 20°C (293 K), and pressure was approximately 0.1 MPa at the time of the experiment. Therefore the estimated water density^b was 0.9983 g/cm^3 . Water atom densities are given in Table 10.

^a "Chart of the Nuclides," Fourteen Edition, General Electric Company, 1989.

^b Ривкин С. Л., Александров А. А. Термодинамические свойства воды и водяного пара: Справочник. М.: Энергоатомиздат, 1984. (Rivkin S. L., Aleksandrov A. A., "Thermodynamic Properties of Water and Steam: Handbook." Moscow. Energoatomizdat, 1984, in Russian.)

Table 10. Atom Densities of Water Reflector.

Element	Atom Density, atoms/(barn \times cm)
H	6.6742×10^{-2}
O	3.3371×10^{-2}

3.3.4 Atom Densities of the Structure - Atom densities of stainless steel constituents were calculated from weight percents given in Table 5 and density of 7.93 g/cm^3 . Data are given in Table 11.

Table 11. Atom Densities for Stainless Steel.

Element	Atom Density, atoms/(barn \times cm)
Fe	5.9088×10^{-2}
Cr	1.6532×10^{-2}
Ni	8.1369×10^{-3}
Mn	1.3039×10^{-3}
Si	1.3603×10^{-3}
Ti	5.9844×10^{-4}

3.4 Temperature Data

All experiments were performed at room temperature (approximately 20°C). The temperature of 300 K was assumed for calculations.

3.5 Experimental and Benchmark-Model k_{eff}

Experimental critical parameters correspond to criticality: $k_{\text{eff}} = 1$. Uncertainties of experimental data result in an uncertainty of k_{eff} , as described in Section 2. Benchmark-model simplifications led to the additional 0.1% uncertainty of k_{eff} . Benchmark-model k_{eff} 's are shown in Table 12.

Table 12. Benchmark-Model k_{eff} Values.

Case Number	Benchmark-Model k_{eff}
1	1.0000 ± 0.0037
2	1.0000 ± 0.0038
3	1.0000 ± 0.0041
4	1.0000 ± 0.0041
5	1.0000 ± 0.0047

4.0 RESULTS OF SAMPLE CALCULATIONS



The results of calculations of k_{eff} of the benchmark models are given in Tables 13.a and 13.b. Details of the calculations including code input listings are provided in Appendix A.

Table 13.a. Sample Calculation Results (Russian Federation).^(a)

Code (Cross Section Set) → Configuration ↓	KENO (ABBN-93)
1	0.9995 ± 0.0007
2	1.0082 ± 0.0007
3	1.0007 ± 0.0007
4	1.0025 ± 0.0007
5	1.0055 ± 0.0007

(a) Results supplied by Yevgeniy Rozhikhin, IPPE.

Table 13.b. Sample Calculation Results (United States).^(a)

Code (Cross Section Set) → Configuration ↓	KENO (27-Group ENDF/B-IV)	MCNP (Continuous Energy ENDF/B-V)
1	1.0005 ± 0.0007	1.0014 ± 0.0007
2	1.0062 ± 0.0007	1.0071 ± 0.0007
3	0.9980 ± 0.0007	1.0027 ± 0.0007
4	1.0012 ± 0.0007	1.0010 ± 0.0007
5	1.0036 ± 0.0007	1.0024 ± 0.0007

(a) Results supplied by Yevgeniy Rozhikhin, IPPE.

5.0 REFERENCES

1. Дубовский Б. Г., Камаев А. В., Кузнецов Ф. М. и др. Критические параметры систем с делющимися веществами и ядерная безопасность. Справочник. Москва, Атомиздат, 1966. (Dubovskii B. G., Kamaev A. V., Kuznetsov F. M. et al., "Critical Parameters of Systems with Fissile Materials and Nuclear Safety. Handbook." Moscow. Atomizdat, 1966, in Russian.)

APPENDIX A: TYPICAL INPUT LISTINGS

A.1 KENO Input Listings

299-Group ABBN-93 Cross Sections

KENO-5A with the ABBN-93 299-group cross sections was run using 1020 generations of neutrons with 1000 neutrons each. Twenty generations were skipped before averaging, so the result is the average of 1,000,000 neutron histories.

27-Group ENDF/B-IV Cross Sections

The 27-group inputs given below are for the KENO-5A code as run with the ORNL SCALE4.3 code system using the CSAS option. The problems were run with 1,000,000 active and 20,000 inactive histories (1000 active generations, twenty skipped generations, and 1000 histories per generation).

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 1 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections).

```
=csas25
leu-sol-therm-006-01
27groupndf4 infhommedium
u-234 1 0 9.5863E-07 end
u-235 1 0 1.0854E-04 end
u-238 1 0 9.5565E-04 end
n 1 0 2.3712E-03 end
o 1 0 3.7970E-02 end
h 1 0 5.7694E-02 end
fe 2 0 5.9088E-02 end
cr 2 0 1.6532E-02 end
ni 2 0 8.1369E-03 end
mn 2 0 1.3039E-03 end
si 2 0 1.3603E-03 end
ti 2 0 5.9844E-04 end
h 3 0 6.6742E-02 end
o 3 0 3.3371E-02 end
b-10 4 0 1.0844E-02 end
b-11 4 0 4.3648E-02 end
c 4 0 1.3623E-02 end
end comp
leu-sol-therm-006-01
read param
tme=360
gen=1020
npg=1000
nsk=20
fdn=yes
far=yes
plt=no
end param
read geometry
unit 11
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 5.3 -5.3 3. -3. 1.7 0.
' . . . . .
unit 61
array 61 0. 0. 0.
unit 71
array 71 0. 0. 0.
unit 81
array 81 0. 0. 0.
unit 91
array 91 0. 0. 0.
unit 101
array 101 0. 0. 0.
' . . . . .
unit 999
cylinder 0 1 55. 248.5 108.
cylinder 2 1 55.6 248.5 108.
' . . . . .
global unit 1000
cylinder 2 1 54.8 1.7 0.
hole 61 -21.2 -49.89934640058 0.
hole 71 -37.1 -39.71947712046 0.
hole 81 -42.4 -30.53960784035 0.
hole 91 -47.7 -21.35973856023 0.
hole 101 -53. -12.17986928012 0.
hole 91 -47.7 -3. 0.
hole 101 -53.0 6.179869280115 0.
hole 91 -47.7 15.35973856023 0.
hole 81 -42.4 24.53960784035 0.
hole 71 -37.1 33.71947712046 0.
hole 61 -21.2 42.89934640058 0.
```

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 1 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```
cylinder 1 1 55.          23.4174 0.
cylinder 0 1 55.          108.  0.
cylinder 2 1 55.6         108.  -1.5
cylinder 3 1 99.2         108.  -37.5
cylinder 0 1 99.2         248.5 -37.5
hole 999  0.  0.  0.
cylinder 2 1 100.         248.5 -38.5
end geometry
read array
ara= 61  nux= 4 nuy= 1 nuz=1  fill f11 end fill
ara= 71  nux= 7 nuy= 1 nuz=1  fill f11 end fill
ara= 81  nux= 8 nuy= 1 nuz=1  fill f11 end fill
ara= 91  nux= 9 nuy= 1 nuz=1  fill f11 end fill
ara=101  nux=10 nuy= 1 nuz=1  fill f11 end fill
end array
read plot
scr=yes
ttl='x-z slice at y=0.0 with x across and z down'
xul=-101.0 yul=  0.0  zul= 250.0
xlr= 101.0 ylr=  0.0  zlr=-40.0
uax=1.0 wdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=10.0 with x across and y down'
xul=-101.0 yul= 101.0  zul= 10.0
xlr= 101.0 ylr=-101.0  zlr= 10.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=1.0 with x across and y down'
xul=-60.0 yul= 60.0  zul= 1.0
xlr= 60.0 ylr=-60.0  zlr= 1.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
end plot
end data
end
```

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 2 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections).

```
=csas25
leu-sol-therm-006-02
27groupndf4 infhommedium
u-234 1 0 9.5863E-07 end
u-235 1 0 1.0854E-04 end
u-238 1 0 9.5565E-04 end
n 1 0 2.3712E-03 end
o 1 0 3.7970E-02 end
h 1 0 5.7694E-02 end
fe 2 0 5.9088E-02 end
cr 2 0 1.6532E-02 end
ni 2 0 8.1369E-03 end
mn 2 0 1.3039E-03 end
si 2 0 1.3603E-03 end
ti 2 0 5.9844E-04 end
h 3 0 6.6742E-02 end
o 3 0 3.3371E-02 end
b-10 4 0 1.0844E-02 end
b-11 4 0 4.3648E-02 end
c 4 0 1.3623E-02 end
end comp
leu-sol-therm-006-02
read param
tme=360
gen=1020
npg=1000
nsk=20
fdn=yes
far=yes
plt=no
end param
read geometry
unit 11
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 5.3 -5.3 3. -3. 1.7 0.
unit 21
cylinder 4 1 2.25 1.7 0.7
cylinder 2 1 2.75 1.7 0.
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 5.3 -5.3 3. -3. 1.7 0.
unit 22
cylinder 4 1 2.25 26.1371 1.7
cylinder 2 1 2.75 26.1371 1.7
unit 23
cylinder 4 1 2.25 108. 26.1371
cylinder 2 1 2.75 108. 26.1371
unit 24
cylinder 4 1 2.25 248.5 108.
cylinder 2 1 2.75 248.5 108.
' . . . . .
unit 61
array 61 0. 0. 0.
unit 71
array 71 0. 0. 0.
unit 81
array 81 0. 0. 0.
unit 91
array 91 0. 0. 0.
unit 101
array 101 0. 0. 0.
unit 111
array 111 0. 0. 0.
' . . . . .
unit 999
```

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 2 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

cylinder 0 1 55.          248.5 108.
hole 24  0.  0.          0.
hole 24 -31.8 0.          0.
hole 24 -15.9 27.53960784035 0.
hole 24  15.9 27.53960784035 0.
hole 24  31.8 0.          0.
hole 24  15.9 -27.53960784035 0.
hole 24 -15.9 -27.53960784035 0.
cylinder 2 1 55.6          248.5 108.
global unit 1000
cylinder 2 1 54.8          1.7 0.
hole 61 -21.2 -49.89934640058 0.
hole 71 -37.1 -39.71947712046 0.
hole 81 -42.4 -30.53960784035 0.
hole 91 -47.7 -21.35973856023 0.
hole 101 -53. -12.17986928012 0.
hole 111 -47.7 -3.          0.
hole 101 -53.0 6.179869280115 0.
hole 91 -47.7 15.35973856023 0.
hole 81 -42.4 24.53960784035 0.
hole 71 -37.1 33.71947712046 0.
hole 61 -21.2 42.89934640058 0.
cylinder 1 1 55.          26.1371 0.
hole 22  0.  0.          0.
hole 22 -31.8 0.          0.
hole 22 -15.9 27.53960784035 0.
hole 22  15.9 27.53960784035 0.
hole 22  31.8 0.          0.
hole 22  15.9 -27.53960784035 0.
hole 22 -15.9 -27.53960784035 0.
cylinder 0 1 55.          108. 0.
hole 23  0.  0.          0.
hole 23 -31.8 0.          0.
hole 23 -15.9 27.53960784035 0.
hole 23  15.9 27.53960784035 0.
hole 23  31.8 0.          0.
hole 23  15.9 -27.53960784035 0.
hole 23 -15.9 -27.53960784035 0.
cylinder 2 1 55.6          108. -1.5
cylinder 3 1 99.2          108. -37.5
cylinder 0 1 99.2          248.5 -37.5
hole 999 0.  0.  0.
cylinder 2 1 100.          248.5 -38.5
end geometry
read array
ara= 61 nux= 4 nuy= 1 nuz=1 fill f11 end fill
ara= 71 nux= 7 nuy= 1 nuz=1 fill f11 end fill
ara= 81 nux= 8 nuy= 1 nuz=1 fill 2r11 21 11 n4 end fill
ara= 91 nux= 9 nuy= 1 nuz=1 fill f11 end fill
ara=101 nux=10 nuy= 1 nuz=1 fill f11 end fill
ara=111 nux= 9 nuy= 1 nuz=1 fill 11 21 11 2q3 end fill
end array
read plot
scr=yes
ttl='x-z slice at y=0.0 with x across and z down'
xul=-101.0 yul= 0.0 zul= 250.0
xlr= 101.0 ylr= 0.0 zlr=-40.0
uax=1.0 wdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=10.0 with x across and y down'
xul=-101.0 yul= 101.0 zul= 10.0
xlr= 101.0 ylr=-101.0 zlr= 10.0
uax=1.0 vdn=-1.0

```

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 2 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```
nax=800 lpi=10 end
ttl='x-y slice at z=1.0 with x across and y down'
xul= -60.0 yul= 60.0 zul= 1.0
xlr= 60.0 ylr= -60.0 zlr= 1.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
end plot
end data
end
```

LEU-SOL-THERM-006

KENO-5A Input Listing for Case 3 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections).

```
=csas25
leu-sol-therm-006-03
27groupndf4 infhommedium
u-234 1 0 9.5863E-07 end
u-235 1 0 1.0854E-04 end
u-238 1 0 9.5565E-04 end
n 1 0 2.3712E-03 end
o 1 0 3.7970E-02 end
h 1 0 5.7694E-02 end
fe 2 0 5.9088E-02 end
cr 2 0 1.6532E-02 end
ni 2 0 8.1369E-03 end
mn 2 0 1.3039E-03 end
si 2 0 1.3603E-03 end
ti 2 0 5.9844E-04 end
h 3 0 6.6742E-02 end
o 3 0 3.3371E-02 end
b-10 4 0 1.0844E-02 end
b-11 4 0 4.3648E-02 end
c 4 0 1.3623E-02 end
end comp
leu-sol-therm-006-03
read param
tme=360
gen=1020
npg=1000
nsk=20
fdn=yes
far=yes
plt=no
end param
read geometry
unit 11
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 3.8 -3.8 3. -3. 1.7 0.
unit 21
cylinder 4 1 2.25 1.7 0.7
cylinder 2 1 2.75 1.7 0.
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 3.8 -3.8 3. -3. 1.7 0.
unit 12
cuboid 1 1 3.8 -3.8 3. -3. 28.7180 1.7
unit 22
cylinder 4 1 2.25 28.7180 1.7
cylinder 2 1 2.75 28.7180 1.7
cuboid 1 1 3.8 -3.8 3. -3. 28.7180 1.7
unit 13
cuboid 0 1 3.8 -3.8 3. -3. 108. 28.7180
unit 23
cylinder 4 1 2.25 108. 28.7180
cylinder 2 1 2.75 108. 28.7180
cuboid 0 1 3.8 -3.8 3. -3. 108. 28.7180
unit 14
cuboid 0 1 3.8 -3.8 3. -3. 248.5 108.
unit 24
cylinder 4 1 2.25 248.5 108.
cylinder 2 1 2.75 248.5 108.
cuboid 0 1 3.8 -3.8 3. -3. 248.5 108.
' . . . . .
unit 81
array 81 0. 0. 0.
unit 91
array 91 0. 0. 0.
unit 101
```

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KENO-5A Input Listing for Case 3 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

array 101	0.	0.	0.
unit 111			
array 111	0.	0.	0.
unit 121			
array 121	0.	0.	0.
unit 131			
array 131	0.	0.	0.
unit 141			
array 141	0.	0.	0.
unit 151			
array 151	0.	0.	0.
unit 191			
array 191	0.	0.	0.
unit 82			
array 82	0.	0.	0.
unit 92			
array 92	0.	0.	0.
unit 102			
array 102	0.	0.	0.
unit 112			
array 112	0.	0.	0.
unit 122			
array 122	0.	0.	0.
unit 132			
array 132	0.	0.	0.
unit 142			
array 142	0.	0.	0.
unit 152			
array 152	0.	0.	0.
unit 192			
array 192	0.	0.	0.
unit 83			
array 83	0.	0.	0.
unit 93			
array 93	0.	0.	0.
unit 103			
array 103	0.	0.	0.
unit 113			
array 113	0.	0.	0.
unit 123			
array 123	0.	0.	0.
unit 133			
array 133	0.	0.	0.
unit 143			
array 143	0.	0.	0.
unit 153			
array 153	0.	0.	0.
unit 193			
array 193	0.	0.	0.
unit 84			
array 84	0.	0.	0.
unit 94			
array 94	0.	0.	0.
unit 104			
array 104	0.	0.	0.
unit 114			
array 114	0.	0.	0.
unit 124			
array 124	0.	0.	0.
unit 134			
array 134	0.	0.	0.
unit 144			
array 144	0.	0.	0.

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KENO-5A Input Listing for Case 3 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

unit 154
array 154  0.  0.  0.
unit 194
array 194  0.  0.  0.
' . . . . .
unit 999
cylinder 0 1 55.                248.5 108.
hole 84  -22.8 -49.07255148133 108.
hole 94  -34.2 -42.49075841257 108.
hole 104 -38.  -35.90896534381 108.
hole 114 -41.8 -29.32717227505 108.
hole 124 -45.6 -22.74537920629 108.
hole 134 -49.4 -16.16358613752 108.
hole 144 -53.2 -9.581793068762 108.
hole 154 -49.4 -3.  108.
hole 144 -53.2  3.581793068762 108.
hole 134 -49.4  10.16358613752 108.
hole 124 -45.6  16.74537920629 108.
hole 114 -41.8  23.32717227505 108.
hole 104 -38.  29.90896534381 108.
hole 194 -34.2  36.49075841257 108.
hole 84  -22.8  43.07255148133 108.
cylinder 2 1 55.6              248.5 108.
' . . . . .
global unit 1000
cylinder 2 1 54.8              1.7  0.
hole 81  -22.8 -49.07255148133  0.
hole 91  -34.2 -42.49075841257  0.
hole 101 -38.  -35.90896534381  0.
hole 111 -41.8 -29.32717227505  0.
hole 121 -45.6 -22.74537920629  0.
hole 131 -49.4 -16.16358613752  0.
hole 141 -53.2 -9.581793068762  0.
hole 151 -49.4 -3.  0.
hole 141 -53.2  3.581793068762  0.
hole 131 -49.4  10.16358613752  0.
hole 121 -45.6  16.74537920629  0.
hole 111 -41.8  23.32717227505  0.
hole 101 -38.  29.90896534381  0.
hole 191 -34.2  36.49075841257  0.
hole 81  -22.8  43.07255148133  0.
cylinder 1 1 55.                28.7180 0.
hole 82  -22.8 -49.07255148133  1.7
hole 92  -34.2 -42.49075841257  1.7
hole 102 -38.  -35.90896534381  1.7
hole 112 -41.8 -29.32717227505  1.7
hole 122 -45.6 -22.74537920629  1.7
hole 132 -49.4 -16.16358613752  1.7
hole 142 -53.2 -9.581793068762  1.7
hole 152 -49.4 -3.  1.7
hole 142 -53.2  3.581793068762  1.7
hole 132 -49.4  10.16358613752  1.7
hole 122 -45.6  16.74537920629  1.7
hole 112 -41.8  23.32717227505  1.7
hole 102 -38.  29.90896534381  1.7
hole 192 -34.2  36.49075841257  1.7
hole 82  -22.8  43.07255148133  1.7
cylinder 0 1 55.                108.  0.
hole 83  -22.8 -49.07255148133 28.7180
hole 93  -34.2 -42.49075841257 28.7180
hole 103 -38.  -35.90896534381 28.7180
hole 113 -41.8 -29.32717227505 28.7180
hole 123 -45.6 -22.74537920629 28.7180
hole 133 -49.4 -16.16358613752 28.7180

```


LEU-SOL-THERM-006

KENO-5A Input Listing for Case 3 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

hole 143 -53.2 -9.581793068762 28.7180
hole 153 -49.4 -3. 28.7180
hole 143 -53.2 3.581793068762 28.7180
hole 133 -49.4 10.16358613752 28.7180
hole 123 -45.6 16.74537920629 28.7180
hole 113 -41.8 23.32717227505 28.7180
hole 103 -38. 29.90896534381 28.7180
hole 193 -34.2 36.49075841257 28.7180
hole 83 -22.8 43.07255148133 28.7180
cylinder 2 1 55.6 108. -1.5
cylinder 3 1 99.2 108. -37.5
cylinder 0 1 99.2 248.5 -37.5
hole 999 0. 0. 0.
cylinder 2 1 100. 248.5 -38.5
end geometry
read array
ara= 81 nux= 6 nuy= 1 nuz=1 fill f11 end fill
ara= 91 nux= 9 nuy= 1 nuz=1 fill 11 21 2r11 q3 21 11 end fill
ara=101 nux=10 nuy= 1 nuz=1 fill f11 end fill
ara=111 nux=11 nuy= 1 nuz=1 fill f11 end fill
ara=121 nux=12 nuy= 1 nuz=1 fill 11 21 2r11 2q3 21 11 end fill
ara=131 nux=13 nuy= 1 nuz=1 fill f11 end fill
ara=141 nux=14 nuy= 1 nuz=1 fill f11 end fill
ara=151 nux=13 nuy= 1 nuz=1 fill 21 2r11 3q3 21 end fill
ara=191 nux= 9 nuy= 1 nuz=1 fill 11 21 5r11 21 11 end fill
ara= 82 nux= 6 nuy= 1 nuz=1 fill f12 end fill
ara= 92 nux= 9 nuy= 1 nuz=1 fill 12 22 2r12 q3 22 12 end fill
ara=102 nux=10 nuy= 1 nuz=1 fill f12 end fill
ara=112 nux=11 nuy= 1 nuz=1 fill f12 end fill
ara=122 nux=12 nuy= 1 nuz=1 fill 12 22 2r12 2q3 22 12 end fill
ara=132 nux=13 nuy= 1 nuz=1 fill f12 end fill
ara=142 nux=14 nuy= 1 nuz=1 fill f12 end fill
ara=152 nux=13 nuy= 1 nuz=1 fill 22 2r12 3q3 22 end fill
ara=192 nux= 9 nuy= 1 nuz=1 fill 12 22 5r12 22 12 end fill
ara= 83 nux= 6 nuy= 1 nuz=1 fill f13 end fill
ara= 93 nux= 9 nuy= 1 nuz=1 fill 13 23 2r13 q3 23 13 end fill
ara=103 nux=10 nuy= 1 nuz=1 fill f13 end fill
ara=113 nux=11 nuy= 1 nuz=1 fill f13 end fill
ara=123 nux=12 nuy= 1 nuz=1 fill 13 23 2r13 2q3 23 13 end fill
ara=133 nux=13 nuy= 1 nuz=1 fill f13 end fill
ara=143 nux=14 nuy= 1 nuz=1 fill f13 end fill
ara=153 nux=13 nuy= 1 nuz=1 fill 23 2r13 3q3 23 end fill
ara=193 nux= 9 nuy= 1 nuz=1 fill 13 23 5r13 23 13 end fill
ara= 84 nux= 6 nuy= 1 nuz=1 fill f14 end fill
ara= 94 nux= 9 nuy= 1 nuz=1 fill 14 24 2r14 q3 24 14 end fill
ara=104 nux=10 nuy= 1 nuz=1 fill f14 end fill
ara=114 nux=11 nuy= 1 nuz=1 fill f14 end fill
ara=124 nux=12 nuy= 1 nuz=1 fill 14 24 2r14 2q3 24 14 end fill
ara=134 nux=13 nuy= 1 nuz=1 fill f14 end fill
ara=144 nux=14 nuy= 1 nuz=1 fill f14 end fill
ara=154 nux=13 nuy= 1 nuz=1 fill 24 2r14 3q3 24 end fill
ara=194 nux= 9 nuy= 1 nuz=1 fill 14 24 5r14 24 14 end fill
end array
read plot
scr=yes
ttl='x-z slice at y=0.0 with x across and z down'
xul=-101.0 yul= 0.0 zul= 250.0
xlr= 101.0 ylr= 0.0 zlr=-40.0
uax=1.0 wdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=10.0 with x across and y down'
xul=-101.0 yul= 101.0 zul= 10.0
xlr= 101.0 ylr=-101.0 zlr= 10.0
uax=1.0 vdn=-1.0

```

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KENO-5A Input Listing for Case 3 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```
nax=800 lpi=10 end
ttl='x-y slice at z=1.0 with x across and y down'
xul= -60.0 yul= 60.0 zul= 1.0
xlr= 60.0 ylr= -60.0 zlr= 1.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
end plot
end data
end
```

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KENO-5A Input Listing for Case 4 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections).

```
=csas25
leu-sol-therm-006-04
27groupndf4 infhommedium
u-234 1 0 9.5863E-07 end
u-235 1 0 1.0854E-04 end
u-238 1 0 9.5565E-04 end
n 1 0 2.3712E-03 end
o 1 0 3.7970E-02 end
h 1 0 5.7694E-02 end
fe 2 0 5.9088E-02 end
cr 2 0 1.6532E-02 end
ni 2 0 8.1369E-03 end
mn 2 0 1.3039E-03 end
si 2 0 1.3603E-03 end
ti 2 0 5.9844E-04 end
h 3 0 6.6742E-02 end
o 3 0 3.3371E-02 end
b-10 4 0 1.0844E-02 end
b-11 4 0 4.3648E-02 end
c 4 0 1.3623E-02 end
end comp
leu-sol-therm-006-04
read param
tme=360
gen=1020
npg=1000
nsk=20
fdn=yes
far=yes
plt=no
end param
read geometry
unit 11
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 3.8 -3.8 3. -3. 1.7 0.
unit 21
cylinder 4 1 2.25 1.7 0.7
cylinder 2 1 2.75 1.7 0.
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 3.8 -3.8 3. -3. 1.7 0.
unit 12
cuboid 1 1 3.8 -3.8 3. -3. 29.2573 1.7
unit 22
cylinder 4 1 2.25 29.2573 1.7
cylinder 2 1 2.75 29.2573 1.7
cuboid 1 1 3.8 -3.8 3. -3. 29.2573 1.7
unit 13
cuboid 0 1 3.8 -3.8 3. -3. 108. 29.2573
unit 23
cylinder 4 1 2.25 108. 29.2573
cylinder 2 1 2.75 108. 29.2573
cuboid 0 1 3.8 -3.8 3. -3. 108. 29.2573
unit 14
cuboid 0 1 3.8 -3.8 3. -3. 248.5 108.
unit 24
cylinder 4 1 2.25 248.5 108.
cylinder 2 1 2.75 248.5 108.
cuboid 0 1 3.8 -3.8 3. -3. 248.5 108.
' . . . . .
unit 81
array 81 0. 0. 0.
unit 91
array 91 0. 0. 0.
unit 101
```

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KENO-5A Input Listing for Case 4 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

array 101  0.  0.  0.
unit 111
array 111  0.  0.  0.
unit 121
array 121  0.  0.  0.
unit 131
array 131  0.  0.  0.
unit 141
array 141  0.  0.  0.
unit 151
array 151  0.  0.  0.
unit 82
array 82  0.  0.  0.
unit 92
array 92  0.  0.  0.
unit 102
array 102  0.  0.  0.
unit 112
array 112  0.  0.  0.
unit 122
array 122  0.  0.  0.
unit 132
array 132  0.  0.  0.
unit 142
array 142  0.  0.  0.
unit 152
array 152  0.  0.  0.
unit 83
array 83  0.  0.  0.
unit 93
array 93  0.  0.  0.
unit 103
array 103  0.  0.  0.
unit 113
array 113  0.  0.  0.
unit 123
array 123  0.  0.  0.
unit 133
array 133  0.  0.  0.
unit 143
array 143  0.  0.  0.
unit 153
array 153  0.  0.  0.
unit 84
array 84  0.  0.  0.
unit 94
array 94  0.  0.  0.
unit 104
array 104  0.  0.  0.
unit 114
array 114  0.  0.  0.
unit 124
array 124  0.  0.  0.
unit 134
array 134  0.  0.  0.
unit 144
array 144  0.  0.  0.
unit 154
array 154  0.  0.  0.
'
unit 999
cylinder 0 1 55.          248.5 108.
hole 84  -22.8 -49.07255148133 108.
hole 94  -34.2 -42.49075841257 108.

```

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KENO-5A Input Listing for Case 4 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

hole 104 -38. -35.90896534381 108.
hole 114 -41.8 -29.32717227505 108.
hole 124 -45.6 -22.74537920629 108.
hole 134 -49.4 -16.16358613752 108.
hole 144 -53.2 -9.581793068762 108.
hole 154 -49.4 -3. 108.
hole 144 -53.2 3.581793068762 108.
hole 134 -49.4 10.16358613752 108.
hole 124 -45.6 16.74537920629 108.
hole 114 -41.8 23.32717227505 108.
hole 104 -38. 29.90896534381 108.
hole 94 -34.2 36.49075841257 108.
hole 84 -22.8 43.07255148133 108.
cylinder 2 1 55.6 248.5 108.

global unit 1000
cylinder 2 1 54.8 1.7 0.
hole 81 -22.8 -49.07255148133 0.
hole 91 -34.2 -42.49075841257 0.
hole 101 -38. -35.90896534381 0.
hole 111 -41.8 -29.32717227505 0.
hole 121 -45.6 -22.74537920629 0.
hole 131 -49.4 -16.16358613752 0.
hole 141 -53.2 -9.581793068762 0.
hole 151 -49.4 -3. 0.
hole 141 -53.2 3.581793068762 0.
hole 131 -49.4 10.16358613752 0.
hole 121 -45.6 16.74537920629 0.
hole 111 -41.8 23.32717227505 0.
hole 101 -38. 29.90896534381 0.
hole 91 -34.2 36.49075841257 0.
hole 81 -22.8 43.07255148133 0.
cylinder 1 1 55. 29.2573 0.
hole 82 -22.8 -49.07255148133 1.7
hole 92 -34.2 -42.49075841257 1.7
hole 102 -38. -35.90896534381 1.7
hole 112 -41.8 -29.32717227505 1.7
hole 122 -45.6 -22.74537920629 1.7
hole 132 -49.4 -16.16358613752 1.7
hole 142 -53.2 -9.581793068762 1.7
hole 152 -49.4 -3. 1.7
hole 142 -53.2 3.581793068762 1.7
hole 132 -49.4 10.16358613752 1.7
hole 122 -45.6 16.74537920629 1.7
hole 112 -41.8 23.32717227505 1.7
hole 102 -38. 29.90896534381 1.7
hole 92 -34.2 36.49075841257 1.7
hole 82 -22.8 43.07255148133 1.7
cylinder 0 1 55. 108. 0.
hole 83 -22.8 -49.07255148133 29.2573
hole 93 -34.2 -42.49075841257 29.2573
hole 103 -38. -35.90896534381 29.2573
hole 113 -41.8 -29.32717227505 29.2573
hole 123 -45.6 -22.74537920629 29.2573
hole 133 -49.4 -16.16358613752 29.2573
hole 143 -53.2 -9.581793068762 29.2573
hole 153 -49.4 -3. 29.2573
hole 143 -53.2 3.581793068762 29.2573
hole 133 -49.4 10.16358613752 29.2573
hole 123 -45.6 16.74537920629 29.2573
hole 113 -41.8 23.32717227505 29.2573
hole 103 -38. 29.90896534381 29.2573
hole 93 -34.2 36.49075841257 29.2573
hole 83 -22.8 43.07255148133 29.2573

```

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KENO-5A Input Listing for Case 4 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

cylinder 2 1 55.6          108.  -1.5
cylinder 3 1 99.2          108.  -37.5
cylinder 0 1 99.2          248.5  -37.5
hole 999  0.  0.  0.
cylinder 2 1 100.          248.5  -38.5
end geometry
read array
ara= 81  nux= 6 nuy= 1 nuz=1  fill f11 end fill
ara= 91  nux= 9 nuy= 1 nuz=1  fill 11 21 2r11 q3 21 11 end fill
ara=101  nux=10 nuy= 1 nuz=1  fill f11 end fill
ara=111  nux=11 nuy= 1 nuz=1  fill f11 end fill
ara=121  nux=12 nuy= 1 nuz=1  fill 11 21 2r11 2q3 21 11 end fill
ara=131  nux=13 nuy= 1 nuz=1  fill f11 end fill
ara=141  nux=14 nuy= 1 nuz=1  fill f11 end fill
ara=151  nux=13 nuy= 1 nuz=1  fill 21 2r11 3q3 21 end fill
ara= 82  nux= 6 nuy= 1 nuz=1  fill f12 end fill
ara= 92  nux= 9 nuy= 1 nuz=1  fill 12 22 2r12 q3 22 12 end fill
ara=102  nux=10 nuy= 1 nuz=1  fill f12 end fill
ara=112  nux=11 nuy= 1 nuz=1  fill f12 end fill
ara=122  nux=12 nuy= 1 nuz=1  fill 12 22 2r12 2q3 22 12 end fill
ara=132  nux=13 nuy= 1 nuz=1  fill f12 end fill
ara=142  nux=14 nuy= 1 nuz=1  fill f12 end fill
ara=152  nux=13 nuy= 1 nuz=1  fill 22 2r12 3q3 22 end fill
ara= 83  nux= 6 nuy= 1 nuz=1  fill f13 end fill
ara= 93  nux= 9 nuy= 1 nuz=1  fill 13 23 2r13 q3 23 13 end fill
ara=103  nux=10 nuy= 1 nuz=1  fill f13 end fill
ara=113  nux=11 nuy= 1 nuz=1  fill f13 end fill
ara=123  nux=12 nuy= 1 nuz=1  fill 13 23 2r13 2q3 23 13 end fill
ara=133  nux=13 nuy= 1 nuz=1  fill f13 end fill
ara=143  nux=14 nuy= 1 nuz=1  fill f13 end fill
ara=153  nux=13 nuy= 1 nuz=1  fill 23 2r13 3q3 23 end fill
ara= 84  nux= 6 nuy= 1 nuz=1  fill f14 end fill
ara= 94  nux= 9 nuy= 1 nuz=1  fill 14 24 2r14 q3 24 14 end fill
ara=104  nux=10 nuy= 1 nuz=1  fill f14 end fill
ara=114  nux=11 nuy= 1 nuz=1  fill f14 end fill
ara=124  nux=12 nuy= 1 nuz=1  fill 14 24 2r14 2q3 24 14 end fill
ara=134  nux=13 nuy= 1 nuz=1  fill f14 end fill
ara=144  nux=14 nuy= 1 nuz=1  fill f14 end fill
ara=154  nux=13 nuy= 1 nuz=1  fill 24 2r14 3q3 24 end fill
end array
read plot
scr=yes
ttl='x-z slice at y=0.0 with x across and z down'
xul=-101.0 yul=  0.0  zul= 250.0
xlr= 101.0 ylr=  0.0  zlr= -40.0
uax=1.0 wdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=10.0 with x across and y down'
xul=-101.0 yul= 101.0  zul=  10.0
xlr= 101.0 ylr=-101.0  zlr=  10.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=1.0 with x across and y down'
xul= -60.0 yul=  60.0  zul=  1.0
xlr=  60.0 ylr= -60.0  zlr=  1.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
end plot
end data
end

```

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KENO-5A Input Listing for Case 5 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections).

```
=csas25
leu-sol-therm-006-05
27groupndf4 infhommedium
u-234 1 0 9.5863E-07 end
u-235 1 0 1.0854E-04 end
u-238 1 0 9.5565E-04 end
n 1 0 2.3712E-03 end
o 1 0 3.7970E-02 end
h 1 0 5.7694E-02 end
fe 2 0 5.9088E-02 end
cr 2 0 1.6532E-02 end
ni 2 0 8.1369E-03 end
mn 2 0 1.3039E-03 end
si 2 0 1.3603E-03 end
ti 2 0 5.9844E-04 end
h 3 0 6.6742E-02 end
o 3 0 3.3371E-02 end
b-10 4 0 1.0844E-02 end
b-11 4 0 4.3648E-02 end
c 4 0 1.3623E-02 end
end comp
leu-sol-therm-006-05
read param
tme=360
gen=1020
npg=1000
nsk=20
fdn=yes
far=yes
plt=no
end param
read geometry
unit 11
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 5.3 -5.3 3. -3. 1.7 0.
unit 21
cylinder 4 1 2.25 1.7 0.7
cylinder 2 1 2.75 1.7 0.
cylinder 1 1 2.775 1.7 0.
cuboid 2 1 5.3 -5.3 3. -3. 1.7 0.
unit 12
cuboid 1 1 5.3 -5.3 3. -3. 36.3693 1.7
unit 22
cylinder 4 1 2.25 36.3693 1.7
cylinder 2 1 2.75 36.3693 1.7
cuboid 1 1 5.3 -5.3 3. -3. 36.3693 1.7
unit 13
cuboid 0 1 5.3 -5.3 3. -3. 108. 36.3693
unit 23
cylinder 4 1 2.25 108. 36.3693
cylinder 2 1 2.75 108. 36.3693
cuboid 0 1 5.3 -5.3 3. -3. 108. 36.3693
unit 14
cuboid 0 1 5.3 -5.3 3. -3. 248.5 108.
unit 24
cylinder 4 1 2.25 248.5 108.
cylinder 2 1 2.75 248.5 108.
cuboid 0 1 5.3 -5.3 3. -3. 248.5 108.
' . . . . .
unit 61
array 61 0. 0. 0.
unit 71
array 71 0. 0. 0.
```

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KENO-5A Input Listing for Case 5 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

unit 81
array 81 0. 0. 0.
unit 91
array 91 0. 0. 0.
unit 101
array 101 0. 0. 0.
unit 62
array 62 0. 0. 0.
unit 72
array 72 0. 0. 0.
unit 82
array 82 0. 0. 0.
unit 92
array 92 0. 0. 0.
unit 102
array 102 0. 0. 0.
unit 63
array 63 0. 0. 0.
unit 73
array 73 0. 0. 0.
unit 83
array 83 0. 0. 0.
unit 93
array 93 0. 0. 0.
unit 103
array 103 0. 0. 0.
unit 64
array 64 0. 0. 0.
unit 74
array 74 0. 0. 0.
unit 84
array 84 0. 0. 0.
unit 94
array 94 0. 0. 0.
unit 104
array 104 0. 0. 0.
' . . . . .
unit 999
cylinder 0 1 55. 248.5 108.
hole 64 -21.2 -49.89934640058 108.
hole 74 -37.1 -39.71947712046 108.
hole 84 -42.4 -30.53960784035 108.
hole 94 -47.7 -21.35973856023 108.
hole 104 -53. -12.17986928012 108.
hole 94 -47.7 -3. 108.
hole 104 -53.0 6.179869280115 108.
hole 94 -47.7 15.35973856023 108.
hole 84 -42.4 24.53960784035 108.
hole 74 -37.1 33.71947712046 108.
hole 64 -21.2 42.89934640058 108.
cylinder 2 1 55.6 248.5 108.
' . . . . .
global unit 1000
cylinder 2 1 54.8 1.7 0.
hole 61 -21.2 -49.89934640058 0.
hole 71 -37.1 -39.71947712046 0.
hole 81 -42.4 -30.53960784035 0.
hole 91 -47.7 -21.35973856023 0.
hole 101 -53. -12.17986928012 0.
hole 91 -47.7 -3. 0.
hole 101 -53.0 6.179869280115 0.
hole 91 -47.7 15.35973856023 0.
hole 81 -42.4 24.53960784035 0.
hole 71 -37.1 33.71947712046 0.

```


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KENO-5A Input Listing for Case 5 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```

hole 61 -21.2 42.89934640058 0.
cylinder 1 1 55. 36.3693 0.
hole 62 -21.2 -49.89934640058 1.7
hole 72 -37.1 -39.71947712046 1.7
hole 82 -42.4 -30.53960784035 1.7
hole 92 -47.7 -21.35973856023 1.7
hole 102 -53. -12.17986928012 1.7
hole 92 -47.7 -3. 1.7
hole 102 -53.0 6.179869280115 1.7
hole 92 -47.7 15.35973856023 1.7
hole 82 -42.4 24.53960784035 1.7
hole 72 -37.1 33.71947712046 1.7
hole 62 -21.2 42.89934640058 1.7
cylinder 0 1 55. 108. 0.
hole 63 -21.2 -49.89934640058 36.3693
hole 73 -37.1 -39.71947712046 36.3693
hole 83 -42.4 -30.53960784035 36.3693
hole 93 -47.7 -21.35973856023 36.3693
hole 103 -53. -12.17986928012 36.3693
hole 93 -47.7 -3. 36.3693
hole 103 -53.0 6.179869280115 36.3693
hole 93 -47.7 15.35973856023 36.3693
hole 83 -42.4 24.53960784035 36.3693
hole 73 -37.1 33.71947712046 36.3693
hole 63 -21.2 42.89934640058 36.3693
cylinder 2 1 55.6 108. -1.5
cylinder 3 1 99.2 108. -37.5
cylinder 0 1 99.2 248.5 -37.5
hole 999 0. 0. 0.
cylinder 2 1 100. 248.5 -38.5
end geometry
read array
ara= 61 nux= 4 nuy= 1 nuz=1 fill 21 11 n2 end fill
ara= 71 nux= 7 nuy= 1 nuz=1 fill 21 2r11 21 1b3 end fill
ara= 81 nux= 8 nuy= 1 nuz=1 fill 2r11 21 11 n4 end fill
ara= 91 nux= 9 nuy= 1 nuz=1 fill 11 21 2r11 21 2r11 21 11 end fill
ara=101 nux=10 nuy= 1 nuz=1 fill 21 2r11 21 2r11 21 2r11 21 end fill
ara= 62 nux= 4 nuy= 1 nuz=1 fill 22 12 n2 end fill
ara= 72 nux= 7 nuy= 1 nuz=1 fill 22 2r12 22 1b3 end fill
ara= 82 nux= 8 nuy= 1 nuz=1 fill 2r12 22 12 n4 end fill
ara= 92 nux= 9 nuy= 1 nuz=1 fill 12 22 2r12 22 2r12 22 12 end fill
ara=102 nux=10 nuy= 1 nuz=1 fill 22 2r12 22 2r12 22 2r12 22 end fill
ara= 63 nux= 4 nuy= 1 nuz=1 fill 23 13 n2 end fill
ara= 73 nux= 7 nuy= 1 nuz=1 fill 23 2r13 23 1b3 end fill
ara= 83 nux= 8 nuy= 1 nuz=1 fill 2r13 23 13 n4 end fill
ara= 93 nux= 9 nuy= 1 nuz=1 fill 13 23 2r13 23 2r13 23 13 end fill
ara=103 nux=10 nuy= 1 nuz=1 fill 23 2r13 23 2r13 23 2r13 23 end fill
ara= 64 nux= 4 nuy= 1 nuz=1 fill 24 14 n2 end fill
ara= 74 nux= 7 nuy= 1 nuz=1 fill 24 2r14 24 1b3 end fill
ara= 84 nux= 8 nuy= 1 nuz=1 fill 2r14 24 14 n4 end fill
ara= 94 nux= 9 nuy= 1 nuz=1 fill 14 24 2r14 24 2r14 24 14 end fill
ara=104 nux=10 nuy= 1 nuz=1 fill 24 2r14 24 2r14 24 2r14 24 end fill
end array
read plot
scr=yes
ttl='x-z slice at y=0.0 with x across and z down'
xul=-101.0 yul= 0.0 zul= 250.0
xlr= 101.0 ylr= 0.0 zlr= -40.0
uax=1.0 wdn=-1.0
nax=800 lpi=10 end
ttl='x-y slice at z=10.0 with x across and y down'
xul=-101.0 yul= 101.0 zul= 10.0
xlr= 101.0 ylr=-101.0 zlr= 10.0
uax=1.0 vdn=-1.0

```

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KENO-5A Input Listing for Case 5 of Table 13.b (27-Energy-Group ENDF/B-IV Cross Sections) (cont'd).

```
nax=800 lpi=10 end
ttl='x-y slice at z=1.0 with x across and y down'
xul= -60.0 yul= 60.0 zul= 1.0
xlr= 60.0 ylr= -60.0 zlr= 1.0
uax=1.0 vdn=-1.0
nax=800 lpi=10 end
end plot
end data
end
```

A.2 MCNP Input Listings

The MCNP4A calculations with continuous-energy ENDF/B-V cross sections were run using 1020 generations of neutrons with 1000 neutrons each. Twenty generations were skipped before averaging, so the result is the average of 1,000,000 neutron histories.

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MCNP Input Listing for Case 1 of Table 13.b.

leu-sol-therm-006-01

```

1 0      4:-5:10      imp:n=0
2 3 8.701954E-02 -4 5 -10 (3:-6)  imp:n=1
3 2 1.001130E-01 -3 6 -9 (2:-7)  imp:n=1
4 0      -3 2 9 -10      imp:n=1
5 3 8.701954E-02 -2 7 -10 (1:-8)  imp:n=1
6 1 9.910035E-02 -1 11 8 -201  imp:n=1
7 0      -1 11 201 -10      imp:n=1
8 0      -11 8 -10      imp:n=1 fill=100
c
100 0 -401 402 -403 404 imp:n=1 lat=1 u=100
    fill=-6:6 -6:6 0:0
        1 18r
            1 2 2 2 2 1      1 5r
            2 2 2 2 2 2      1 4r
            2 2 2 2 2 2 2      1 3r
            2 2 2 2 2 2 2 2      1 2r
            2 2 2 2 2 2 2 2 2      1 1r
            1 2 2 2 2 2 2 2 2 1      1 1r
            2 2 2 2 2 2 2 2 2      1 2r
            2 2 2 2 2 2 2 2      1 3r
            2 2 2 2 2 2 2      1 4r
            2 2 2 2 2 2      1 5r
            1 2 2 2 2 1      1 18r
c
1000 3 8.701954E-02      -200 imp:n=1 u=1
1001 1 9.910035E-02      200 -201 imp:n=1 u=1
1002 0      201      imp:n=1 u=1
c
2000 1 9.910035E-02      -300 -200 imp:n=1 u=2
2001 3 8.701954E-02 300      -200 imp:n=1 u=2
2003 1 9.910035E-02      200 -201 imp:n=1 u=2
2004 0      201      imp:n=1 u=2
c
3000 3 8.701954E-02      -302 -303 imp:n=1 u=3
3001 4 6.811600E-02      -302 303      imp:n=1 u=3
3002 3 8.701954E-02 302 -301      imp:n=1 u=3
3003 1 9.910035E-02 301 -300      -200 imp:n=1 u=3
3004 3 8.701954E-02 300      -200 imp:n=1 u=3
3005 1 9.910035E-02 301      200 -201 imp:n=1 u=3
3006 0      301      201      imp:n=1 u=3

1 cz 55.
2 cz 55.6
3 cz 99.2
4 cz 100.
c
5 pz -38.5
6 pz -37.5
7 pz -1.5
8 pz 0.
9 pz 108.    $ water level
10 pz 248.5
c
11 cz 54.8
c
200 pz 1.7
201 pz 23.4174    $ solution level
c
300 cz 2.775
301 cz 2.75
302 cz 2.25
303 pz 0.7
c

```

LEU-SOL-THERM-006

MCNP Input Listing for Case 1 of Table 13.b (cont'd).

```
401 p 1.732050807569 -1 0 9.179869280115
402 p 1.732050807569 -1 0 -9.179869280115
403 py 4.589934640058
404 py -4.589934640058

m1 92234.50c 9.5863E-07
    92235.50c 1.0854E-04
    92238.50c 9.5565E-04
    7014.50c 2.3712E-03
    8016.50c 3.7970E-02
    1001.50c 5.7694E-02
m2 8016.50c 3.3371E-02
    1001.50c 6.6742E-02
m3 26000.50c 5.9088E-02
    24000.50c 1.6532E-02
    28000.50c 8.1369E-03
    25055.50c 1.3039E-03
    14000.50c 1.3603E-03
    22000.50c 5.9844E-04
m4 5010.50c 1.0844E-02
    5011.50c 4.3649E-02
    6012.50c 1.3623E-02
mt1 lwtr.01t
mt2 lwtr.01t
kcode 1000 1.0 20 1020
sdef pos=0. 0. 11.8 axs=0 0 1 rad=d1 ext=d2
si1 0. 55.
si2 11.8
prdmp 3j 1
print
```

LEU-SOL-THERM-006

MCNP Input Listing for Case 2 of Table 13.b.

```

leu-sol-therm-006-02
 1 0      4:-5:10      imp:n=0
 2 3 8.701954E-02 -4 5 -10 (3:-6)  imp:n=1
 3 2 1.001130E-01 -3 6 -9 (2:-7)  imp:n=1
 4 0      -3 2 9 -10      imp:n=1
 5 3 8.701954E-02 -2 7 -10 (1:-8)  imp:n=1
 6 1 9.910035E-02 -1 11 8 -201  imp:n=1
 7 0      -1 11 201 -10      imp:n=1
 8 0      -11 8 -10      imp:n=1 fill=100
c
100 0 -401 402 -403 404 imp:n=1 lat=1 u=100
    fill=-6:6 -6:6 0:0
          1 18r
          1 2 2 2 2 1      1 5r
          2 2 2 2 2 2      1 4r
          2 2 3 2 2 3 2 2      1 3r
          2 2 2 2 2 2 2 2      1 2r
          2 2 2 2 2 2 2 2 2      1 1r
          1 2 3 2 2 3 2 2 3 2 1      1 1r
          2 2 2 2 2 2 2 2 2      1 2r
          2 2 2 2 2 2 2 2 2      1 3r
          2 2 3 2 2 3 2 2      1 4r
          2 2 2 2 2 2 2      1 5r
          1 2 2 2 2 1      1 18r
c
1000 3 8.701954E-02      -200 imp:n=1 u=1
1001 1 9.910035E-02      200 -201 imp:n=1 u=1
1002 0      201      imp:n=1 u=1
c
2000 1 9.910035E-02      -300 -200 imp:n=1 u=2
2001 3 8.701954E-02 300      -200 imp:n=1 u=2
2003 1 9.910035E-02      200 -201 imp:n=1 u=2
2004 0      201      imp:n=1 u=2
c
3000 3 8.701954E-02      -302 -303 imp:n=1 u=3
3001 4 6.811600E-02      -302 303      imp:n=1 u=3
3002 3 8.701954E-02 302 -301      imp:n=1 u=3
3003 1 9.910035E-02 301 -300      -200 imp:n=1 u=3
3004 3 8.701954E-02 300      -200 imp:n=1 u=3
3005 1 9.910035E-02 301      200 -201 imp:n=1 u=3
3006 0      301      201      imp:n=1 u=3

 1 cz 55.
 2 cz 55.6
 3 cz 99.2
 4 cz 100.
c
 5 pz -38.5
 6 pz -37.5
 7 pz -1.5
 8 pz 0.
 9 pz 108.    $ water level
10 pz 248.5
c
11 cz 54.8
c
200 pz 1.7
201 pz 26.1371    $ solution level
c
300 cz 2.775
301 cz 2.75
302 cz 2.25
303 pz 0.7
c

```

LEU-SOL-THERM-006

MCNP Input Listing for Case 2 of Table 13.b (cont'd).

```
401 p 1.732050807569 -1 0 9.179869280115
402 p 1.732050807569 -1 0 -9.179869280115
403 py 4.589934640058
404 py -4.589934640058

m1 92234.50c 9.5863E-07
    92235.50c 1.0854E-04
    92238.50c 9.5565E-04
    7014.50c 2.3712E-03
    8016.50c 3.7970E-02
    1001.50c 5.7694E-02
m2 8016.50c 3.3371E-02
    1001.50c 6.6742E-02
m3 26000.50c 5.9088E-02
    24000.50c 1.6532E-02
    28000.50c 8.1369E-03
    25055.50c 1.3039E-03
    14000.50c 1.3603E-03
    22000.50c 5.9844E-04
m4 5010.50c 1.0844E-02
    5011.50c 4.3649E-02
    6012.50c 1.3623E-02
mt1 lwtr.01t
mt2 lwtr.01t
kcode 1000 1.0 20 1020
sdef pos=0. 0. 13.1 axs=0 0 1 rad=d1 ext=d2
si1 0. 55.
si2 13.1
prdmp 3j 1
print
```

LEU-SOL-THERM-006

MCNP Input Listing for Case 3 of Table 13.b.

leu-sol-therm-006-03

```

1 0      4:-5:10      imp:n=0
2 3 8.701954E-02 -4 5 -10 (3:-6)  imp:n=1
3 2 1.001130E-01 -3 6 -9 (2:-7)  imp:n=1
4 0      -3 2 9 -10      imp:n=1
5 3 8.701954E-02 -2 7 -10 (1:-8)  imp:n=1
6 1 9.910035E-02 -1 11 8 -201  imp:n=1
7 0      -1 11 201 -10      imp:n=1
8 0      -11 8 -10      imp:n=1 fill=100
c
100 0 -401 402 -403 404 imp:n=1 lat=1 u=100
    fill=-8:8 -8:8 0:0
        1 24r
            1 2 2 2 2 2 2 1      1 7r
            2 3 2 2 3 2 2 3 2      1 6r
            2 2 2 2 2 2 2 2 2      1 5r
            2 2 2 2 2 2 2 2 2 2      1 4r
            2 3 2 2 3 2 2 3 2 2 3 2      1 3r
            2 2 2 2 2 2 2 2 2 2 2 2      1 2r
            2 2 2 2 2 2 2 2 2 2 2 2      1 1r
            1 3 2 2 3 2 2 3 2 2 3 2 1      1 1r
            2 2 2 2 2 2 2 2 2 2 2 2      1 2r
            2 2 2 2 2 2 2 2 2 2 2 2      1 3r
            2 3 2 2 3 2 2 3 2 2 3 2      1 4r
            2 2 2 2 2 2 2 2 2 2      1 5r
            2 2 2 2 2 2 2 2 2 2      1 6r
            2 3 2 2 2 2 3 2      1 7r
            1 2 2 2 2 2 1      1 24r
c
1000 3 8.701954E-02      -200 imp:n=1 u=1
1001 1 9.910035E-02      200 -201 imp:n=1 u=1
1002 0      201      imp:n=1 u=1
c
2000 1 9.910035E-02      -300      -200 imp:n=1 u=2
2001 3 8.701954E-02      300      -200 imp:n=1 u=2
2003 1 9.910035E-02      200 -201 imp:n=1 u=2
2004 0      201      imp:n=1 u=2
c
3000 3 8.701954E-02      -302      -303 imp:n=1 u=3
3001 4 6.811600E-02      -302 303      imp:n=1 u=3
3002 3 8.701954E-02      302 -301      imp:n=1 u=3
3003 1 9.910035E-02      301 -300      -200 imp:n=1 u=3
3004 3 8.701954E-02      300      -200 imp:n=1 u=3
3005 1 9.910035E-02      301      200 -201 imp:n=1 u=3
3006 0      301      201      imp:n=1 u=3

1 cz 55.
2 cz 55.6
3 cz 99.2
4 cz 100.
c
5 pz -38.5
6 pz -37.5
7 pz -1.5
8 pz 0.
9 pz 108. $ water level
10 pz 248.5
c
11 cz 54.8
c
200 pz 1.7
201 pz 28.7180 $ solution level
c
300 cz 2.775

```


LEU-SOL-THERM-006

MCNP Input Listing for Case 3 of Table 13.b (cont'd).

```
301 cz 2.75
302 cz 2.25
303 pz 0.7
c
401 p 1.732050807569 -1 0 6.581793068762
402 p 1.732050807569 -1 0 -6.581793068762
403 py 3.290896534381
404 py -3.290896534381

m1 92234.50c 9.5863E-07
    92235.50c 1.0854E-04
    92238.50c 9.5565E-04
    7014.50c 2.3712E-03
    8016.50c 3.7970E-02
    1001.50c 5.7694E-02
m2 8016.50c 3.3371E-02
    1001.50c 6.6742E-02
m3 26000.50c 5.9088E-02
    24000.50c 1.6532E-02
    28000.50c 8.1369E-03
    25055.50c 1.3039E-03
    14000.50c 1.3603E-03
    22000.50c 5.9844E-04
m4 5010.50c 1.0844E-02
    5011.50c 4.3649E-02
    6012.50c 1.3623E-02
mt1 lwtr.01t
mt2 lwtr.01t
kcode 1000 1.0 20 1020
sdef pos=0, 0, 14.4 axs=0 0 1 rad=d1 ext=d2
si1 0.55.
si2 14.4
prdmp 3j 1
print
```

LEU-SOL-THERM-006

MCNP Input Listing for Case 4 of Table 13.b.

leu-sol-therm-006-04

```

1 0      4:-5:10      imp:n=0
2 3 8.701954E-02 -4 5 -10 (3:-6)  imp:n=1
3 2 1.001130E-01 -3 6 -9 (2:-7)  imp:n=1
4 0      -3 2 9 -10      imp:n=1
5 3 8.701954E-02 -2 7 -10 (1:-8)  imp:n=1
6 1 9.910035E-02 -1 11 8 -201  imp:n=1
7 0      -1 11 201 -10  imp:n=1
8 0      -11 8 -10      imp:n=1 fill=100
c
100 0 -401 402 -403 404 imp:n=1 lat=1 u=100
    fill=-8:8 -8:8 0:0
        1 24r
        1 2 2 2 2 2 2 1      1 7r
        2 3 2 2 3 2 2 3 2      1 6r
        2 2 2 2 2 2 2 2 2      1 5r
        2 2 2 2 2 2 2 2 2 2      1 4r
        2 3 2 2 3 2 2 3 2 2 3 2      1 3r
        2 2 2 2 2 2 2 2 2 2 2 2 2 2      1 2r
        2 2 2 2 2 2 2 2 2 2 2 2 2 2      1 1r
        1 3 2 2 3 2 2 3 2 2 3 2 2 3 1      1 1r
        2 2 2 2 2 2 2 2 2 2 2 2 2 2      1 2r
        2 2 2 2 2 2 2 2 2 2 2 2 2 2      1 3r
        2 3 2 2 3 2 2 3 2 2 3 2      1 4r
        2 2 2 2 2 2 2 2 2 2 2 2      1 5r
        2 2 2 2 2 2 2 2 2 2      1 6r
        2 3 2 2 3 2 2 3 2      1 7r
        1 2 2 2 2 2 2 1      1 24r
c
1000 3 8.701954E-02      -200 imp:n=1 u=1
1001 1 9.910035E-02      200 -201 imp:n=1 u=1
1002 0      201      imp:n=1 u=1
c
2000 1 9.910035E-02      -300      -200 imp:n=1 u=2
2001 3 8.701954E-02      300      -200 imp:n=1 u=2
2003 1 9.910035E-02      200 -201 imp:n=1 u=2
2004 0      201      imp:n=1 u=2
c
3000 3 8.701954E-02      -302      -303 imp:n=1 u=3
3001 4 6.811600E-02      -302 303      imp:n=1 u=3
3002 3 8.701954E-02      302 -301      imp:n=1 u=3
3003 1 9.910035E-02      301 -300      -200 imp:n=1 u=3
3004 3 8.701954E-02      300      -200 imp:n=1 u=3
3005 1 9.910035E-02      301      200 -201 imp:n=1 u=3
3006 0      301      201      imp:n=1 u=3

1 cz 55.
2 cz 55.6
3 cz 99.2
4 cz 100.
c
5 pz -38.5
6 pz -37.5
7 pz -1.5
8 pz 0.
9 pz 108. $ water level
10 pz 248.5
c
11 cz 54.8
c
200 pz 1.7
201 pz 29.2573 $ solution level
c
300 cz 2.775

```

LEU-SOL-THERM-006

MCNP Input Listing for Case 4 of Table 13.b (cont'd).

```
301 cz 2.75
302 cz 2.25
303 pz 0.7
c
401 p 1.732050807569 -1 0 6.581793068762
402 p 1.732050807569 -1 0 -6.581793068762
403 py 3.290896534381
404 py -3.290896534381

m1 92234.50c 9.5863E-07
    92235.50c 1.0854E-04
    92238.50c 9.5565E-04
    7014.50c 2.3712E-03
    8016.50c 3.7970E-02
    1001.50c 5.7694E-02
m2 8016.50c 3.3371E-02
    1001.50c 6.6742E-02
m3 26000.50c 5.9088E-02
    24000.50c 1.6532E-02
    28000.50c 8.1369E-03
    25055.50c 1.3039E-03
    14000.50c 1.3603E-03
    22000.50c 5.9844E-04
m4 5010.50c 1.0844E-02
    5011.50c 4.3649E-02
    6012.50c 1.3623E-02
mt1 lwtr.01t
mt2 lwtr.01t
kcode 1000 1.0 20 1020
sdef pos=0, 0, 14.7 axs=0 0 1 rad=d1 ext=d2
si1 0.55.
si2 14.7
prdmp 3j 1
print
```

LEU-SOL-THERM-006

MCNP Input Listing for Case 5 of Table 13.b.

leu-sol-therm-006-05

```

1 0      4:-5:10      imp:n=0
2 3 8.701954E-02 -4 5 -10 (3:-6)  imp:n=1
3 2 1.001130E-01 -3 6 -9 (2:-7)  imp:n=1
4 0      -3 2 9 -10      imp:n=1
5 3 8.701954E-02 -2 7 -10 (1:-8)  imp:n=1
6 1 9.910035E-02 -1 11 8 -201  imp:n=1
7 0      -1 11 201 -10      imp:n=1
8 0      -11 8 -10      imp:n=1 fill=100
c
100 0 -401 402 -403 404 imp:n=1 lat=1 u=100
    fill=-6:6 -6:6 0:0
        1 18r
            1 3 2 2 3 1      1 5r
            3 2 2 3 2 2 3    1 4r
            2 2 3 2 2 3 2 2    1 3r
            2 3 2 2 3 2 2 3 2    1 2r
            3 2 2 3 2 2 3 2 2 3    1 1r
            1 2 3 2 2 3 2 2 3 2 1    1 1r
            3 2 2 3 2 2 3 2 2 3    1 2r
            2 3 2 2 3 2 2 3 2    1 3r
            2 2 3 2 2 3 2 2    1 4r
            3 2 2 3 2 2 3    1 5r
            1 3 2 2 3 1      1 18r
c
1000 3 8.701954E-02      -200 imp:n=1 u=1
1001 1 9.910035E-02      200 -201 imp:n=1 u=1
1002 0      201      imp:n=1 u=1
c
2000 1 9.910035E-02      -300 -200 imp:n=1 u=2
2001 3 8.701954E-02 300      -200 imp:n=1 u=2
2003 1 9.910035E-02      200 -201 imp:n=1 u=2
2004 0      201      imp:n=1 u=2
c
3000 3 8.701954E-02      -302 -303 imp:n=1 u=3
3001 4 6.811600E-02      -302 303      imp:n=1 u=3
3002 3 8.701954E-02 302 -301      imp:n=1 u=3
3003 1 9.910035E-02 301 -300      -200 imp:n=1 u=3
3004 3 8.701954E-02 300      -200 imp:n=1 u=3
3005 1 9.910035E-02 301      200 -201 imp:n=1 u=3
3006 0      301      201      imp:n=1 u=3

1 cz 55.
2 cz 55.6
3 cz 99.2
4 cz 100.
c
5 pz -38.5
6 pz -37.5
7 pz -1.5
8 pz 0.
9 pz 108.    $ water level
10 pz 248.5
c
11 cz 54.8
c
200 pz 1.7
201 pz 36.3693    $ solution level
c
300 cz 2.775
301 cz 2.75
302 cz 2.25
303 pz 0.7
c

```

LEU-SOL-THERM-006

MCNP Input Listing for Case 5 of Table 13.b (cont'd).

```
401 p 1.732050807569 -1 0 9.179869280115
402 p 1.732050807569 -1 0 -9.179869280115
403 py 4.589934640058
404 py -4.589934640058

m1 92234.50c 9.5863E-07
    92235.50c 1.0854E-04
    92238.50c 9.5565E-04
    7014.50c 2.3712E-03
    8016.50c 3.7970E-02
    1001.50c 5.7694E-02
m2 8016.50c 3.3371E-02
    1001.50c 6.6742E-02
m3 26000.50c 5.9088E-02
    24000.50c 1.6532E-02
    28000.50c 8.1369E-03
    25055.50c 1.3039E-03
    14000.50c 1.3603E-03
    22000.50c 5.9844E-04
m4 5010.50c 1.0844E-02
    5011.50c 4.3649E-02
    6012.50c 1.3623E-02
mt1 lwtr.01t
mt2 lwtr.01t
kcode 1000 1.0 20 1020
sdef pos=0. 0. 18.2 axs=0 0 1 rad=d1 ext=d2
si1 0. 55.
si2 18.2
prdmp 3j 1
print
```