

## General description for SHETRAN code and Benchmark problem calculation methods

### 1. Code name

SHETRAN

#### References:

K. Sakai, et al., "Development of Ultra Fine Energy Spectrum Analysis Code for PWR Assembly," Proc. of Autumn meeting, the Atomic Energy Society of Japan, F47, (1997), [in Japanese]

K. Sakai, et al., "Improvement of SEPCO Core Management Code System," Proc. of Annual meeting, the Atomic Energy Society of Japan, C6, (2000), [in Japanese]

### 2. Organization

Shikoku Electric Power Co., Inc. (SEPCO)

### 3. Nuclear data library

ENDF/B-VI rev.3

### 4. Code used for cross section data processing

NJOY94.105

### 5. Spectrum calculation and energy condensation methods

The numerical solution method for the transport equation in heterogeneous geometries is based on the method of characteristics. Multi-group constants are used for the fast energy region and the thermal energy region, and ultra-fine-group constants are used for the resonance energy region. One rod cell and 3x3 rod cells are calculated in 1/8 two-dimensional heterogeneous geometry. A non-fuel cell (for example: a water hole or a burnable poison rod) is assigned at center and surrounded by fuel rods in 3x3 rod cells geometry.

70 energy group constants are generated by condensing the flux weighted multi-group constants and ultra-fine-group constants.

For assembly geometry, 70 energy group transport equations are solved by the method of characteristics in heterogeneous system.

## 6. Energy groups

- Group constants calculation:

Fast energy region	56 groups (10MeV – about 10keV)
Resonance energy region	about 60 thousands groups (about 10keV – 2.1eV)
Thermal energy region	160 groups (2.1eV – 1.0E-5eV)
Cut-off energy is 2.1eV.	

- Assembly calculation: 70 groups (10MeV – 1.0E-5eV)

## 7. Assumption for cross section generation

Pseudo-FPs are not assumed.

## 8. Nuclides self-shielding and mutual shielding considered

For resonance energy region, all nuclides are accurately treated by ultra-fine-group cross sections. For fast and thermal energy region, self-shielding is considered by NR approximation.

## 9. Cell calculation geometry

Mesh number for Gd fuel pin : The number of radial division is 10.

The number of azimuthal division is 4.

## 10. Miscellaneous

- Burn-up steps (GWd/t)

UO<sub>2</sub> pin cell, MOX pin cell, MOX assembly:

0,0.1,0.5,1,2,4,5,6,8,10, and every 2.5 for 10-70 (total 34 steps)

Gd assembly:

0,0.1,0.25, every 0.25 for 0.25–30, and every 2.5 for 30–70 (total 138 steps)

- Burn-up chain

101 nuclides (18 actinides, 83 FP nuclides) are considered.