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Design Analysis Cover Sheet

Complete only applicable items.

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8. Checker	David Henderson	<i>[Signature]</i>	9-2-97
9. Lead Design Engineer	Dan Thomas	<i>[Signature]</i>	09/02/97
10. Department Manager	Hugh Benton	<i>[Signature]</i>	9/2/97

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1. Purpose

The purpose of this design analysis is to determine the accuracy of the SAS2H module of SCALE 4.3 in predicting isotopic concentrations of spent fuel assemblies. The objective is to develop a methodology for modeling assemblies similar to those evaluated within this analysis and to establish the consistency of SAS2H predictions. The results of this analysis may then be applied to future depletion calculations using SAS2H in which no measurements are available.

2. Quality Assurance

The Quality Assurance (QA) program applies to this analysis. The work reported in this document is part of the Waste Package Design analysis that will eventually support the License Application Design phase. This activity, when appropriately confirmed, can impact the proper functioning of the Mined Geologic Disposal System (MGDS) waste package; the waste package has been identified as an MGDS Q-List item important to safety and waste isolation (pp. 4, 15, Reference 5.1). The waste package is on the Q-List by direct inclusion by the Department of Energy (DOE), without conducting a QAP-2-3 evaluation. The Waste Package Development Department (WPDD) responsible manager has evaluated this activity in accordance with QAP-2-0, *Conduct of Activities*. The *Perform Criticality, Thermal, Structural, and Shielding Analyses* (Reference 5.2) evaluation has determined the preparation and review of this design analysis is subject to *Quality Assurance Requirements and Description* (Reference 5.3) requirements. As specified in NLP-3-18, this activity is subject to QA controls.

The analysis described in this document supports development of the disposal criticality analysis methodology. No designs were analyzed in this document. This document will not directly support any construction, fabrication, or procurement activity and therefore is not required to be procedurally controlled as TBV (to be verified). The calculation design inputs or information used in this document come from data accepted by the Nuclear Regulatory Commission and by the scientific and engineering community as established fact. The specific references are listed in Section 5 and identified in Section 7. The information is therefore not treated as unqualified data.

3. Method

The analytical model employed for this analysis was the SAS2H module of the SCALE sequence. Based upon fuel design, power history, and operating data for specific assemblies in the Trino Vercelles pressurized water reactor (PWR), a computational model was developed for use with the SAS2H module of SCALE. The SAS2H module is used to perform a fuel depletion analysis to predict the isotopic concentrations in localized areas of assembly pins (pellet samples) subsequent to irradiation and cooling time. The isotopic concentrations predicted by the SAS2H module are then compared with measured concentrations of the same localized areas (axial locations) of the assembly pins to determine the accuracy of the developed model. The measured isotopic concentrations used for comparisons in the analysis are obtained from a separate report (Reference 5.5).

4. Design Inputs

The sources for the design parameters are References 5.4 through 5.7. Reference 5.4 provides information on the molar masses; the assembly design, power history and operating parameters are obtained from Reference 5.5; and a list of trace elements in the fuel is derived from Reference 5.6; and light elements from Reference 5.7.

4.1 Design Parameters

The molar masses of selected elements are obtained from Reference 5.4 and are provided below within three significant figures. Precision beyond three significant figures is not necessary since the calculated isotopic concentrations from SCALE are only to three significant figures. Also, the weight per mole of enriched uranium is approximated by the weight per mole of natural uranium since the weight percent of enrichment is small.

Mole of natural uranium = 238 g,
 6.02×10^{23} atoms per mole (Reference 5.6)

General spent fuel characteristics for each sample are presented in Table 4-1 and include the initial ^{235}U enrichment, final burnup and the cooling time (Table 1, Reference 5.5). The initial enrichment ranges from 3.13 to 3.87 wt% ^{235}U and the burnup ranges from 11.529 to 24.548 GWd/MTU. The cooling time reflects that the measured isotopic concentrations were back-calculated to a "zero time". However, it is unknown whether the "zero time" is the time of shutdown or the time of discharge. Therefore, cooling time for the final cycle is specified as 10 days, since a shorter decay period would effect isotopes that come in part from the decay of short lived parent isotopes (pp. 7 and 10, Reference 5.5). Most notable is the decay of ^{239}U ($T_{1/2} = 23.5$ m) and ^{239}Np ($T_{1/2} = 2.35$ d) to form ^{239}Pu , which increases by roughly 1% during the first 10 days (pp. 7 and 10, Reference 5.5). The predicted concentration of ^{239}Pu relative to that which is measured, is used as an indication as to whether the zero day cooling time is appropriate.

Assembly design parameters are presented in Table 4-2 (Tables 2 and 3, Reference 5.5). Trino Vercelles utilizes two assembly types, the first is a Westinghouse 15 x 15 which has 16 outer cell positions removed so that the cruciform assemblies may fit between the square assemblies. A cross section of the square and cruciform assemblies as they fit together is presented in Figure 4-1, while the rod locations for the samples from each of the three square assemblies are included in Figures 4-2 through 4-4 (Figure 1, Reference 5.5). The cruciform assembly contains either Ag-In-Cd absorber rods, with moderator in the absorber cell positions when the absorber rods are withdrawn, or fuel rods with an enrichment of 2.72 wt % ^{235}U . The initial enrichment of the fuel rods in the cruciform assembly and the square assembly are given in Table 4-3 for each of the three assemblies (Table 4, Reference 5.5). The absorber rod composition was not known and therefore is approximated by the composition of an absorber rod from a similar design (Yankee Rowe) and is presented in Table 4-4 (Table 3 and p. 14, Reference 5.5). The fuel stack density was calculated for

the square assembly using the pellet diameter, active fuel length, number of fuel rods and weight of uranium per assembly.

General operating data is given in Table 4-5 and power histories for each assembly during the two cycles of irradiation are presented in Table 4-6 (Tables 5 and 8, Reference 5.5). The assemblies 509-104 and 509-032 were irradiated for 1 cycle while assembly 509-069 was irradiated for two cycles. The first cycle contained two intervals of downtime and therefore is divided into three periods. The boron concentration for cycle 2 was taken as the average for cycle 1, which ranged from 1300 ppm to 0 ppm. The cumulative burnup was only available for assembly 509-069, Table 4-5, and then applied to each sample's final burnup and length of cycles to determine the specific powers.

The operating temperatures are included in Tables 4-5 and 4-7 for the moderator and fuel (Tables 5 and 7, Reference 5.5). The moderator temperatures were calculated by a sinusoidal function of axial height, and inlet and outlet temperatures. The densities for the moderator were derived from correlations using the temperature and pressure data. Cladding and fuel temperatures were approximated using data from Yankee Rowe which is of a similar design (p. 13, Reference 5.5).

A list of trace elements in the fuel used in updating cross sections during the depletion analysis is presented in Table 4-8 and developed with consideration of elements used in (Table 1, Reference 5.6). A generic set of light element weights for PWRs that is typically used in depletion analyses is included in Table 4-9 (Table 17, Reference 5.7). Variations in light element masses per unit fuel in different PWRs are small when compared to this generic set (p. 2-2, Reference 5.9). This data is provided in units of kg/MTUO₂ or kg/MTU depending on the units required in the analysis.

Measured isotopic concentrations are presented in Table 4-10 and are given in mg/gU (Appendix D, Reference 5.5). The radiochemical analyses were performed for 14 samples coming from three different assemblies, and five different axial heights.

Table 4-1. Spent Fuel Characteristic Parameters for Trino Vercelles PWR

Test Assembly (Rod Number)	Axial Level Number	Axial Location from Bottom, cm	Enrichment, wt % ²³⁵ U	Burnup, GWd/MTU	Cooling Time, days
509-104 (M11)	7	79.2	3.87	12.042	0
509-032 (E11)	4	158.5	3.13	15.377	0
	7	79.2		15.898	0
	9	26.4		11.529	0
509-069 (E11)	1	237.7	3.13	12.859	0
	2	211.3		20.602	0
	4	158.5		23.718	0
	7	79.2		24.304	0
509-069 (E05)	4	158.5		23.867	0
	7	79.2		24.548	0
509-069 (L11)	4	158.5		23.928	0
	7	79.2		24.362	0
509-069 (L5)	4	158.5		24.330	0
	7	79.2		24.313	0

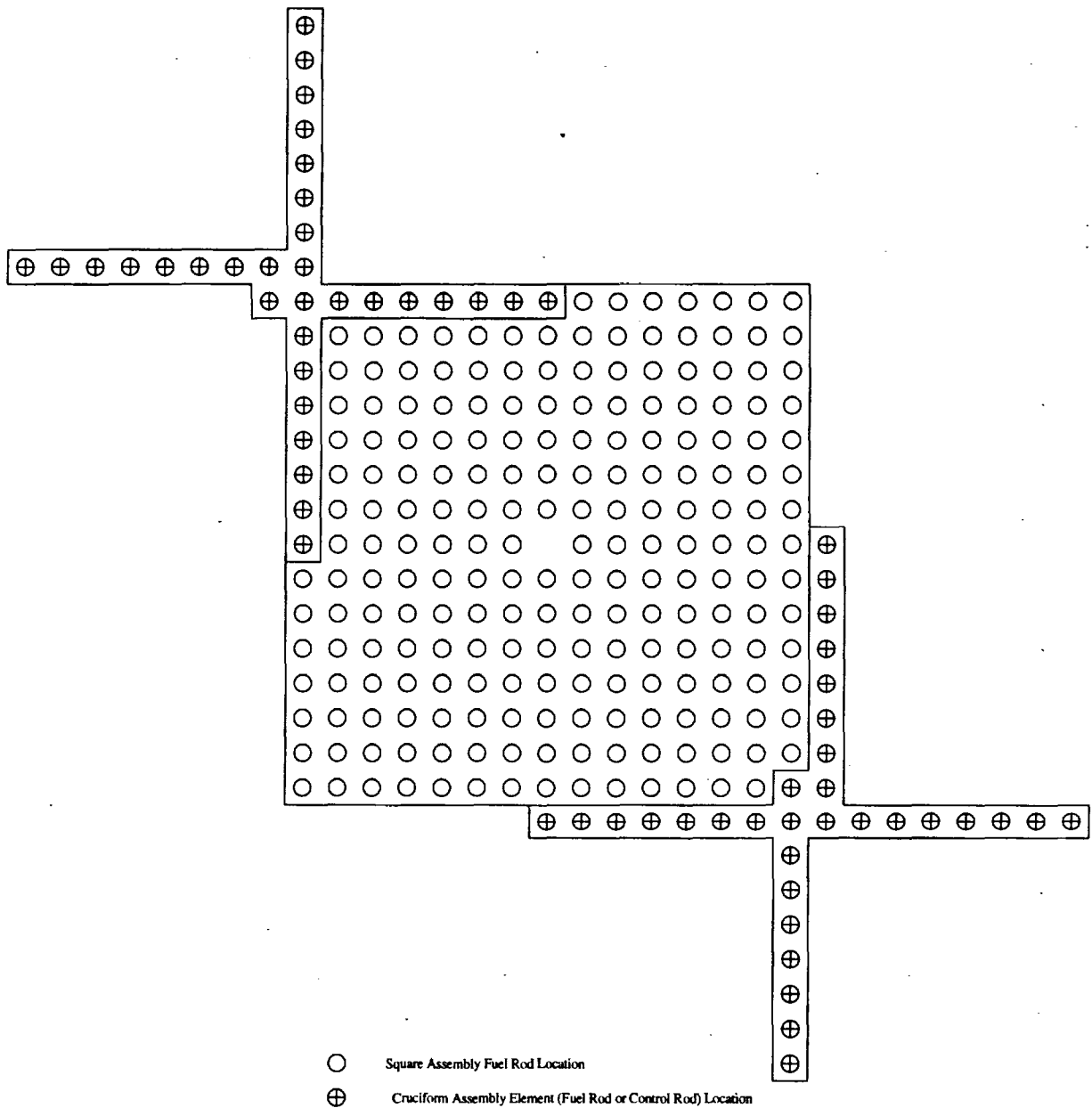
Reference 5.5

Table 4-2. Assembly Design Parameters for Trino Vercelles PWR

Parameter	Data
Square Fuel Assembly Data:	
Designer	Westinghouse
Lattice	15 x 15
Square Assemblies/Core, Cycle 1 (Cycle 2)	120 (112)
Number of Fuel Rods	208
Number of Inner Holes	1
Assembly Fuel, kg UO ₂	353.81
Assembly Tube Outer Side, cm	20.1
Channel Tube Thickness, cm	0.06
Channel Tube Material	SS-304
Fuel Rod Data:	
Type of Fuel Pellet	UO ₂
Pellet Stack Density, g/cm ³	10.079
Rod Pitch, cm	1.303
Rod Outside Diameter (OD), cm	0.9786
Rod Inside Diameter (ID), cm	0.9020
Active Fuel Length, cm	264.1
Clad Material	SS-304
Cruciform Fuel Assembly Data:	
Number of Cruciform Assemblies/Core	24
Fuel Weight/Assembly, kg UO ₂	44
Cruciform Control Rod Data:	
Number of Control Assemblies/Core	28
Number of Absorber Rods/Control Rod	32
Absorber Rod OD, cm	1.001
Absorber Rod Clad Thickness, cm	0.0432

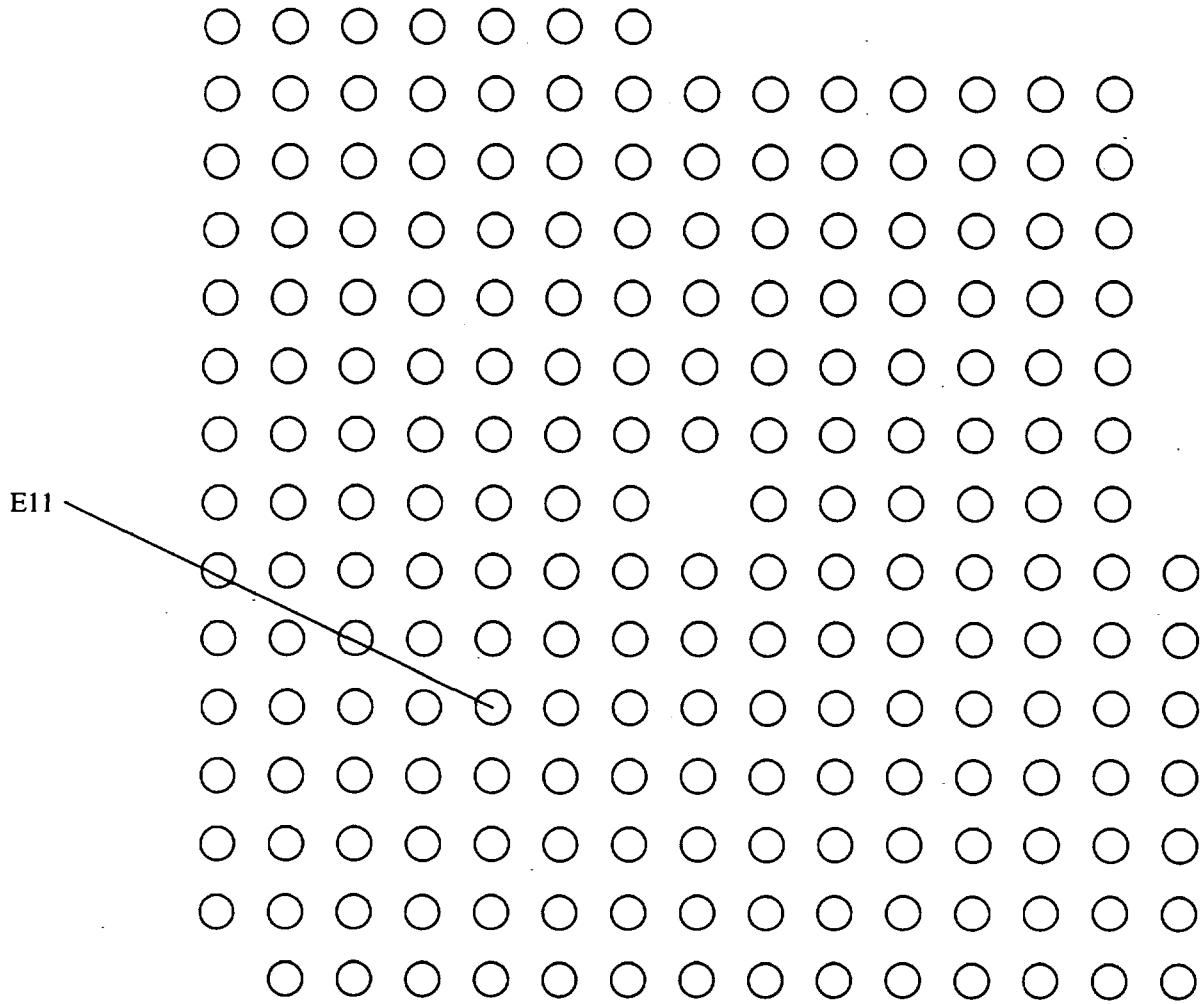
Reference 5.5

Figure 4-1. Cross Section of the Square and Cruciform Assemblies for Trino Vercelles



Reference 5.5

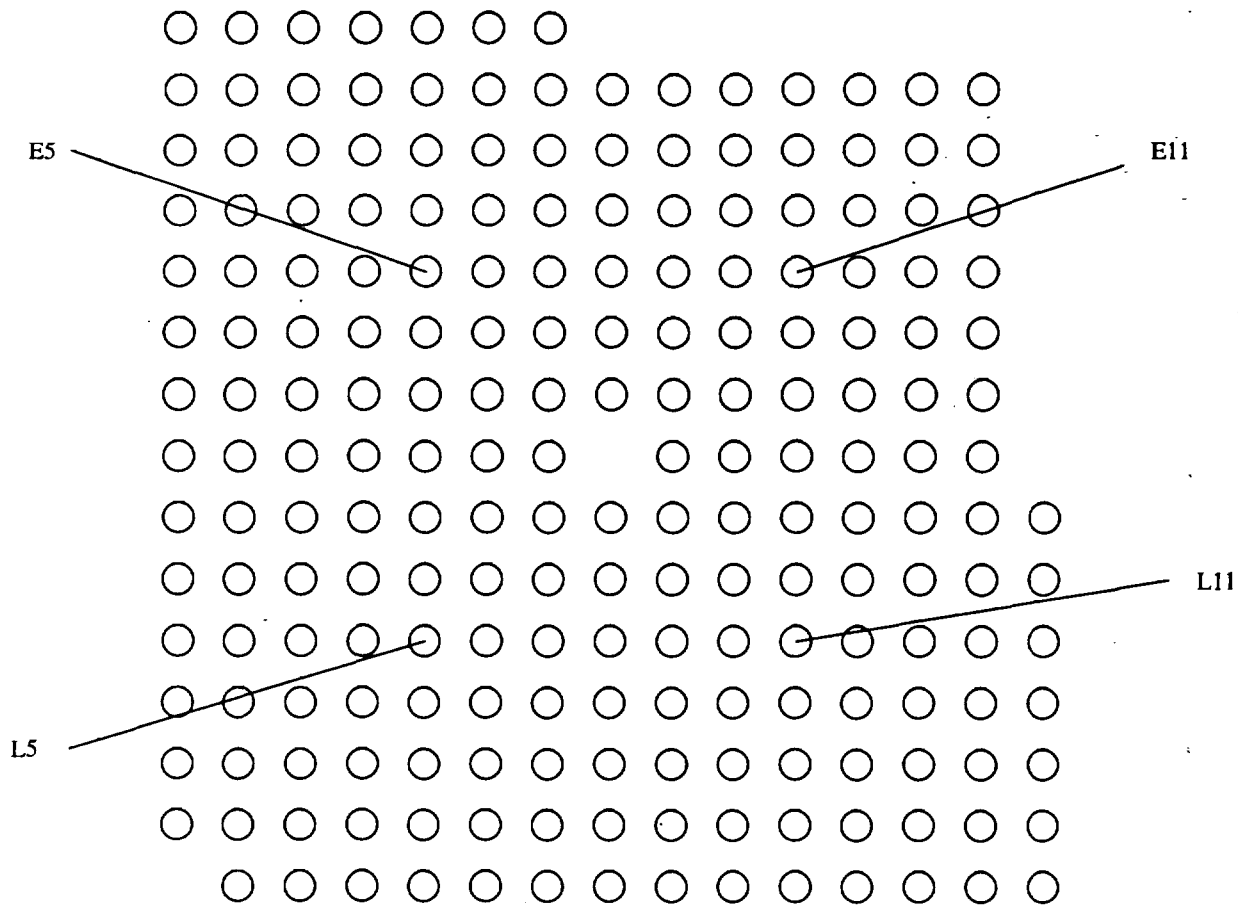
Figure 4-2. Fuel Rod Locations for Assembly 509-032



509-032

Reference 5.5

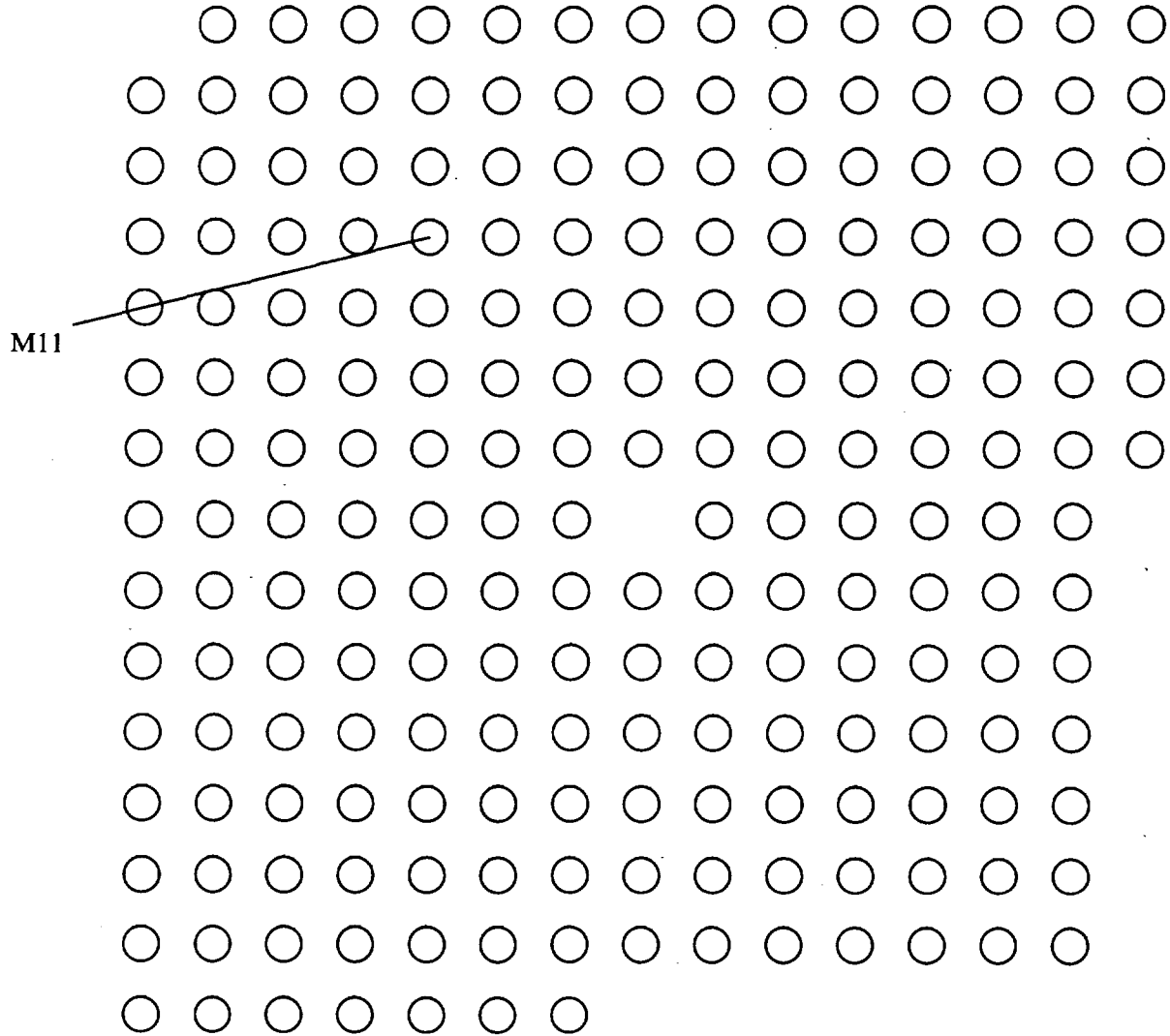
Figure 4-3. Fuel Rod Locations for Assembly 509-069



509-069

Reference 5.5

Figure 4-4. Fuel Rod Locations for Assembly 509-104



509-104

Reference 5.5

Table 4-3. Initial Fuel Composition for Trino Vercelles PWR

Uranium Isotope	Assembly Identification Number		
	509-104	509-32	509-069
wt % ²³⁵ U	3.897	3.13	3.13
wt % ²³⁴ U	0.035	0.028	0.028
wt % ²³⁶ U	0.018	0.014	0.014
wt % ²³⁸ U	96.050	96.828	96.828

Reference 5.5

Table 4-4. Absorber Rod Composition

Material	Weight Percent
Ag	80
In	15
Cd	5
Density, g/cm ³	10.159

Reference 5.5

Table 4-5. Core Operating Data for Trino Vercelles PWR

Parameter	Cycle/Period Data			
	1	1	1	2
Cycle	I	II	III	I
Starting Date	10/23/64	8/31/65	7/11/66	5/20/70
Uptime, days	226	263	292	416
Downtime, days	86	51	1117	See Table 4-1
Cladding Temperature, K	570	570	570	570
Coolant Average Temperature, K	282	278	278	269
Control Rod Insertion, %	30	4	4	0
Average Boron, ppm	1175	850	325	650
509-069 Burnup, GWd/MTU	2.726	4.927	6.327	7.720
Core Burnup, GWd/MTU	2.260	4.085	5.245	NA

Reference 5.5

Table 4-6. Operating Data for Trino Vercelles PWR

Assembly Identification Number	509-104	509-032	509-032	509-032	509-069	509-069	509-069
Fuel Rod Location	M11	E11	E11	E11	E11	E11	E11
Axial Level No.	7	4	7	9	1	2	4
Burnup, GWd/MTU	12.042	15.377	15.898	11.529	12.859	20.602	23.718
Specific Power, MW/MTU							
Cycle 1, I	10.390	13.268	13.717	9.947	7.148	11.452	13.184
Cycle 1, II	16.138	20.607	21.309	15.451	11.101	17.786	20.476
Cycle 1, III	18.663	23.831	24.639	17.868	12.840	20.571	23.683
Cycle 2	-	-	-	-	10.997	17.619	20.283
Assembly Identification Number							
Assembly Identification Number	509-069	509-069	509-069	509-069	509-069	509-069	509-069
Fuel Rod Location	E11	E5	R5	L11	L11	L5	L5
Axial Level No.	7	4	7	4	7	4	7
Burnup, GWd/MTU	24.304	23.867	24.548	23.928	24.362	24.330	24.313
Specific Power, MW/MTU							
Cycle 1, I	13.509	13.266	13.645	13.300	13.542	13.524	13.514
Cycle 1, II	20.982	20.605	21.193	20.657	21.032	21.004	20.990
Cycle 1, III	24.268	23.832	24.512	23.893	24.326	24.294	24.277
Cycle 2	20.785	20.411	20.993	20.463	20.834	20.807	20.792

Reference 5.5

Table 4-7. Operating Temperatures for Trino Vercelles PWR

Level No.	Effective Fuel Temperature, K		Effective Water Temperature, K		Water Density, g/cm ³	
	Cycle 1	Cycles 1 and 2	Cycle 1	Cycles 1 and 2	Cycle 1	Cycle 2
1	915	915	567	563	0.7365	0.7555
2	968	968	565	561	0.7407	0.7595
4	1015	1015	557	553	0.7554	0.7730
7	1001	1001	543	540	0.7795	0.7953
9	927	927	537	534	0.7885	0.8032

Reference 5.5

Table 4-8. Nuclides Updated in SAS2H

⁸³ Kr	⁸⁵ Kr	⁸⁹ Y	⁹⁰ Sr	⁹⁵ Mo	⁹³ Zr
⁹⁴ Zr	⁹⁴ Nb	⁹⁵ Zr	⁹⁹ Tc	¹⁰¹ Ru	¹⁰³ Rh
¹⁰⁵ Rh	¹⁰⁶ Ru	¹⁰⁵ Pd	¹⁰⁸ Pd	¹⁰⁹ Ag	¹²⁴ Sb
¹³¹ Xe	¹³² Xe	¹³⁴ Cs	¹³⁵ Xe	¹³⁵ Cs	¹³⁶ Xe
¹³⁶ Ba	¹³⁷ Cs	¹³⁹ La	¹⁴¹ Pr	¹⁴³ Pr	¹⁴³ Nd
¹⁴⁴ Ce	¹⁴⁵ Nd	¹⁴⁷ Nd	¹⁴⁷ Pm	¹⁴⁷ Sm	¹⁴⁸ Pm
¹⁴⁹ Sm	¹⁵⁰ Sm	¹⁵¹ Sm	¹⁵² Sm	¹⁵³ Eu	¹⁵⁴ Eu
¹⁵⁵ Gd	¹⁵⁵ Eu				

Reference 5.6

Table 4-9. Light Element Mass per Unit of Fuel for a Typical PWR

Element	kg/MTU
O	135.0
Cr	5.9
Mn	0.33
Fe	13.0
Co	0.075
Ni	9.9
Zr	221.0
Nb	0.71
Sn	3.6

Reference 5.7

Table 4-10. Measured Isotopic Concentrations (mg/gU)

Assembly	509-104 M11	509-032 E11	509-032 E11	509-032 E11	509-069 E11	509-069 E11	509-069 E11
Location	7	4	7	9	1	2	4
Burnup, GWd/MTU	12.042	15.377	15.898	11.529	12.859	20.602	23.718
²³⁵ U	2.662E01	1.728E01	1.661E01	2.017E01	1.946E01	1.436E01	1.248E01
²³⁶ U	2.736E00	2.834E00	2.739E00	2.502E00	2.453E00	3.317E00	3.610E00
²³⁸ U	9.513E02	9.558E02	9.558E02	9.595E02	9.587E02	9.518E02	9.493E02
²³⁹ Pu	4.586E00	5.266E00	5.234E00	4.418E00	4.580E00	5.755E00	5.895E00
²⁴⁰ Pu	7.165E-01	1.118E00	1.137E00	7.750E-01	8.400E-01	1.520E00	1.755E00
²⁴¹ Pu	3.475E-01	6.140E-01	6.180E-01	3.690E-01	4.000E-01	8.850E-01	1.030E00
²⁴² Pu	3.135E-02	8.638E-02	9.487E-02	3.803E-02	4.600E-02	1.720E-01	2.435E-01
²⁴³ Am	-	-	-	-	-	2.394E-02	4.533E-02
Assembly	509-069 E11	509-069 E5	509-069 E5	509-069 L11	509-069 L11	509-069 L5	509-069 L5
Location	7	4	7	4	7	4	7
Burnup, GWd/MTU	24.304	23.867	24.548	23.928	24.362	24.330	24.313
²³⁵ U	1.235E01	1.291E01	1.221E01	1.282E01	1.225E01	1.297E01	1.231E01
²³⁶ U	3.638E00	3.520E00	3.540E00	3.753E00	3.465E00	3.471E00	3.569E00
²³⁸ U	9.492E02	9.492E02	9.483E02	9.485E02	9.482E02	9.486E02	9.472E02
²³⁹ Pu	6.070E00	5.950E00	5.980E00	6.060E00	5.995E00	6.060E00	5.970E00
²⁴⁰ Pu	1.825E00	1.760E00	1.785E00	1.790E00	1.810E00	1.770E00	1.790E00
²⁴¹ Pu	1.060E00	1.050E00	1.055E00	1.050E00	1.055E00	1.060E00	1.060E00
²⁴² Pu	2.575E-01	2.400E-01	2.540E-01	2.470E-01	2.590E-01	2.440E-01	2.500E-01
²⁴³ Am	4.584E-02	-	4.615E-02	4.452E-02	4.247E-02	-	-

Reference 5.5

4.2 Criteria

The design of the waste package will depend on waste package configuration criticality analyses performed using an acceptable disposal criticality analysis methodology. Criteria that relate to the development and design of repository and engineered barrier components are derived from the applicable requirements and planning documents. The Engineered Barrier Design Requirements Document (EBDRD, Reference 5.10) provides requirements for engineered barrier segment design. The Repository Design Requirements Document (RDRD, Reference 5.11) provides requirements for repository design. The Controlled Design Assumptions Document (Reference 5.12) provides guidance for requirements listed in the EBDRD and RDRD which have unqualified or unconfirmed data associated with the requirement.

This analysis supports the disposal criticality analysis methodology by providing input, in the form of fuel depletion results, to benchmark calculations which address the prediction of both spent fuel isotopic compositions and their associated reactivity. These benchmark calculations will contribute to the determination of bias values in the method of critical multiplication factor calculation that is implemented by the analytic tools to be used in the disposal criticality methodology. The requirements for utilizing the bias in the method of calculation of the critical multiplication factor for disposal configurations containing spent nuclear fuel are located in Section 3.2.2.5 of the RDRD and Section 3.2.2.6 of the EBDRD. This analysis does not satisfy these requirements, but the results from this analysis will be used as input to subsequent analyses which will satisfy these requirements.

4.3 Assumptions

4.3.1 Measured isotopic concentrations from Reference 5.5 are calculated and corrected to either the time of shutdown or the time of discharge. Since it is unclear which time the concentrations are back-calculated to, a cooling time of 10 days is assumed. The basis for this assumption is that a decay period longer than 10 days does not significantly effect isotopic concentrations, but a shorter decay period effects concentrations of those isotopes with short lived parents, specifically ^{239}Pu which increases by approximately 1% in the first ten days. This assumption is used in Section 7.5.

4.4 Codes and Standards

There are no applicable codes or standards for this design analysis.

5. References

- 5.1 *Yucca Mountain Site Characterization Project Q-List*, YMP/90-55Q REV 4, Yucca Mountain Site Characterization Project.
- 5.2 *QAP-2-0 Activity Evaluations: ID #WP-20, Perform Criticality, Thermal, Structural, and Shielding Analyses*, Civilian Radioactive Waste Management System (CRWMS) Management and Operating Contractor (M&O), August 3, 1997.
- 5.3 *Quality Assurance Requirements and Description*, DOE/RW-0333P REV 7, U.S. Department of Energy (DOE) Office of Civilian Radioactive Waste Management (OCRWM).
- 5.4 *Nuclides and Isotopes*, General Electric Company, 14ed., 1989.
- 5.5 *An Extension of the Validation of SCALE (SAS2H) Isotopic Prediction for PWR Spent Fuel*, ORNL/TM-13317, Oak Ridge National Laboratory, September 1996.
- 5.6 *SCALE-4 Analysis of Pressurized Water Reactor Critical Configurations: Volume 2- Sequoyah Unit 2 Cycle 3*, ORNL/TM-12294/V2, March 1995.
- 5.7 *Validation of the Scale System for PWR Spent Fuel Isotopic Composition Analyses*, ORNL/TM-12667, Oak Ridge National Laboratory, March 1995.
- 5.8 *SCALE 4.3, A Modular Code System for Performing Standardized Computer Analyses for Licensing Evaluation for Workstations and Personal Computers*, NUREG/CR-0200 REV 5, ORNL/NUREG/CSD-2/R5, Volumes 1-3, Oak Ridge National Laboratory.
- 5.9 *Software Qualification Report for the SCALE Modular Code System*, Document Identifier Number: 30011-2002 REV 01, CRWMS M&O.
- 5.10 *Engineered Barrier Design Requirements Document*, YMP/CM-0024, REV 00, ICN 01, DOE OCRWM.
- 5.11 *Repository Design Requirements Document*, YMP/CM-0023, REV 00, ICN 01, DOE OCRWM.
- 5.12 *Controlled Design Assumptions Document*, DI#: B00000000-01717-4600-00032 REV 04, ICN 01, CRWMS M&O.

6. Use of Computer Software

- A.** Reference 5.8 describes the SAS2H module of SCALE 4.3 that is used with the 44GROUPNDF5 cross section library to calculate the isotopic concentrations for the specified burnup and cooling time. The computer code's spatially independent point depletion model is appropriate for comparison with pellet sample measurements, and is used within the range of validation, as described in Reference 5.9, in accordance with the QAP-SI series procedures. SCALE is obtained from the Software Configuration Management in accordance with appropriate procedures. SCALE's CSCI number is 30011 V4.3 and is installed on the WPDD HP 9000, 700 Workstation with CRWMS M&O tag number 110433.

- B.** *Lotus 1-2-3 Release 5 for Windows 95* is an Acquired Software spreadsheet program as defined in QAP-SI-0. User defined formulas and/or algorithms, inputs and results, are documented in the appropriate sections.

7. Design Analysis

The SAS2H module of SCALE 4.3 is used to perform one-dimensional (1-D) neutron transport and point depletion analyses on the Trino Vercelles samples using the preferred 44GROUPNDF5 cross-section library. To properly model the neutron flux spectrum and the nuclide composition changes, it is necessary to define the materials, temperatures, and geometry of the fuel assembly. This is accomplished with the use of data blocks in which similar parameters are grouped together.

7.1 SCALE Input Data Blocks 1, 2, and 3

Data blocks 1 through 3 define the SCALE module to be used, the title of the input file, the cross section library to be used, and the lattice type to be modeled. The module used is SAS2H and the cross sectional library is 44GROUPNDF5, abbreviated as 44GROUP. The 44GROUP cross section library is recommended by Oak Ridge National Laboratory. Since SAS2H is only to be used for isotopic depletion/generation, the 'parm=skipshipdata' option is used so that a shipping cask shielding analysis is not performed. The title is arbitrary and should contain information that is sample specific, while the lattice type is "latticecell" to reflect the array characteristic of the assembly.

7.2 SCALE Input Data Block 4

Data block 4 defines the material compositions present in the assembly. A unique mixture number is assigned to each composition, and follows the form of mixture 1 for fuel, mixture 2 for cladding and mixture 3 for moderator.

The fuel mixture is specified as UO_2 using the density and isotopic weight percentages from Tables 4-2 and 4-3 and the fuel temperature is obtained from Table 4-7. Isotopes which are selected as needing their cross sections updated during the depletion analysis are included in the fuel mixture. A standard list of trace fuel elements is given in Table 4-8 and defined in the fuel mixture to have a concentration of 10^{-20} atoms/barn-cm. The cladding material of stainless steel 304 is defined by 'ss304' with the temperature given in Table 4-7. Finally, the absorber material is defined as mixture 4 with the density and composition given in Table 4-4.

It was discovered that the temperature is incorrectly specified for the trace elements in the input file for the sample at axial height 4 of rod L11 from assembly 509-069. A temperature of 1001 K is specified when a temperature of 1015 K should have been specified. However, the correct temperature was specified for the fuel component of the same mixture. When different temperatures are specified for the same mixture, the SAS2H code uses the higher temperature of the two, in this case 1015 K. Therefore, the correct temperature was used during the calculation.

The moderator temperature, density and boron concentration are given in Tables 4-5 and 4-7 and is composed of H_2O and boron. The boron is defined as an arbitrary material with the moderator

density and temperature, with a volume fraction equal to the average boron concentration of Cycle 1 Period I, and a standard boron composition of from the Standard Composition Library designated as 5000.

7.3 SCALE Input Data Blocks 5 Through 7

The unit fuel rod cell geometry is defined in data block 5. The 'squarepitch' designation for the type of lattice is appropriate since the fuel assembly consists of a square array of fuel rods. Fuel rod, cladding and pitch dimensions are given in Table 4-2 with the mixture number for each composition defined in Section 7.2. The gap mixture is defined as 0.

Data block 6 allows the user to specify such parameters as the spatial mesh, angular quadrature and the convergence criteria. It is determined that the default values are sufficient and such options are not used in this design model.

In data block 7 the user defines general assembly data and determines the level of detail in which the assembly is to be modeled. The number of fuel rods per assembly is given in Table 4-2 and the length is calculated so that an assembly contains 1 MTU, using the following equation:

$$Length = \frac{1}{\frac{\pi}{4}(POD)^2(PDen)(NFR)} * \frac{270gUO_2}{238gU} * \frac{10^6gU}{1MTU} \quad \text{Equation 7-1}$$

Where:

Length = Length Required for an Assembly to Contain 1 MTU (cm)

POD = Fuel Pellet Diameter (cm)

PDen = Fuel Pellet Stack Density (gUO_2/cm^3)

NFR = Number of Fuel Rods

Since, measured isotopic concentrations are presented in milligrams of isotope per grams of uranium and SCALE presents concentrations in grams of isotope per assembly, it is possible to alter the length so that the assembly contains 1 MTU. This is possible since the 1-D transport calculation is axially independent, consequently the length of the assembly does not impact the neutron flux spectrum nor the nuclide cross sections. The resulting length for Trino Vercelles assemblies is 823.10 cm.

Each assembly was irradiated for one or two cycles. It is determined that approximately 80 days per cross section library is sufficient to accurately model the change in nuclide cross sections with increasing burnup, without over-burdening the SAS2H code. Therefore, the number of libraries per cycle are specified as five. To obtain the concentrations of all interested nuclides, a print level of 5 is chosen, while an input level of 2 is defined so that a Path B model may be utilized. The number of light elements is nine and is determined from Table 4-9, while the number of zones is three or

five, depending on the presence of the absorber rod, which is determined by the Path B model described in Section 7.4.

SAS2H calculates a linear boron concentration let down over each cycle when the number of libraries per cycle are specified as greater than one. The methodology calculates an initial boron concentration equal to 1.9 times the average boron concentration and a final boron concentration calculated as 0.1 times the average boron concentration. These two end points are used to determine a linear fit for the boron concentration let down, which is used to calculate the boron concentrations for each library in the cycle. However, a problem arises when the cycle specified is not a full cycle but rather an interval of a cycle. (It is sometimes necessary to specify an interval as a cycle in SAS2H so that either a period of downtime during the actual cycle or a power change during the actual cycle may be modeled.) In this case SAS2H would let down the boron concentration to 0.1 times the average boron concentration for the first interval and jump up to 1.9 times the average boron concentration for the beginning of the second cycle. In reality, this jump does not exist and, therefore, the modeling of the boron concentration let down is inexact. A solution to this problem would be to split all cycles into approximately 80 day intervals and change the number of libraries per cycle to one. However, experience with the SAS2H code indicates that the error due to an inexact boron concentration let down function, such as described here, is usually less than 0.5%. Therefore, allowing the SAS2H code to perform a boron concentration over an interval of a cycle is an acceptable approximation to the actual boron concentration let down.

7.4 SCALE Input Data Block 8

Unfortunately, the exact locations of control rod cruciform assemblies and fuel rod cruciform assemblies are not known. However, it is known that there are 32 fuel rods per fuel cruciform assembly and 24 fuel cruciform assemblies in a core containing 120 square assemblies. Thus, it is possible to find an average number of fuel rods from the fuel cruciform assembly to be added to a square assembly. The result is an average of 6 fuel rods from a fuel cruciform assembly and 10 absorber rods from the absorber cruciform assembly to be added to the square assembly. Furthermore, the fuel rods from the fuel cruciform assemblies are approximated as having the same properties, including initial enrichments, as the fuel rods from the square assembly. This creates an effective square assembly in which there are 214 fuel rods, one centrally located empty fuel cell, and 10 absorber rod cells on the outside of the assembly, that contain moderator when the absorber rods are withdrawn.

The Path B model for Trino Vercelles is dependent upon whether the control rods are inserted to the axial height of the sample, since the effects of the absorber rods are determined to be negligible if the absorber rods are not inserted to the axial height of the sample. The control rods were inserted 30% for period I of cycle 1, 4% for the rest of the first cycle, and were not inserted for cycle 2. Therefore, the effects of the absorber rods are only modeled at an axial height of 185 cm or greater, from the bottom of the assembly during period I of cycle 1. For the remainder of cycle 1 and cycle 2 the effects of the absorber rods are not experienced by any of the samples.

If it is determined that the sample does not experience the effects of the absorber rods, then the model centralizes an empty unit cell surrounded by an homogenized fuel and moderator mixture, that conserves the fuel to moderator volume ratio, and further surrounded by the assembly casing. The equation below is used to determine the number of fuel unit cells that surround the empty unit cell. All of the following equations used to calculate the Path B model dimensions are derived. The results of the fuel-unit-cell calculations are presented in Table 7-1, and the resulting Path B model dimensions are presented in Table 7-2.

$$x = \frac{\left(\frac{F}{M}\right)(CUCMV)}{(FV) - \left(\frac{F}{M}\right)(MV)} \quad \text{Equation 7-2}$$

$$\frac{F}{M} = \frac{(NFR)\left(\frac{\pi}{4}\right)(POD)^2}{(NFR)[RP^2 - \left(\frac{\pi}{4}\right)(COD)^2] + (NHole)[RP^2]} \quad \text{Equation 7-3}$$

$$CUCMV = RP^2 \quad \text{Equation 7-4}$$

$$FV = \left(\frac{\pi}{4}\right)(POD)^2 \quad \text{Equation 7-5}$$

$$MV = RP^2 - \left(\frac{\pi}{4}\right)(COD)^2 \quad \text{Equation 7-6}$$

Where:

- x = Number of Unit Fuel Cells per Empty Unit Cell
- F/M = Fuel to Moderator Volume Ratio
- NFR = Number of Fuel Rods
- POD = Fuel Pellet Outer Diameter
- RP = Rod Pitch
- COD = Cladding Outer Diameter
- NHole = Number of Empty Unit Cells
- CUCMV = Central Unit Cell Moderator Volume
- FV = Fuel Volume of One Fuel Unit Cell
- MV = Moderator Volume of One Fuel Unit Cell

Once the number of fuel cells per empty unit cell is determined the geometry of the Path B model is determined. Since the empty unit cell is centralized, the radius for the moderator is calculated with the following equation:

$$R_1 = \sqrt{\left(\frac{1}{\pi}\right)RP^2} \quad \text{Equation 7-7}$$

Where:

R_1 = Radius of Moderator Surrounding Empty Unit Cell

The area of an annular region is calculated by the difference between the outer circular area and the inner circular area. Equation 7-8 is the basis for the Equations 7-9 and 7-10 which determine the radii of the homogenized fuel zone and the outer moderator zone.

$$ARA = \pi(ORAR^2 - IRAR^2) \quad \text{Equation 7-8}$$

Where:

ARA = Annular Region Area

ORAR = Outer Radius of Annular Region

IRAR = Inner Radius of Annular Region

The area of the homogenized fuel zone surrounding the empty unit cell is equal to the number of fuel unit cells surrounding the empty unit cell multiplied by the area of a fuel unit cell. Consequently, the radius of the homogenized fuel zone is computed with the following equation:

$$R_2 = \sqrt{\left(\frac{x}{\pi}\right)RP^2 + R_1^2} \quad \text{Equation 7-9}$$

Where:

R_2 = Radius of Homogenized Fuel and Moderator Zone

The mixture number of the homogenized fuel and moderator mixture must be specified as 500. The code then determines the composition of the region using cell averages or homogenized densities of the fuel rod cell.

The Trino Vercelles reactor assembly is surrounded by a stainless steel channel casing. The total casing volume is determined by the thickness and the outer width. The radius of the stainless steel casing is calculated from the following equation:

$$R_3 = \sqrt{\frac{(x+1)}{\pi * N_{Cell}} [AW^2 - (AW - thick)^2] + R_2^2} \tag{Equation 7-10}$$

Where:

R_3 = Radius of Assembly Casing Zone

AW = Assembly Outside Width

thick = Thickness of Casing

If the sample is located at an axial height in which the control rod is inserted, then the model includes the absorber rods. In this model, the absorber rod unit cell is centralized with a surrounding homogenized fuel and moderator zone and surrounding assembly casing. The radii for the first two zones are the radii for the absorber rod, with corresponding absorber and cladding material. The third zone is the moderator surrounding the absorber rod and the radius is calculated with Equation 7-7.

The radii for the surrounding homogenized fuel and moderator, and the surrounding assembly casing are determined using the same Equations 7-9 and 7-10, respectively. However, the calculation of 'x' is different in that the equations for the fuel to moderator ratio and the central unit cell moderator volume are slightly altered, as shown below in equations 7-11 and 7-12.

$$\frac{F}{M} = \frac{(NFR)(\frac{\pi}{4})(POD)^2}{(NFR)[RP^2 - (\frac{\pi}{4})(COD)^2] + (NA)[RP^2 - (\frac{\pi}{4})(AOD)^2] + [RP^2]} \tag{Equation 7-11}$$

Where:

NA = Number of Absorber Rods

AOD = Absorber Rod Outer Diameter

$$CUCMV = RP^2 - (\frac{\pi}{4})(AOD)^2 \tag{Equation 7-12}$$

Table 7-1. Calculation of Fuel Unit Cell per Guide Tube Unit Cell

Centralized Region	F/M	CUCMV, cm ²	FV, cm ²	MV, cm ²	x
Empty Unit Cell	0.6186	1.6978	0.6390	0.9457	19.4545
Absorber Rod	0.6415	0.9108	0.6390	0.9457	18.0337

Table 7-2. Path B Model Dimensions

	Empty Unit Cell		Absorber Rod	
	Radius, cm	Composition	Radius, cm	Composition
R ₁	0.7351	Moderator	0.4573	Absorber
R ₂	3.3248	Fuel/Mod.	0.5005	SS-304
R ₃	3.3353	SS-304	0.7351	Moderator
R ₄	NA	NA	3.2076	Fuel/Mod.
R ₅	NA	NA	3.2177	SS-304

7.5 SCALE Input Data Blocks 9 Through 16

Data block 9 is used to describe the power history of the reactor. The specific power, fuel irradiation period, the length of downtime, the fraction of boron and moderator density, and the temperature during the cycle may all be defined. The specific power is in units of MW/MTU while the irradiation period and length of downtime are both defined in days and are found in Tables 4-5 and 4-6. The boron and moderator density fractions are determined by dividing the cycle, or period, value by the value specified in data block 4, using the following command options: 'bfrac=' for boron fraction, and 'h2ofrac=' for the moderator density fraction. The fuel temperatures were provided as an average over both cycles and therefore the option to specify the cycle specific temperature is not used.

Light elements and their effective weight, in kg per assembly, are entered in data block 10. Table 4-9 provides a generic set of light elements and their weights, in kg per MTU. Since the fuel length is altered so that there is 1 MTU per assembly the use of light elements with weights of kg per MTU is appropriate.

Data blocks 11 through 15 describe parameters used in the radial shielding analysis of a shipping cask and are not necessary in performing the depletion analysis. Data block 16 denotes the end of the SCALE input.

7.6 Comparison of Calculated and Measured Concentrations

Comparisons of corresponding calculated and measured concentrations are performed on a percent difference basis. The difference between the measured and the calculated value is divided by the measured value to determine the accuracy of the SAS2H calculation. A positive percent difference represents an over-prediction by the code, while a negative percent difference represents an under-prediction by the code.

Measured concentrations presented in Table 4-10 have units of mg of isotope per g of U, while calculated concentrations presented in Table 7-3 have units of g of isotope per Metric Ton of U. Therefore, the measured concentrations must be multiplied by 10^3 to obtain similar units. Percent differences are presented in Table 7-4.

7.7 Results

SAS2H predicted isotopic concentrations are presented in Table 7-3. The calculated concentrations are obtained through the methodology described in Sections 7.1 through 7.5, and with the input parameters defined in Section 4.1. Calculated concentrations are compared with measured concentrations as described in Section 7.6 to determine the accuracy of the SAS2H module. Results of the comparison, in the form of percent differences, are presented in Table 7-4.

Table 7-3. Calculated concentrations (g/MTU)

Assembly	509-104 M11	509-032 E11	509-032 E11	509-032 E11	509-069 E11	509-069 E11	509-069 E11
Location	7	4	7	9	1	2	4
Burnup, GWD/MTU	12.042	15.377	15.898	11.529	12.859	20.602	23.718
²³⁵ U	2.69E4	1.77E4	1.73E4	2.03E4	1.99E4	1.51E4	1.29E4
²³⁶ U	2.46E3	2.62E3	2.68E3	2.15E3	2.27E3	3.09E3	3.41E3
²³⁸ U	9.52E5	9.56E5	9.56E5	9.59E5	9.58E5	9.52E5	9.49E5
²³⁹ Pu	4.45E3	5.25E3	5.24E3	4.37E3	4.71E3	5.93E3	6.22E3
²⁴⁰ Pu	6.77E2	1.11E3	1.14E3	7.53E2	8.68E2	1.51E3	1.78E3
²⁴¹ Pu	3.06E2	5.82E2	5.98E2	3.38E2	3.94E2	8.39E2	1.03E3
²⁴² Pu	2.72E1	8.26E1	8.86E1	3.52E1	4.55E1	1.62E2	2.40E2
²⁴³ Am	-	-	-	-	-	2.32E1	4.04E1
Assembly	509-069 E11	509-069 E5	509-069 E5	509-069 L11	509-069 L11	509-069 L5	509-069 L5
Location	7	4	7	4	7	4	7
Burnup, GWD/MTU	24.304	23.867	24.548	23.928	24.362	24.330	24.313
²³⁵ U	1.25E4	1.28E4	1.24E4	1.28E4	1.25E4	1.26E4	1.25E4
²³⁶ U	3.46E3	3.42E3	3.47E3	3.42E3	3.46E3	3.45E3	3.46E3
²³⁸ U	9.49E5	9.49E5	9.48E5	9.49E5	9.49E5	9.48E5	9.49E5
²³⁹ Pu	6.16E3	6.23E3	6.18E3	6.24E3	6.16E3	6.27E3	6.16E3
²⁴⁰ Pu	1.80E3	1.79E3	1.82E3	1.79E3	1.81E3	1.82E3	1.80E3
²⁴¹ Pu	1.04E3	1.04E3	1.05E3	1.04E3	1.04E3	1.06E3	1.04E3
²⁴² Pu	2.52E2	2.43E2	2.58E2	2.45E2	2.53E2	2.55E2	2.52E2
²⁴³ Am	4.26E1	-	4.41E1	4.16E1	4.30E1	-	-

Table 7-4. Percent Difference Between Measured and Calculated $[(C/M-1)*100]$

Assembly	509-104 M11	509-032 E11	509-032 E11	509-032 E11	509-069 E11	509-069 E11	509-069 E11
Location	7	4	7	9	1	2	4
Burnup, GWd/MTU	12.042	15.377	15.898	11.529	12.859	20.602	23.718
²³⁵ U	1.05	2.43	4.15	0.64	2.26	5.15	3.37
²³⁶ U	-10.09	-7.55	-2.15	-14.07	-7.46	-6.84	-5.54
²³⁸ U	0.07	0.02	0.02	-0.05	-0.07	0.02	-0.03
²³⁹ Pu	-2.97	-0.30	0.11	-1.09	2.84	3.04	5.51
²⁴⁰ Pu	-5.51	-0.72	0.26	-2.84	3.33	-0.66	1.42
²⁴¹ Pu	-11.94	-5.21	-3.24	-8.40	-1.50	-5.20	0.00
²⁴² Pu	-13.24	-4.38	-6.61	-7.44	-1.09	-5.81	-1.44
²⁴³ Am	-	-	-	-	-	-3.09	-10.88
Assembly	509-069 E11	509-069 E5	509-069 E5	509-069 L11	509-069 L11	509-069 L5	509-069 L5
Location	7	4	7	4	7	4	7
Burnup, GWd/MTU	24.304	23.867	24.548	23.928	24.362	24.330	24.313
²³⁵ U	1.21	-0.85	1.56	-0.16	2.04	-2.85	1.54
²³⁶ U	-4.89	-2.84	-1.98	-8.87	-0.14	-0.61	-3.05
²³⁸ U	-0.02	-0.02	-0.03	0.05	0.08	-0.06	0.19
²³⁹ Pu	1.48	4.71	3.34	2.97	2.75	3.47	3.18
²⁴⁰ Pu	-1.37	1.70	1.96	0.00	0.00	2.82	0.56
²⁴¹ Pu	-1.89	-0.95	-0.47	-0.95	-1.42	0.00	-1.89
²⁴² Pu	-2.14	1.25	1.57	-0.81	-2.32	4.51	0.80
²⁴³ Am	-7.07	-	-4.44	-6.56	1.25	-	-

8. Conclusions

The accuracy in which the SAS2H module is able to predict isotopic concentrations is indicated by the percent differences presented in Table 7-4. Inspection of such results reveals that the code has a tendency to under-predict ^{236}U . Percent differences from this analysis are compared with results from Reference 5.6, in which similar calculations were performed with a previous version of SCALE and the 27burnplib cross section library. The concentrations calculated in Reference 5.5 for the most part agree with the concentrations calculated in this analysis; however, significant differences are seen for the plutonium isotopes. However, since there is a significant difference in the Path B model from Reference 5.5 and the model contained within, it cannot be determined what is responsible for the differences between calculated concentrations.

The assumption of a 10 day cooling time does not significantly effect isotopic concentrations. Since the concentration of ^{239}Pu is over-predicted for only some samples, it cannot be conclusively determined whether the assumed cooling time is inappropriate. However, the effects of such a short cooling time are minimal in comparison to the approximations of operating parameters.

The SAS2H code normally predicts isotopic concentrations as a radial assembly average; however, measurements are performed on individual pellet samples. Therefore, local pellet conditions are modeled as closely as possible in this analysis so that a more realistic pellet composition can be determined. However, approximations made to obtain local pellet conditions will influence the calculated isotopic concentrations.

In general, the SAS2H module of SCALE does well in predicting isotopic concentrations for samples from Trino Vercelles, using the methodology presented. The majority of calculated concentrations are close to the measured concentrations; however, more detailed operating data are expected to improve the accuracy of the calculated concentrations in relation to the corresponding measurements. It is recommended that future analyses use more detailed data if possible.

9. Attachments

Attachment I includes 14 pages and contains the input files used in the modeling of the Trino Vercelles samples. A description of the parameters contained within the input files is found in Sections 7.1 through 7.5.

Included in Attachment II is an extraction from each of the output files, containing the following information:

- echo of the SAS2H input deck,
- time/date stamp for when the SAS2H depletion calculation was performed,
- the output extraction of information pertinent to the Radiochemical Assay evaluations from the final ORIGEN calculation of the SAS2H depletion calculation.

t04m1lh7.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-104, rod m11 7, 12.042 gwd/mtu, June 97
```

```
mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1001
  92234 0.035 92235 3.897 92236 0.018 92238 96.050 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7795 1 543 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 543 end
```

```
1175 ppm boron (wt) in moderator
```

```
end comp
```

```
fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=10.390 burn=226 down=86 end
power=16.138 burn=263 down=51 bfrac=0.7234 end
power=18.663 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
```

```
end
```

t32e1h4.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-032, rod e11, level 4, 15.377 gwd/mtu, June 97
```

```
-----
mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1015
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-135 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7554 1 557 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 557 end
```

```
1175 ppm boron (wt) in moderator at start (1st segment)
```

```
-----
end comp
```

```
fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
-----
assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.268 burn=226 down=86 end
power=20.607 burn=263 down=51 bfrac=0.7234 end
power=23.831 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
```

```
-----
end
```

t32e1h7.input

=sas2h parm=skipshipdata
trino vercellese pwr, 509-032, rod e11 7, 15.898 gwd/mtu, June 97

mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 1001
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7795 1 543 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 543 end

1175 ppm boron (wt) in moderator

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.717 burn=226 down=86 end
power=21.309 burn=263 down=51 bfrac=0.7234 end
power=24.639 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

end

t32e11h9.input

=sas2h parm=skipshipdata

trino vercellese pwr, 509-032, rod e11 9, 11.529 gwd/mtu, June 97

* mixtures of fuel-pin-unit-cell:

44group latticecell

```

uo2 1 den=10.079 1 927
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 927 end
kr-85 1 0 1-20 927 end
y-89 1 0 1-20 927 end
sr-90 1 0 1-20 927 end
zr-93 1 0 1-20 927 end
zr-94 1 0 1-20 927 end
zr-95 1 0 1-20 927 end
nb-94 1 0 1-20 927 end
mo-95 1 0 1-20 927 end
tc-99 1 0 1-20 927 end
ru-101 1 0 1-20 927 end
ru-106 1 0 1-20 927 end
rh-103 1 0 1-20 927 end
rh-105 1 0 1-20 927 end
pd-105 1 0 1-20 927 end
pd-108 1 0 1-20 927 end
ag-109 1 0 1-20 927 end
sb-124 1 0 1-20 927 end
xe-131 1 0 1-20 927 end
xe-132 1 0 1-20 927 end
xe-135 1 0 1-20 927 end
xe-136 1 0 1-20 927 end
cs-134 1 0 1-20 927 end
cs-135 1 0 1-20 927 end
cs-137 1 0 1-20 927 end
ba-136 1 0 1-20 927 end
la-139 1 0 1-20 927 end
pr-141 1 0 1-20 927 end
pr-143 1 0 1-20 927 end
ce-144 1 0 1-20 927 end
nd-143 1 0 1-20 927 end
nd-145 1 0 1-20 927 end
nd-147 1 0 1-20 927 end
pm-147 1 0 1-20 927 end
pm-148 1 0 1-20 927 end
sm-147 1 0 1-20 927 end
sm-149 1 0 1-20 927 end
sm-150 1 0 1-20 927 end
sm-151 1 0 1-20 927 end
sm-152 1 0 1-20 927 end
eu-153 1 0 1-20 927 end
eu-154 1 0 1-20 927 end
eu-155 1 0 1-20 927 end
gd-155 1 0 1-20 927 end

```

ss304 2 1 570 end

h2o 3 den=0.7885 1 537 end

arbm-bormod 0.7885 1 1 0 0 5000 100 3 1175.0e-6 537 end

* 1175 ppm boron (wt) in moderator

end comp

* fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

* assembly and cycle parameters:

```

npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=9.947 burn=226 down=86 end
power=15.451 burn=263 down=51 bfrac=0.7234 end
power=17.868 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

```

end

t69e11h1.input

=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod e11, level 1, 12.859 gwd/mtu, June 97

mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 915
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 915 end
kr-85 1 0 1-20 915 end
y-89 1 0 1-20 915 end
sr-90 1 0 1-20 915 end
zr-93 1 0 1-20 915 end
zr-94 1 0 1-20 915 end
zr-95 1 0 1-20 915 end
nb-94 1 0 1-20 915 end
mo-95 1 0 1-20 915 end
tc-99 1 0 1-20 915 end
ru-101 1 0 1-20 915 end
ru-106 1 0 1-20 915 end
rh-103 1 0 1-20 915 end
rh-105 1 0 1-20 915 end
pd-105 1 0 1-20 915 end
pd-108 1 0 1-20 915 end
ag-109 1 0 1-20 915 end
sb-124 1 0 1-20 915 end
xe-131 1 0 1-20 915 end
xe-132 1 0 1-20 915 end
xe-135 1 0 1-20 915 end
xe-136 1 0 1-20 915 end
cs-134 1 0 1-20 915 end
cs-135 1 0 1-20 915 end
cs-137 1 0 1-20 915 end
ba-136 1 0 1-20 915 end
pr-141 1 0 1-20 915 end
pr-143 1 0 1-20 915 end
la-139 1 0 1-20 915 end
ce-144 1 0 1-20 915 end
nd-143 1 0 1-20 915 end
nd-145 1 0 1-20 915 end
nd-147 1 0 1-20 915 end
pm-147 1 0 1-20 915 end
pm-148 1 0 1-20 915 end
sm-147 1 0 1-20 915 end
sm-149 1 0 1-20 915 end
sm-150 1 0 1-20 915 end
sm-151 1 0 1-20 915 end
sm-152 1 0 1-20 915 end
eu-153 1 0 1-20 915 end
eu-154 1 0 1-20 915 end
eu-155 1 0 1-20 915 end
gd-155 1 0 1-20 915 end

ss304 2 1 570 end

h2o 3 den=0.7365 1 563 end
arbm-bormod 0.7365 1 1 0 0 5000 100 3 1175.0e-6 563 end

arbm-ag 10.159 1 1 0 1 47000 100 4 0.80 563 end
arbm-in 10.159 1 1 0 1 49000 100 4 0.15 563 end
arbm-cd 10.159 8 0 0 1
48108 0.85 48110 12.21 48111 12.63 48112 24.02
48113 12.27 48114 29.11 48601 1-20 48116 7.72
4 0.05 563 end

1175 ppm boron (wt) in moderator at start (1st segment)

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=5 mxrepeats=0
end

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

power=7.148 burn=226 down=86 end
power=11.101 burn=263 down=51 bfrac=0.7234 end
power=12.840 burn=292 down=1117 bfrac=0.2766 end
power=10.997 burn=416 down=10 bfrac=0.5532
h2ofrac=1.026 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

end

t69e11h2.input

=sas2h parm=skipshipdata

trino vercellese pwr, 509-069, rod e11, level 2, 20.602 gwd/mtu, June 97

mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 968

92234 0.028 92235 3.13 92236 0.014 92238 96.828 end

kr-83 1 0 1-20 968 end

kr-85 1 0 1-20 968 end

y-89 1 0 1-20 968 end

sr-90 1 0 1-20 968 end

zr-93 1 0 1-20 968 end

zr-94 1 0 1-20 968 end

zr-95 1 0 1-20 968 end

nb-94 1 0 1-20 968 end

mo-95 1 0 1-20 968 end

tc-99 1 0 1-20 968 end

ru-101 1 0 1-20 968 end

ru-106 1 0 1-20 968 end

rh-103 1 0 1-20 968 end

rh-105 1 0 1-20 968 end

pd-105 1 0 1-20 968 end

pd-108 1 0 1-20 968 end

ag-109 1 0 1-20 968 end

sb-124 1 0 1-20 968 end

xe-131 1 0 1-20 968 end

xe-132 1 0 1-20 968 end

xe-135 1 0 1-20 968 end

xe-136 1 0 1-20 968 end

cs-134 1 0 1-20 968 end

cs-135 1 0 1-20 968 end

cs-137 1 0 1-20 968 end

ba-136 1 0 1-20 968 end

la-139 1 0 1-20 968 end

pr-141 1 0 1-20 968 end

pr-143 1 0 1-20 968 end

ce-144 1 0 1-20 968 end

nd-143 1 0 1-20 968 end

nd-145 1 0 1-20 968 end

nd-147 1 0 1-20 968 end

pm-147 1 0 1-20 968 end

pm-148 1 0 1-20 968 end

sm-147 1 0 1-20 968 end

sm-149 1 0 1-20 968 end

sm-150 1 0 1-20 968 end

sm-151 1 0 1-20 968 end

sm-152 1 0 1-20 968 end

eu-153 1 0 1-20 968 end

eu-154 1 0 1-20 968 end

eu-155 1 0 1-20 968 end

gd-155 1 0 1-20 968 end

ss304 2 1 570 end

h2o 3 den=0.7407 1 561 end

arbm-bormod 0.7407 1 1 0 0 5000 100 3 1175.0e-6 561 end

arbm-ag 10.159 1 1 0 1 47000 100 4 0.80 561 end

arbm-in 10.159 1 1 0 1 49000 100 4 0.15 561 end

arbm-cd 10.159 8 0 0 1

48108 0.85 48110 12.21 48111 12.63 48112 24.02

48113 12.27 48114 29.11 48601 1-20 48116 7.72

4 0.05 561 end

1175 ppm boron (wt) in moderator at start (1st segment)

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=5 mxrepeats=0
end

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

power=11.452 burn=226 down=86 end

power=17.786 burn=263 down=51 bfrac=0.7234 end

power=20.571 burn=292 down=1117 bfrac=0.2766 end

power=17.619 burn=416 down=10 bfrac=0.5532

h2frac=1.025 end

o 135 cr 5.9 mn 0.33

fe 13.0 co 0.075 ni 9.9

zr 221.0 nb 0.71 sn 3.6

end

t69e11h4.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod e11, level 4, 23.718 gwd/mtu, June
97
```

```
-----
mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1015
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-135 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7554 1 553 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
```

```
1175 ppm boron (wt) in moderator at start (1st segment)
```

```
-----
end comp
-----
```

```
fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
-----
assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=4
nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.184 burn=226 down=86 end
power=20.476 burn=263 down=51 bfrac=0.7234
end
power=23.683 burn=292 down=1117 bfrac=0.2766
end
power=20.283 burn=416 down=10 bfrac=0.5532
h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
```

```
-----
end
```

t69e1 lh7.input

=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod e11, level 7, 24.304 gwd/mtu, June 97

' mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 1001
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7795 1 540 end
arbm-bormod 0.7795 1 1 0 5000 100 3 1175.0e-6 540 end

' 1175 ppm boron (wt) in moderator at start (1st segment)

' end comp

' fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

' assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inlevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.509 burn=226 down=86 end
power=20.982 burn=263 down=51 bfrac=0.7234 end
power=24.268 burn=292 down=1117 bfrac=0.2766 end
power=20.785 burn=416 down=10 bfrac=0.5532
h2ofrac=1.020 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

' end

t69e5h4.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod e5, level 4, 23.867 gwd/mtu, June 97
```

```
' mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1015
```

```
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
```

```
kr-83 1 0 1-20 1015 end
```

```
kr-85 1 0 1-20 1015 end
```

```
y-89 1 0 1-20 1015 end
```

```
sr-90 1 0 1-20 1015 end
```

```
zr-93 1 0 1-20 1015 end
```

```
zr-94 1 0 1-20 1015 end
```

```
zr-95 1 0 1-20 1015 end
```

```
nb-94 1 0 1-20 1015 end
```

```
mo-95 1 0 1-20 1015 end
```

```
tc-99 1 0 1-20 1015 end
```

```
ru-101 1 0 1-20 1015 end
```

```
ru-106 1 0 1-20 1015 end
```

```
rh-103 1 0 1-20 1015 end
```

```
rh-105 1 0 1-20 1015 end
```

```
pd-105 1 0 1-20 1015 end
```

```
pd-108 1 0 1-20 1015 end
```

```
ag-109 1 0 1-20 1015 end
```

```
sb-124 1 0 1-20 1015 end
```

```
xe-131 1 0 1-20 1015 end
```

```
xe-132 1 0 1-20 1015 end
```

```
xe-135 1 0 1-20 1015 end
```

```
xe-136 1 0 1-20 1015 end
```

```
cs-134 1 0 1-20 1015 end
```

```
cs-135 1 0 1-20 1015 end
```

```
cs-137 1 0 1-20 1015 end
```

```
ba-136 1 0 1-20 1015 end
```

```
la-139 1 0 1-20 1015 end
```

```
pr-141 1 0 1-20 1015 end
```

```
pr-143 1 0 1-20 1015 end
```

```
ce-144 1 0 1-20 1015 end
```

```
nd-143 1 0 1-20 1015 end
```

```
nd-145 1 0 1-20 1015 end
```

```
nd-147 1 0 1-20 1015 end
```

```
pm-147 1 0 1-20 1015 end
```

```
pm-148 1 0 1-20 1015 end
```

```
sm-147 1 0 1-20 1015 end
```

```
sm-149 1 0 1-20 1015 end
```

```
sm-150 1 0 1-20 1015 end
```

```
sm-151 1 0 1-20 1015 end
```

```
sm-152 1 0 1-20 1015 end
```

```
eu-153 1 0 1-20 1015 end
```

```
eu-154 1 0 1-20 1015 end
```

```
eu-155 1 0 1-20 1015 end
```

```
gd-155 1 0 1-20 1015 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7554 1 553 end
```

```
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
```

```
' 1175 ppm boron (wt) in moderator at start (1st segment)
```

```
end comp
```

```
' fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
```

```
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
```

```
3 0.7351 500 3.3248 2 3.3353
```

```
power=13.266 burn=226 down=86 end
```

```
power=20.605 burn=263 down=51 bfrac=0.7234 end
```

```
power=23.832 burn=292 down=1117 bfrac=0.2766 end
```

```
power=20.411 burn=416 down=10 bfrac=0.5532
```

```
h2ofrac=1.023 end
```

```
o 135 cr 5.9 mn 0.33
```

```
fe 13.0 co 0.075 ni 9.9
```

```
zr 221.0 nb 0.71 sn 3.6
```

```
end
```

t69e5h7.input

=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod e5, level 7, 24.548 gwd/mtu, June 97

' mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 1001
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7795 1 540 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 540 end

' 1175 ppm boron (wt) in moderator at start (1st segment)

end comp

' fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

' assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.645 burn=226 down=86 end
power=21.193 burn=263 down=51 bfrac=0.7234 end
power=24.512 burn=292 down=1117 bfrac=0.2766 end
power=20.993 burn=416 down=10 bfrac=0.5532
h2ofrac=1.020 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

end

t6911h4.input

=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod 111, level 4, 23.928 gwd/mtu, June 97

' mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 1015
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7554 1 553 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end

' 1175 ppm boron (wt) in moderator at start (1st segment)

end comp

' fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

' assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.300 burn=226 down=86 end
power=20.657 burn=263 down=51 bfrac=0.7234 end
power=23.893 burn=292 down=1117 bfrac=0.2766 end
power=20.463 burn=416 down=10 bfrac=0.5532
h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

end

t6911h7.input

sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod 111, level 7, 24.362 gwd/mtu, June 97

mixtures of fuel-pin-unit-cell:

44group latticecell

uo2 1 den=10.079 1 1001
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7795 1 540 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 540 end

1175 ppm boron (wt) in moderator at start (1st segment)

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inlevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.542 burn=226 down=86 end
power=21.032 burn=263 down=51 bfrac=0.7234 end
power=24.326 burn=292 down=1117 bfrac=0.2766 end
power=20.834 burn=416 down=10 bfrac=0.5532
h2ofrac=1.020 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

end

t6915h4.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod 15, level 4, 24.330 gwd/mtu, June 97
```

```
-----
' mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1015
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-135 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7554 1 553 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
```

```
' 1175 ppm boron (wt) in moderator at start (1st segment)
```

```
-----
end comp
```

```
-----
' fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
-----
' assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.524 burn=226 down=86 end
power=21.004 burn=263 down=51 bfrac=0.7234 end
power=24.294 burn=292 down=1117 bfrac=0.2766 end
power=20.807 burn=416 down=10 bfrac=0.5532
h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
```

```
-----
end
```

t6915h7.input

```
=sas2h parm=skipshipdata
trino vercellese pwr, 509-069, rod 15, level 7, 24.313 gwd/mtu, June 97
```

```
-----
' mixtures of fuel-pin-unit-cell:
```

```
44group latticecell
```

```
uo2 1 den=10.079 1 1001
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end
```

```
ss304 2 1 570 end
```

```
h2o 3 den=0.7795 1 540 end
arbm-bormod 0.7795 1 1 0 5000 100 3 1175.0e-6 540 end
```

```
' 1175 ppm boron (wt) in moderator at start (1st segment)
```

```
-----
end comp
```

```
-----
' fuel-pin-cell geometry:
```

```
squarepitch 1.303 0.9020 1 3 0.9786 2 end
```

```
-----
' assembly and cycle parameters:
```

```
npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.514 burn=226 down=86 end
power=20.990 burn=263 down=51 bfrac=0.7234 end
power=24.277 burn=292 down=1117 bfrac=0.2766 end
power=20.792 burn=416 down=10 bfrac=0.5532
h2ofrac=1.020 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
```

```
-----
end
```


t04m11h7.sum

```

0.....
|
|                SCALE4.3  Bulletin  Board
|                -----
|
|                Welcome to SCALE-4.3.
|
|.....

```

```

1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  crino vercellese pwr, 509-104, rod mll 7, 12.042 gwd/mtu, June 97

```

mixtures of fuel-pin-unit-cell:

44group latticecell

```

uo2 1 den=10.079 1 1001
    92234 0.035 92235 3.897 92236 0.018 92238 96.050 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

```

ss304 2 1 570 end

```

h2o 3 den=0.7795 1 543 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 543 end

```

1175 ppm boron (wt) in moderator

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

```

npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=10.390 burn=226 down=86 end
power=16.138 burn=263 down=51 bfrac=0.7234 end
power=18.663 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

```

```

1  SSSSSSSSSS      aaaaaaaaaa      SSSSSSSSSS      2222222222      hh      hh
   SSSSSSSSSSSS      aaaaaaaaaa      SSSSSSSSSS      2222222222      hh      hh
   ss      ss      aa      aa      ss      ss      22      22      hh      hh
   ss      ss      aa      aa      ss      ss      22      22      hh      hh
   ss      ss      aa      aa      ss      ss      22      22      hh      hh
   SSSSSSSSSSSS      aaaaaaaaaa      SSSSSSSSSS      22      hhhhhhhhhhhh
   SSSSSSSSSSSS      aaaaaaaaaa      SSSSSSSSSS      22      hhhhhhhhhhhh
   ss      ss      aa      aa      ss      ss      22      hh      hh
   ss      ss      aa      aa      ss      ss      22      hh      hh
   ss      ss      aa      aa      ss      ss      22      hh      hh
   SSSSSSSSSSSS      aa      aa      SSSSSSSSSS      2222222222      hh      hh
   SSSSSSSSSSSS      aa      aa      SSSSSSSSSS      2222222222      hh      hh

```

0

```

nn      nn      iiiiiiiiiiiii      cccccccccc      hh      hh      ooooooooooooo      11
nnnn    nn      iiiiiiiiiiiii      cccccccccc      hh      hh      ooooooooooooo      11
nnnnn   nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hhhhhhhhhhhhh      oo      oo      11
nn      nn      ii            cc      cc      hhhhhhhhhhhhh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      ii            cc      cc      hh      hh      oo      oo      11
nn      nn      iiiiiiiiiiiii      cccccccccc      hh      hh      ooooooooooooo      1111111111111111
nn      nn      iiiiiiiiiiiii      cccccccccc      hh      hh      ooooooooooooo      1111111111111111

```

0

```

00000000      77777777777777      //      33333333333      11      //      99999999999      77777777777777
0000000000    77777777777777      //      3333333333333      111      //      9999999999999      77777777777777
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
00      00      77      //      33      //      99      //
0000000000    77      //      3333333333333      11111111      //      9999999999999      77
00000000      77      //      33333333333      11111111      //      99999999999      77

```

0

```

11      00000000      44      6666666666666      33333333333      00000000
111      000000000      444      6666666666666      3333333333333      000000000
1111     00      4444      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11      00      44      66      //      33      00      00
11111111    000000000      4444444444444      66      //      33333333333      000000000
11111111    00000000      4444444444444      66      //      33333333333      00000000

```

1

0

```

ssssssssss      cccccccccc      11      eeeeeeeeeee
ssssssssssss    cccccccccc      aa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ssssssssssss    cc      aaaaaaaaaa      ee
ss      ss      cc      aaaaaaaaaa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ss      ss      cc      aa      ee
ssssssssssss    cccccccccc      aa      ee
ssssssssssss    cccccccccc      aa      ee

```

```

.....
program verification information
code system:  scale version:  4.3
.....

program:  sas2
creation date:  03/07/97
library:  /opt/neut/Scaled.3/bin

this is not a  scale configuration controlled code
jobname:  nichol
date of execution:  07/31/97
time of execution:  10:46:30
.....

```

1

0

```

initial 1E-18 d
o 16 1.35E+05 1.35E+05
total 3.90E+05 3.90E+05

```

0

```

initial 1E-18 d
u234 3.50E+02 3.50E+02
u235 3.90E+04 3.90E+04
u236 1.80E+02 1.80E+02
u238 9.60E+05 9.60E+05
total 1.00E+06 1.00E+06

```

0

0

0

```

basis =
initial 11.3 d 22.6 d 33.9 d 45.2 d 45.2 d
initial 11.3 d 22.6 d 33.9 d 45.2 d 45.2 d
nuclide concentrations, grams
basis =single reactor assembly
initial 1.7 d 3.3 d 5.0 d 6.7 d 8.3 d 10.0 d

```


t32e1lh4.sum

```

0.....
:
:          SCALE4.3  Bulletin  Board
:          -----
:
:          Welcome to SCALE-4.3.
:
:.....

```

```

1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-032, rod ell, level 4, 15.377 gwd/mtu, June 97

```

```

-----
: mixtures of fuel-pin-unit-cell:

```

```

44group latticecell

```

```

uo2 1 den=10.079 1 1015
    92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-135 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end

```

```

ss304 2 1 570 end

```

```

h2o 3 den=0.7554 1 557 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 557 end

```

```

: 1175 ppm boron (wt) in moderator at start (1st segment)

```

```

-----
: end comp

```

```

: fuel-pin-cell geometry:

```

```

squarepitch 1.303 0.9020 1 3 0.9786 2 end

```

```

-----
: assembly and cycle parameters:

```

```

npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.268 burn=226 down=86 end
power=20.607 burn=263 down=51 bfrac=0.7234 end
power=23.831 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

```

```

1 ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ssssssssss aa aa ssssssssss 2222222222 hh hh

```


				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
o 16	initial	1.7 d	3.3 d	1.35E+05	1.35E+05	1.35E+05	1.35E+05
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
mo 95	initial	1.7 d	3.3 d	7.18E+00	7.20E+00	7.22E+00	7.25E+00
tc 99				7.87E-06	7.88E-06	7.89E-06	7.90E-06
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
total	initial	1.7 d	3.3 d	3.90E+05	3.90E+05	3.90E+05	3.90E+05
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
u233	initial	1.7 d	3.3 d	1.51E-03	1.51E-03	1.51E-03	1.51E-03
u234				2.17E+02	2.17E+02	2.17E+02	2.17E+02
u235				1.77E+04	1.77E+04	1.77E+04	1.77E+04
u236				2.62E+03	2.62E+03	2.62E+03	2.62E+03
u238				9.56E+05	9.56E+05	9.56E+05	9.56E+05
np237				1.74E+02	1.75E+02	1.75E+02	1.77E+02
pu236				1.31E-04	1.32E-04	1.32E-04	1.31E-04
pu238				2.80E+01	2.81E+01	2.82E+01	2.83E+01
pu239				5.19E+03	5.21E+03	5.23E+03	5.24E+03
pu240				1.11E+03	1.11E+03	1.11E+03	1.11E+03
pu241				5.83E+02	5.83E+02	5.83E+02	5.82E+02
pu242				8.26E+01	8.26E+01	8.26E+01	8.26E+01
am241				1.45E+01	1.46E+01	1.48E+01	1.53E+01
am242m				2.67E-01	2.67E-01	2.67E-01	2.67E-01
am243				8.53E+00	8.55E+00	8.55E+00	8.55E+00
total				9.84E+05	9.84E+05	9.84E+05	9.84E+05
				element concentrations, grams			
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
mo 95	initial	1.7 d	3.3 d	3.03E+02	3.04E+02	3.05E+02	3.06E+02
tc 99				3.95E+02	3.96E+02	3.97E+02	3.97E+02
tc 99m				2.02E-01	1.45E-01	9.55E-02	6.27E-02
ru101				3.61E+02	3.61E+02	3.61E+02	3.61E+02
rh103				2.19E+02	2.20E+02	2.21E+02	2.23E+02
rh103m				2.84E-02	2.76E-02	2.68E-02	2.60E-02
ag109				3.19E+01	3.20E+01	3.20E+01	3.20E+01
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
nd143	initial	1.7 d	3.3 d	4.57E+02	4.59E+02	4.60E+02	4.61E+02
nd145				3.45E+02	3.45E+02	3.45E+02	3.45E+02
sml47				3.66E+01	3.68E+01	3.69E+01	3.71E+01
sml49				2.56E+00	2.89E+00	3.09E+00	3.20E+00
sml50				1.27E+02	1.27E+02	1.27E+02	1.27E+02
				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
sml51	initial	1.7 d	3.3 d	1.23E+01	1.24E+01	1.24E+01	1.24E+01
eul51				2.22E-02	2.26E-02	2.30E-02	2.35E-02
sml52				6.36E+01	6.36E+01	6.36E+01	6.36E+01
eul53				4.03E+01	4.05E+01	4.06E+01	4.07E+01
gd155				2.09E-02	2.21E-02	2.33E-02	2.45E-02
total				1.58E+04	1.58E+04	1.58E+04	1.58E+04

t32e11h7.sum

```

.....
.
.          SCALE4.3 Bulletin Board
.          -----
.          Welcome to SCALE-4.3.
.
.....

```

```

1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-032, rod ell 7, 15.898 gwd/mtu, June 97

```

mixtures of fuel-pin-unit-cell:

44group latticecell

```

uo2 1 den=10.079 1 1001
    92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

```

ss304 2 1 570 end

```

h2o 3 den=0.7795 1 543 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 543 end
  1175 ppm boron (wt) in moderator

```

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

```

npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.717 burn=226 down=86 end
power=21.309 burn=263 down=51 bfrac=0.7234 end
power=24.639 burn=292 down=10 bfrac=0.2766 end
  o 135 cr 5.9 mn 0.33
  fe 13.0 co 0.075 ni 9.9
  zr 221.0 nb 0.71 sn 3.6

```

```

1 ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ss ss aa aa ss ss 22 22 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ss ss aa aa ss ss 22 hh hh

```


t32e11h9.sum

```

.....
SCALE4.3 Bulletin Board
-----
Welcome to SCALE-4.3.
.....
1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-032, rod ell 9, 11.529 gwd/mtu, June 97

```

```

-----
mixtures of fuel-pin-unit-cell:
44group latticecell
uo2 1 den=10.079 1 927
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 927 end
kr-85 1 0 1-20 927 end
y-89 1 0 1-20 927 end
sr-90 1 0 1-20 927 end
zr-93 1 0 1-20 927 end
zr-94 1 0 1-20 927 end
zr-95 1 0 1-20 927 end
nb-94 1 0 1-20 927 end
mo-95 1 0 1-20 927 end
tc-99 1 0 1-20 927 end
ru-101 1 0 1-20 927 end
ru-106 1 0 1-20 927 end
rh-103 1 0 1-20 927 end
rh-105 1 0 1-20 927 end
pd-105 1 0 1-20 927 end
pd-108 1 0 1-20 927 end
ag-109 1 0 1-20 927 end
sb-124 1 0 1-20 927 end
xe-131 1 0 1-20 927 end
xe-132 1 0 1-20 927 end
xe-135 1 0 1-20 927 end
xe-136 1 0 1-20 927 end
cs-134 1 0 1-20 927 end
cs-135 1 0 1-20 927 end
cs-137 1 0 1-20 927 end
ba-136 1 0 1-20 927 end
la-139 1 0 1-20 927 end
pr-141 1 0 1-20 927 end
pr-143 1 0 1-20 927 end
ce-144 1 0 1-20 927 end
nd-143 1 0 1-20 927 end
nd-145 1 0 1-20 927 end
nd-147 1 0 1-20 927 end
pm-147 1 0 1-20 927 end
pm-148 1 0 1-20 927 end
sm-147 1 0 1-20 927 end
sm-149 1 0 1-20 927 end
sm-150 1 0 1-20 927 end
sm-151 1 0 1-20 927 end
sm-152 1 0 1-20 927 end
eu-153 1 0 1-20 927 end
eu-154 1 0 1-20 927 end
eu-155 1 0 1-20 927 end
gd-155 1 0 1-20 927 end

ss304 2 1 570 end

h2o 3 den=0.7885 1 537 end
arhm-bormod 0.7885 1 1 0 0 5000 100 3 1175.0e-6 537 end
: 1175 ppm boron (wt) in moderator
-----

```

```

end comp

fuel-pin-cell geometry:
squarepitch 1.303 0.9020 1 3 0.9786 2 end
-----

```

```

assembly and cycle parameters:
npin/assm=214 fuelnght=823.10 ncycles=3 nlib/cyc=5
printlevel=5 lightel=9 implevel=2 numxtotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=9.947 burn=226 down=86 end
power=15.451 burn=263 down=51 bfrac=0.7214 end
power=17.868 burn=292 down=10 bfrac=0.2766 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
-----

```

```

1  SSSSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSSSS  222222222222  hh  hh
SSSSSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSSSS  222222222222  hh  hh
SS  SS  AA  AA  SS  SS  22  22  hh  hh
SS  SS  AA  AA  SS  SS  22  22  hh  hh
SS  SS  AA  AA  SS  SS  22  22  hh  hh
SSSSSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSSSS  22  hhhhhhhhhhhh
SSSSSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSSSS  22  hhhhhhhhhhhh
SS  SS  AA  AA  SS  SS  22  hh  hh
SS  SS  AA  AA  SS  SS  22  hh  hh
SS  SS  AA  AA  SS  SS  22  hh  hh
SSSSSSSSSSSSS  AA  AA  SSSSSSSSSSSS  222222222222  hh  hh

```


				nuclide concentrations, grams			
				basis =single reactor assembly			
				5.0 d	6.7 d	8.3 d	10.0 d
0	o 16	initial	1.7 d 3.3 d	1.35E+05	1.35E+05	1.35E+05	1.35E+05
0							
0	mo 95	initial	1.7 d 3.3 d	5.15E+00	5.17E+00	5.19E+00	5.20E+00
0	tc 99			2.83E-06	2.84E-06	2.84E-06	2.84E-06
0							
0	total	initial	1.7 d 3.3 d	3.90E+05	3.90E+05	3.90E+05	3.90E+05
0							
0	u233	initial	1.7 d 3.3 d	1.21E-03	1.21E-03	1.21E-03	1.21E-03
0	u234			2.32E+02	2.32E+02	2.32E+02	2.32E+02
0	u235			2.03E+04	2.03E+04	2.03E+04	2.03E+04
0	u236			2.15E+03	2.15E+03	2.15E+03	2.15E+03
0	u238			9.59E+05	9.59E+05	9.59E+05	9.59E+05
0	np237			1.13E+02	1.13E+02	1.13E+02	1.13E+02
0	pu236			5.99E-05	5.99E-05	5.99E-05	5.99E-05
0	pu238			1.36E+01	1.37E+01	1.37E+01	1.38E+01
0	pu239			4.33E+03	4.35E+03	4.36E+03	4.37E+03
0	pu240			7.53E+02	7.53E+02	7.53E+02	7.53E+02
0	pu241			3.38E+02	3.38E+02	3.38E+02	3.38E+02
0	pu242			3.52E+01	3.52E+01	3.52E+01	3.52E+01
0	am241			8.47E+00	8.54E+00	8.62E+00	8.69E+00
0	am242m			1.36E-01	1.36E-01	1.36E-01	1.36E-01
0	am243			2.55E+00	2.56E+00	2.56E+00	2.56E+00
0	total			9.88E+05	9.88E+05	9.88E+05	9.88E+05
0							
0							
0	mo 95	initial	1.7 d 3.3 d	2.33E+02	2.33E+02	2.34E+02	2.35E+02
0							
0	tc 99	initial	1.7 d 3.3 d	3.03E+02	3.03E+02	3.03E+02	3.04E+02
0	tc 99m			1.52E-01	1.09E-01	7.20E-02	4.73E-02
0	ru101			2.71E+02	2.71E+02	2.71E+02	2.71E+02
0	rh103			1.65E+02	1.65E+02	1.66E+02	1.67E+02
0	rh103m			2.01E-02	1.95E-02	1.90E-02	1.84E-02
0	ag109			2.01E+01	2.02E+01	2.02E+01	2.02E+01
0							
0							
0	nd143	initial	1.7 d 3.3 d	3.60E+02	3.61E+02	3.62E+02	3.63E+02
0	nd145			2.67E+02	2.67E+02	2.67E+02	2.67E+02
0	sm147			3.01E+01	3.02E+01	3.03E+01	3.05E+01
0	sm149			2.26E+00	2.49E+00	2.63E+00	2.71E+00
0	sm150			9.14E+01	9.14E+01	9.14E+01	9.14E+01
0	sm151			1.04E+01	1.04E+01	1.05E+01	1.05E+01
0							
0							
0	eu151	initial	1.7 d 3.3 d	2.50E-02	2.54E-02	2.58E-02	2.61E-02
0	sm152			4.75E+01	4.75E+01	4.75E+01	4.75E+01
0	eu153			2.63E+01	2.64E+01	2.64E+01	2.64E+01
0	gd155			1.71E-02	1.79E-02	1.87E-02	1.95E-02
0	total			1.19E+04	1.19E+04	1.19E+04	1.19E+04

t69el1hl.sum

```

0.....
*
*          SCALE4.3  Bulletin Board
*          -----
*          Welcome to SCALE-4.3.
*
*.....
1  primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
-  module sas2h will be called
   trino vercellese pwr, 509-069, rod ell, level 1, 12.859 gwd/mtu, June 97
   .
   .-----
   . mixtures of fuel-pin-unit-cell:
   .
   .44group latticecell
   .
   uo2 1 den=10.079 1 915
     92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
   kr-83 1 0 1-20 915 end
   kr-85 1 0 1-20 915 end
   y-89 1 0 1-20 915 end
   sr-90 1 0 1-20 915 end
   zr-93 1 0 1-20 915 end
   zr-94 1 0 1-20 915 end
   zr-95 1 0 1-20 915 end
   nb-94 1 0 1-20 915 end
   mo-95 1 0 1-20 915 end
   tc-99 1 0 1-20 915 end
   ru-101 1 0 1-20 915 end
   ru-106 1 0 1-20 915 end
   rh-103 1 0 1-20 915 end
   rh-105 1 0 1-20 915 end
   pd-105 1 0 1-20 915 end
   pd-108 1 0 1-20 915 end
   ag-109 1 0 1-20 915 end
   sb-124 1 0 1-20 915 end
   xe-131 1 0 1-20 915 end
   xe-132 1 0 1-20 915 end
   xe-135 1 0 1-20 915 end
   xe-136 1 0 1-20 915 end
   cs-134 1 0 1-20 915 end
   cs-135 1 0 1-20 915 end
   cs-137 1 0 1-20 915 end
   ba-136 1 0 1-20 915 end
   pr-141 1 0 1-20 915 end
   pr-143 1 0 1-20 915 end
   la-139 1 0 1-20 915 end
   ce-144 1 0 1-20 915 end
   nd-143 1 0 1-20 915 end
   nd-145 1 0 1-20 915 end
   nd-147 1 0 1-20 915 end
   pm-147 1 0 1-20 915 end
   pm-148 1 0 1-20 915 end
   sm-147 1 0 1-20 915 end
   sm-149 1 0 1-20 915 end
   sm-150 1 0 1-20 915 end
   sm-151 1 0 1-20 915 end
   sm-152 1 0 1-20 915 end
   eu-153 1 0 1-20 915 end
   eu-154 1 0 1-20 915 end
   eu-155 1 0 1-20 915 end
   qd-155 1 0 1-20 915 end

   ss304 2 1 570 end

   h2o 3 den=0.7365 1 563 end
   arbm-bormod 0.7365 1 1 0 0 5000 100 3 1175.0e-6 563 end

   arbm-ag 10.159 1 1 0 1 47000 100 4 0.80 563 end
   arbm-in 10.159 1 1 0 1 49000 100 4 0.15 563 end
   arbm-cd 10.159 8 0 0 1
   48108 0.85 48110 12.21 48111 12.63 48112 24.02
   48113 12.27 48114 29.11 48601 1-20 48116 7.72
   4 0.05 563 end

   .
   . 1175 ppm boron (wt) in moderator at start (1st segment)
   .-----
end comp

   .-----
   . fuel-pin-cell geometry:
   .
   squarepitch 1.303 0.9020 1 3 0.9786 2 end

   .-----
   . assembly and cycle parameters:
   .
   npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
   printlevel=5 lightel=9 implevel=2 numtotal=5 mxrepeats=0 end
   4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
   4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
   4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
   4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
   4 0.4573 2 0.5005 3 0.7351 500 3.2076 2 3.2177
   .
   3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
   3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
   3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
   3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353
   3 0.4573 3 0.5005 3 0.7351 500 3.3248 2 3.3353

```


t69e11h4.sum

```

0.....
  |
  |          SCALE4.3  Bulletin  Board
  |          -----
  |
  |          Welcome to SCALE-4.3.
  |
  |.....

```

```

1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-069, rod ell, level 4, 23.718 gwd/mtu, June 97

```

mixtures of fuel-pin-unit-cell:

44group latticecell

```

uo2 1 den=10.079 1 1015
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
cs-136 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end

ss304 2 1 570 end

```

```

h2o 3 den=0.7554 1 553 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
  1175 ppm boron (wt) in moderator at start (1st segment)
-----

```

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

```

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lighttel=9 implevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.184 burn=226 down=86 end
power=20.476 burn=263 down=51 bfrac=0.7234 end
power=23.683 burn=292 down=1117 bfrac=0.2766 end
power=20.283 burn=416 down=10 bfrac=0.5532 h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 an 3.6

```

```

1 #####          aaaaaaaaaa          #####          222222222222  hh          hh
#####          aaaaaaaaaa          #####          222222222222  hh          hh
SS          SS          aa          aa          SS          SS          22          22          hh          hh
SS          aa          aa          aa          SS          SS          22          22          hh          hh
SS          aa          aa          aa          SS          SS          22          22          hh          hh
#####          aaaaaaaaaa          #####          22          hh          hh          hh          hh
#####          aaaaaaaaaa          #####          22          hh          hh          hh          hh
SS          aa          aa          aa          SS          SS          22          hh          hh          hh
SS          aa          aa          aa          SS          SS          22          hh          hh          hh

```


t69e11h7.sum

```

0.....
:
:          SCALE4.3 Bulletin Board
:          -----
:
:          Welcome to SCALE-4.3.
:
:.....
1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr. 509-069, rod e11, level 7, 24.304 gwd/mtu, June 97

```

mixtures of fuel-pin-unit-cell:

44group latticecell

```

uo2 1 den=10.079 1 1001
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
qd-155 1 0 1-20 1001 end

```

ss304 2 1 570 end

```

h2o 3 den=0.7795 1 540 end
arpm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 540 end
  1175 ppm boron (wt) in moderator at start (1st segment)

```

end comp

fuel-pin-cell geometry:

squarepitch 1.303 0.9020 1 3 0.9786 2 end

assembly and cycle parameters:

```

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numxtotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.509 burn=226 down=86 end
power=20.982 burn=263 down=51 bfrac=0.7234 end
power=24.268 burn=292 down=1117 bfrac=0.2766 end
power=20.785 burn=416 down=10 bfrac=0.5532 h2ofrac=1.020 end
  o 135 cr 5.9 mn 0.33
  fe 13.0 co 0.075 ni 9.9
  zr 221.0 nb 0.71 sn 3.6

```

```

1 sssssssssss aaaaaaaaaa sssssssssss 2222222222 hh hh
  sssssssssss aaaaaaaaaa sssssssssss 2222222222 hh hh
  ss aa aa aa ss 22 hh hh
  ss aa aa aa ss 22 hh hh
  ss aa aa aa ss 22 hh hh
  sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
  sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh

```


t69e5h4.sum

```

.....
*
*          SCALE4.3  Bulletin  Board
*          -----
*          Welcome to SCALE-4.3.
*
.....
1  primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
-  module sas2h will be called
  trino vercellese pwr, 509-069, rod e5, level 4, 23.867 gwd/mtu, June 97
  mixtures of fuel-pin-unit-cell:
  44group latticecell
  uo2 1 den=10.079 1 1015
      92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
  kr-83 1 0 1-20 1015 end
  kr-85 1 0 1-20 1015 end
  y-89 1 0 1-20 1015 end
  sr-90 1 0 1-20 1015 end
  zr-93 1 0 1-20 1015 end
  zr-94 1 0 1-20 1015 end
  zr-95 1 0 1-20 1015 end
  nb-94 1 0 1-20 1015 end
  mo-95 1 0 1-20 1015 end
  tc-99 1 0 1-20 1015 end
  ru-101 1 0 1-20 1015 end
  ru-106 1 0 1-20 1015 end
  rh-103 1 0 1-20 1015 end
  rh-105 1 0 1-20 1015 end
  pd-105 1 0 1-20 1015 end
  pd-108 1 0 1-20 1015 end
  ag-109 1 0 1-20 1015 end
  sb-124 1 0 1-20 1015 end
  xe-131 1 0 1-20 1015 end
  xe-132 1 0 1-20 1015 end
  xe-135 1 0 1-20 1015 end
  xe-136 1 0 1-20 1015 end
  cs-134 1 0 1-20 1015 end
  cs-135 1 0 1-20 1015 end
  cs-137 1 0 1-20 1015 end
  ba-136 1 0 1-20 1015 end
  la-139 1 0 1-20 1015 end
  pr-141 1 0 1-20 1015 end
  pr-143 1 0 1-20 1015 end
  ce-144 1 0 1-20 1015 end
  nd-143 1 0 1-20 1015 end
  nd-145 1 0 1-20 1015 end
  nd-147 1 0 1-20 1015 end
  pm-147 1 0 1-20 1015 end
  pm-148 1 0 1-20 1015 end
  sm-147 1 0 1-20 1015 end
  sm-149 1 0 1-20 1015 end
  sm-150 1 0 1-20 1015 end
  sm-151 1 0 1-20 1015 end
  sm-152 1 0 1-20 1015 end
  eu-153 1 0 1-20 1015 end
  eu-154 1 0 1-20 1015 end
  eu-155 1 0 1-20 1015 end
  gd-155 1 0 1-20 1015 end
  ss304 2 1 570 end
  h2o 3 den=0.7554 1 553 end
  arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
  1175 ppm boron (wt) in moderator at start (1st segment)
  -----
  end comp
  -----
  fuel-pin-cell geometry:
  squarepitch 1.303 0.9020 1 3 0.9786 2 end
  -----
  assembly and cycle parameters:
  npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
  printlevel=5 lightel=9 inplevel=2 numxtotal=3 end
  3 0.7351 500 3.3248 2 3.3353
  power=13.266 burn=226 down=86 end
  power=20.605 burn=263 down=51 bfrac=0.7234 end
  power=23.832 burn=292 down=1117 bfrac=0.2766 end
  power=20.411 burn=416 down=10 bfrac=0.5532 h2ofrac=1.023 end
  o 135 cr 5.9 mn 0.33
  fe 13.0 co 0.075 ni 9.9
  zr 221.0 nb 0.71 sn 3.6
  -----
1  sssssssssss aaaaaaaaaa sssssssssss 2222222222 hh hh
  sssssssssss aaaaaaaaaa sssssssssss 222222222222 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
  sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh
  sssssssssss aa aa sssssssssss 222222222222 hh hh

```


t69e5h7.sum

```

0.....
:
:          SCALE4.3  Bulletin  Board
:          -----
:
:          Welcome to SCALE-4.3.
:
:.....
1  primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
-  module sas2h will be called
   trino vercellese pwr, 509-069, rod e5, level 7, 24.548 gwd/mtu, June 97

```

mixtures of fuel-pin-unit-cell:

```

44group latticecell
uo2 1 den=10.079 1 1001
92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 i 570 end

```

```

h2o 3 den=0.7795 1 540 end
arbm-bormod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 540 end
1175 ppm boron (wt) in moderator at start (1st segment)

```

end comp

fuel-pin-cell geometry:

```

squarepitch 1.303 0.9020 1 3 0.9786 2 end

```

assembly and cycle parameters:

```

npin/assm=214 fuelnigt=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 implevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.645 burn=226 down=86 end
power=21.193 burn=263 down=51 bfrac=0.7234 end
power=24.512 burn=292 down=1117 bfrac=0.2766 end
power=20.993 burn=416 down=10 bfrac=0.5532 h2ofrac=1.020 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6

```

```

1  sssssssssss aaaaaaaaaa sssssssssss 2222222222 hh hh
   sssssssssss aaaaaaaaaa sssssssssss 2222222222 hh hh
ss  ss aa aaaa aa ss ss 22 22 hh hh
ss  ss aa aa aa ss ss 22 22 hh hh
ss  ss aa aa aa ss ss 22 22 hh hh
   sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
   sssssssssss aaaaaaaaaa sssssssssss 22 hhhhhhhhhhhh
   ss aa aa ss 22 hh hh
   ss aa aa ss 22 hh hh

```


t69111h4.sum

```

0.....
:
:          SCALE4.3  Bulletin  Board
:          -----
:
:          Welcome to SCALE-4.3.
:
:.....
1  primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
-  module sas2h will be called
   trino vercellese pwr, 509-069, rod 111, level 4, 23.928 gwd/mtu, June 97

```

```

-----
: mixtures of fuel-pin-unit-cell:
:
44group latticecell
:
uo2 1 den=10.079 1 1015
   92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
gd-155 1 0 1-20 1001 end

ss304 2 1 570 end

h2o 3 den=0.7554 1 553 end
arbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
: 1175 ppm boron (wt) in moderator at start (1st segment)
:-----
end comp

```

```

-----
: fuel-pin-cell geometry:
squarepitch 1.303 0.9020 1 3 0.9786 2 end
:-----

```

```

-----
: assembly and cycle parameters:
npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 implevel=2 numxtotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=11.300 burn=226 down=86 end
power=20.657 burn=263 down=51 bfrac=0.7234 end
power=23.893 burn=292 down=1117 bfrac=0.2766 end
power=20.463 burn=416 down=10 bfrac=0.5532 h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
:-----

```

```

1  ssssssssss  aaaaaaaaaa  ssssssssss  2222222222  hh  hh
   ssssssssss  aaaaaaaaaa  ssssssssss  222222222222  hh  hh
ss  ss  aa  aa  aa  ss  ss  22  22  hh  hh
ss  aa  aa  aa  ss  ss  22  22  hh  hh
ss  aa  aa  aa  ss  ss  22  22  hh  hh
   ssssssssss  aaaaaaaaaa  ssssssssss  22  hhhhhhhhhhhh
   ssssssssss  aaaaaaaaaa  ssssssssss  22  hhhhhhhhhhhh
ss  aa  aa  ss  22  hh  hh
ss  aa  aa  ss  22  hh  hh

```


t6915h4.sum

```

.....
SCALE4.3 Bulletin Board
-----
Welcome to SCALE-4.3.
.....
1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-069, rod 15, level 4, 24.330 gwd/mtu, June 97

```

```

-----
mixtures of fuel-pin-unit-cell:
44group latticecell
uo2 1 den=10.079 1 1015
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1015 end
kr-85 1 0 1-20 1015 end
y-89 1 0 1-20 1015 end
sr-90 1 0 1-20 1015 end
zr-93 1 0 1-20 1015 end
zr-94 1 0 1-20 1015 end
zr-95 1 0 1-20 1015 end
nb-94 1 0 1-20 1015 end
mo-95 1 0 1-20 1015 end
tc-99 1 0 1-20 1015 end
ru-101 1 0 1-20 1015 end
ru-106 1 0 1-20 1015 end
rh-103 1 0 1-20 1015 end
rh-105 1 0 1-20 1015 end
pd-105 1 0 1-20 1015 end
pd-108 1 0 1-20 1015 end
ag-109 1 0 1-20 1015 end
sb-124 1 0 1-20 1015 end
xe-131 1 0 1-20 1015 end
xe-132 1 0 1-20 1015 end
xe-135 1 0 1-20 1015 end
xe-136 1 0 1-20 1015 end
cs-134 1 0 1-20 1015 end
cs-135 1 0 1-20 1015 end
cs-137 1 0 1-20 1015 end
ba-136 1 0 1-20 1015 end
la-139 1 0 1-20 1015 end
pr-141 1 0 1-20 1015 end
pr-143 1 0 1-20 1015 end
ce-144 1 0 1-20 1015 end
nd-143 1 0 1-20 1015 end
nd-145 1 0 1-20 1015 end
nd-147 1 0 1-20 1015 end
pm-147 1 0 1-20 1015 end
pm-148 1 0 1-20 1015 end
sm-147 1 0 1-20 1015 end
sm-149 1 0 1-20 1015 end
sm-150 1 0 1-20 1015 end
sm-151 1 0 1-20 1015 end
sm-152 1 0 1-20 1015 end
eu-153 1 0 1-20 1015 end
eu-154 1 0 1-20 1015 end
eu-155 1 0 1-20 1015 end
gd-155 1 0 1-20 1015 end
ss304 2 1 570 end
h2o 3 den=0.7554 1 553 end
axbm-bormod 0.7554 1 1 0 0 5000 100 3 1175.0e-6 553 end
1175 ppm boron (wt) in moderator at start (1st segment)
-----

```

```

end comp
-----
fuel-pin-cell geometry:
squarepitch 1.303 0.9020 1 3 0.9786 2 end
-----

```

```

assembly and cycle parameters:
npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numtotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.524 burn=226 down=86 end
power=21.004 burn=263 down=51 bfrac=0.7234 end
power=24.294 burn=292 down=1117 bfrac=0.2766 end
power=20.807 burn=416 down=10 bfrac=0.5532 h2ofrac=1.023 end
o 135 cr 5.9 mn 0.33
fe 13.0 co 0.075 ni 9.9
zr 221.0 nb 0.71 sn 3.6
-----

```

```

1  SSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSS  2222222222  hh  hh
  SSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSS  222222222222  hh  hh
  SS  SS  AA  AA  SS  SS  22  22  hh  hh
  SS  SS  AA  AA  SS  SS  22  22  hh  hh
  SS  SS  AA  AA  SS  SS  22  22  hh  hh
  SSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSS  22  hhhhhhhhhhhh
  SSSSSSSSSS  AAAAAAAAAA  SSSSSSSSSS  22  hhhhhhhhhhhh
  SS  SS  AA  AA  SS  SS  22  hh  hh
  SS  SS  AA  AA  SS  SS  22  hh  hh

```


t6915h7.sum

```

.....
          SCALE4.3 Bulletin Board
          -----
          Welcome to SCALE-4.3.
.....

```

```

1 primary module access and input record ( scale driver - 95/03/29 - 09:06:37 )
- module sas2h will be called
  trino vercellese pwr, 509-069, rod 15, level 7, 24.313 gwd/mtu, June 97

```

```

-----
mixtures of fuel-pin-unit-cell:

```

```

44group latticecell

```

```

uo2 1 den=10.079 1 1001
  92234 0.028 92235 3.13 92236 0.014 92238 96.828 end
kr-83 1 0 1-20 1001 end
kr-85 1 0 1-20 1001 end
y-89 1 0 1-20 1001 end
sr-90 1 0 1-20 1001 end
zr-93 1 0 1-20 1001 end
zr-94 1 0 1-20 1001 end
zr-95 1 0 1-20 1001 end
nb-94 1 0 1-20 1001 end
mo-95 1 0 1-20 1001 end
tc-99 1 0 1-20 1001 end
ru-101 1 0 1-20 1001 end
ru-106 1 0 1-20 1001 end
rh-103 1 0 1-20 1001 end
rh-105 1 0 1-20 1001 end
pd-105 1 0 1-20 1001 end
pd-108 1 0 1-20 1001 end
ag-109 1 0 1-20 1001 end
sb-124 1 0 1-20 1001 end
xe-131 1 0 1-20 1001 end
xe-132 1 0 1-20 1001 end
xe-135 1 0 1-20 1001 end
xe-136 1 0 1-20 1001 end
cs-134 1 0 1-20 1001 end
cs-135 1 0 1-20 1001 end
cs-137 1 0 1-20 1001 end
ba-136 1 0 1-20 1001 end
la-139 1 0 1-20 1001 end
pr-141 1 0 1-20 1001 end
pr-143 1 0 1-20 1001 end
ce-144 1 0 1-20 1001 end
nd-143 1 0 1-20 1001 end
nd-145 1 0 1-20 1001 end
nd-147 1 0 1-20 1001 end
pm-147 1 0 1-20 1001 end
pm-148 1 0 1-20 1001 end
sm-147 1 0 1-20 1001 end
sm-149 1 0 1-20 1001 end
sm-150 1 0 1-20 1001 end
sm-151 1 0 1-20 1001 end
sm-152 1 0 1-20 1001 end
eu-153 1 0 1-20 1001 end
eu-154 1 0 1-20 1001 end
eu-155 1 0 1-20 1001 end
qd-155 1 0 1-20 1001 end

```

```

ss304 2 1 570 end
h2o 3 den=0.7795 1 540 end
arpm-boromod 0.7795 1 1 0 0 5000 100 3 1175.0e-6 540 end
  1175 ppm boron (wt) in moderator at start (1st segment)

```

```

end comp

```

```

-----
fuel-pin-cell geometry:

```

```

squarepitch 1.303 0.9020 1 3 0.9786 2 end

```

```

-----
assembly and cycle parameters:

```

```

npin/assm=214 fuelnght=823.10 ncycles=4 nlib/cyc=5
printlevel=5 lightel=9 inplevel=2 numztotal=3 end
3 0.7351 500 3.3248 2 3.3353
power=13.514 burn=226 down=86 end
power=20.990 burn=263 down=51 bfrac=0.7234 end
power=24.277 burn=292 down=1117 bfrac=0.2766 end
power=20.792 burn=416 down=10 bfrac=0.5532 h2ofrac=1.020 end
  o 135 cr 5.9 mm 0.33
  fe 13.0 co 0.075 ni 9.9
  zr 221.0 nb 0.71 sn 3.6

```

```

1 ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 2222222222 hh hh
  ss aa aa ss 22 22 hh hh
  ss aa aa ss 22 22 hh hh
  ss aa aa ss 22 22 hh hh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ssssssssss aaaaaaaaaa ssssssssss 22 hhhhhhhhhhhh
  ss aa aa ss 22 hh hh
  ss aa aa ss 22 hh hh

```