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The development and preliminary V&V of the MOC transport module in new lattice code LATC

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Introduction

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Summary



Introduction of SNPSDC

- Founded in May 2010, a subsidiary of SNPTC
- Take the leading role in the **State Energy Key Lab of NPP software** (proved by DOE of China in Sep, 2011)
- The **sole** state energy key lab specialized in self-reliance NPP
- >80 staff and 13 in reactor physics division



01

**NPP design codes
self-reliance R&D**

02

**Management, maintenance
& technical support for the
3rd generation of NPP
software**

Main Business

03

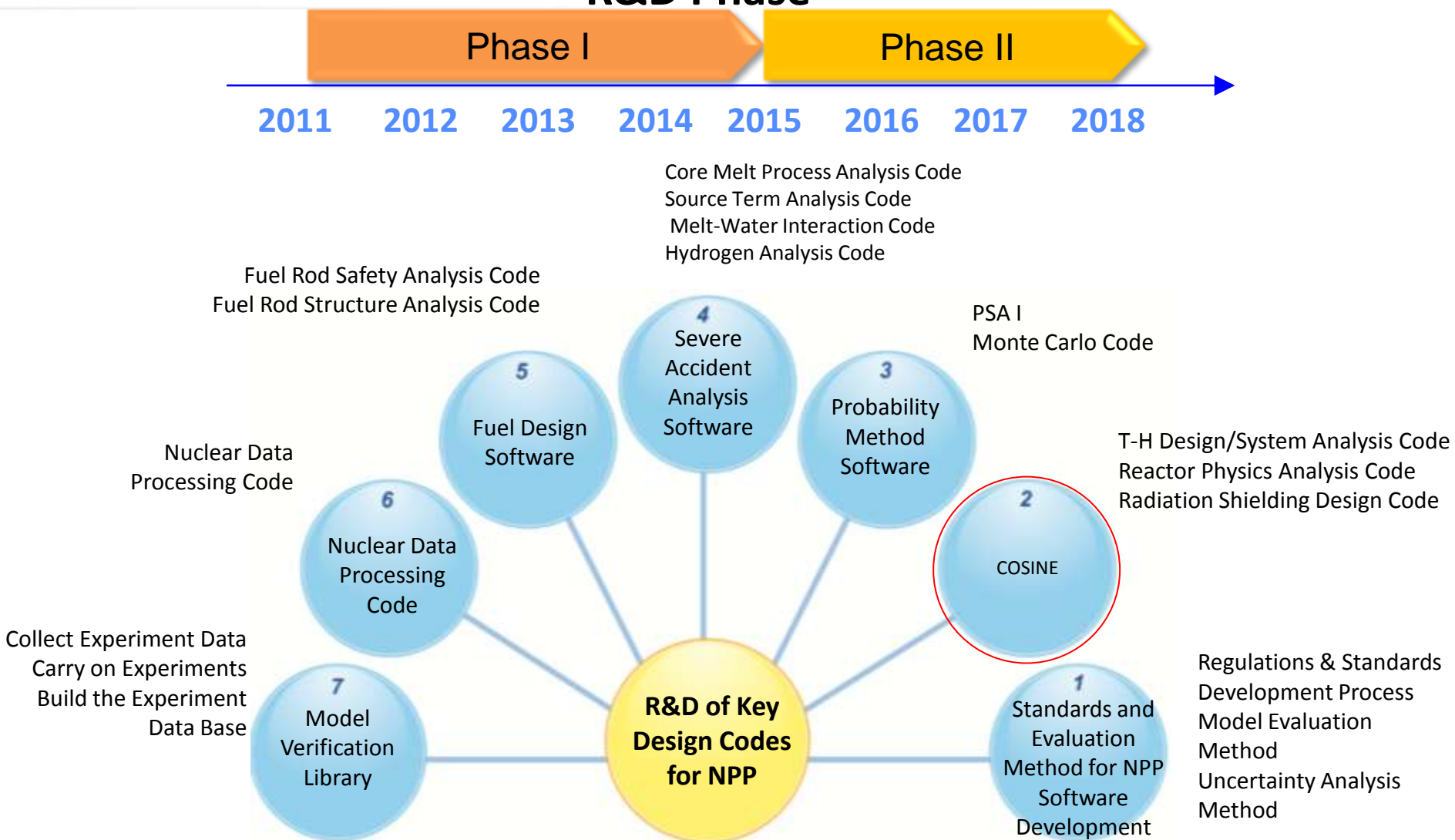
**V&V technologies
research for NPP
software**

04

**The research on engineering
application of advanced
simulation technologies**

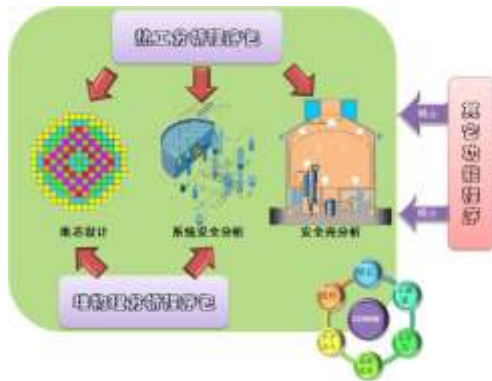


R&D Phase



COSINE

***C**Ore and **S**ystem **I**ntegrated **E**ngine
for design and analysis*



- ✓ An integrated platform
- ✓ Modular code system
- ✓ Multi-physics, multi-scale
- ✓ Sensitivity/Uncertainty analysis
- ✓ User friendly GUIs
- ✓ Parallel computations
- ✓ Software engineering QA
- ✓

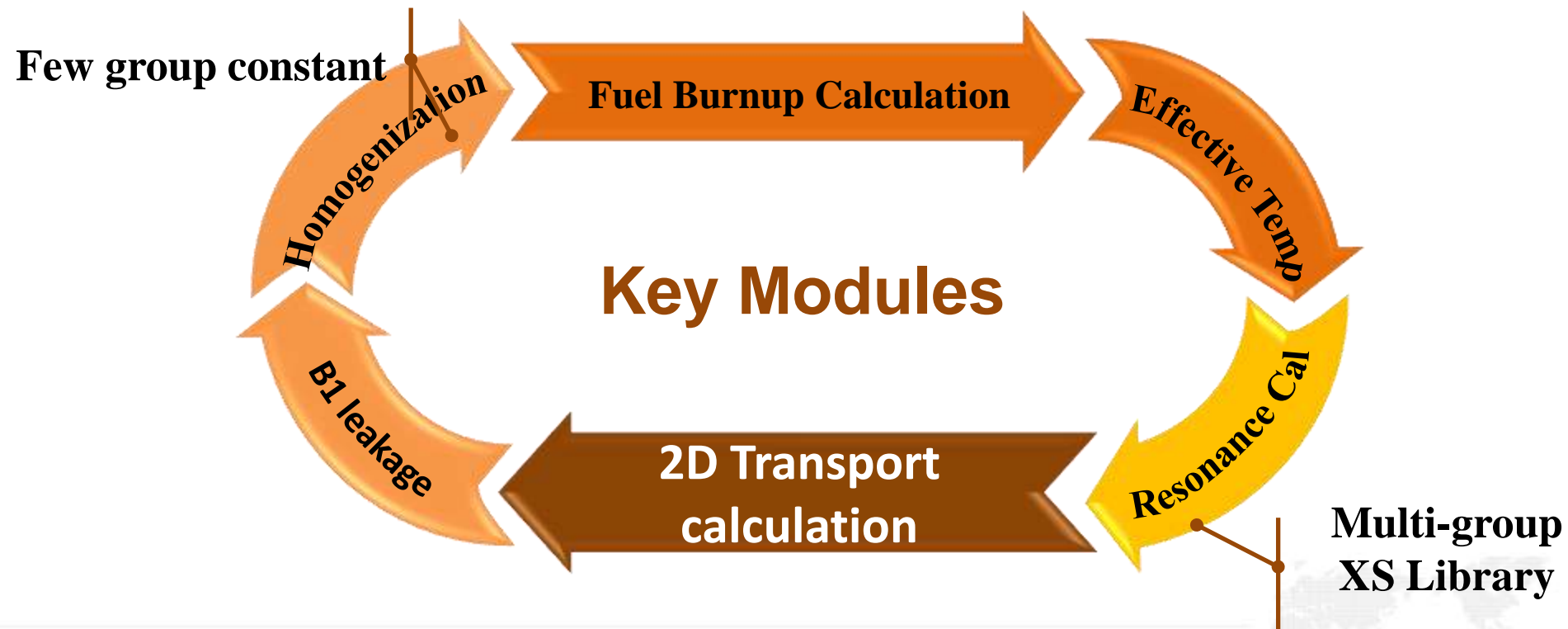
THERMAL CODE

- Subchannel code(SUBC)
- System code (SYST)
- Containment code(CONT)

PHYSICS CODE

- ***Lattice code(LATC)***
- Core code (CORE)
- Kinetics code(KIND)

- ✓ A multi-group two-dimensional lattice transport code.
- ✓ Mainly used on PWR assemblies or pin cells.
- ✓ Capable of satisfying most of the needs for PWR lattice analysis.

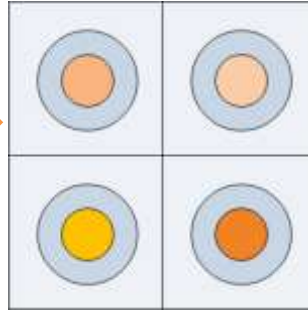


- Present
 - WIMSD-IAEA 69/172
 - Published by IAEA in WLUP
 - Adding the burn-up data of Wolfram isotope

- Future
 - A more accuracy library on develop
 - Energy Structure: SHEM281

Total Group No.	Total Nuclide No.	Fission Product No.	Resonance Nuclide No.
281	176	85	20

Temperature
Distribution



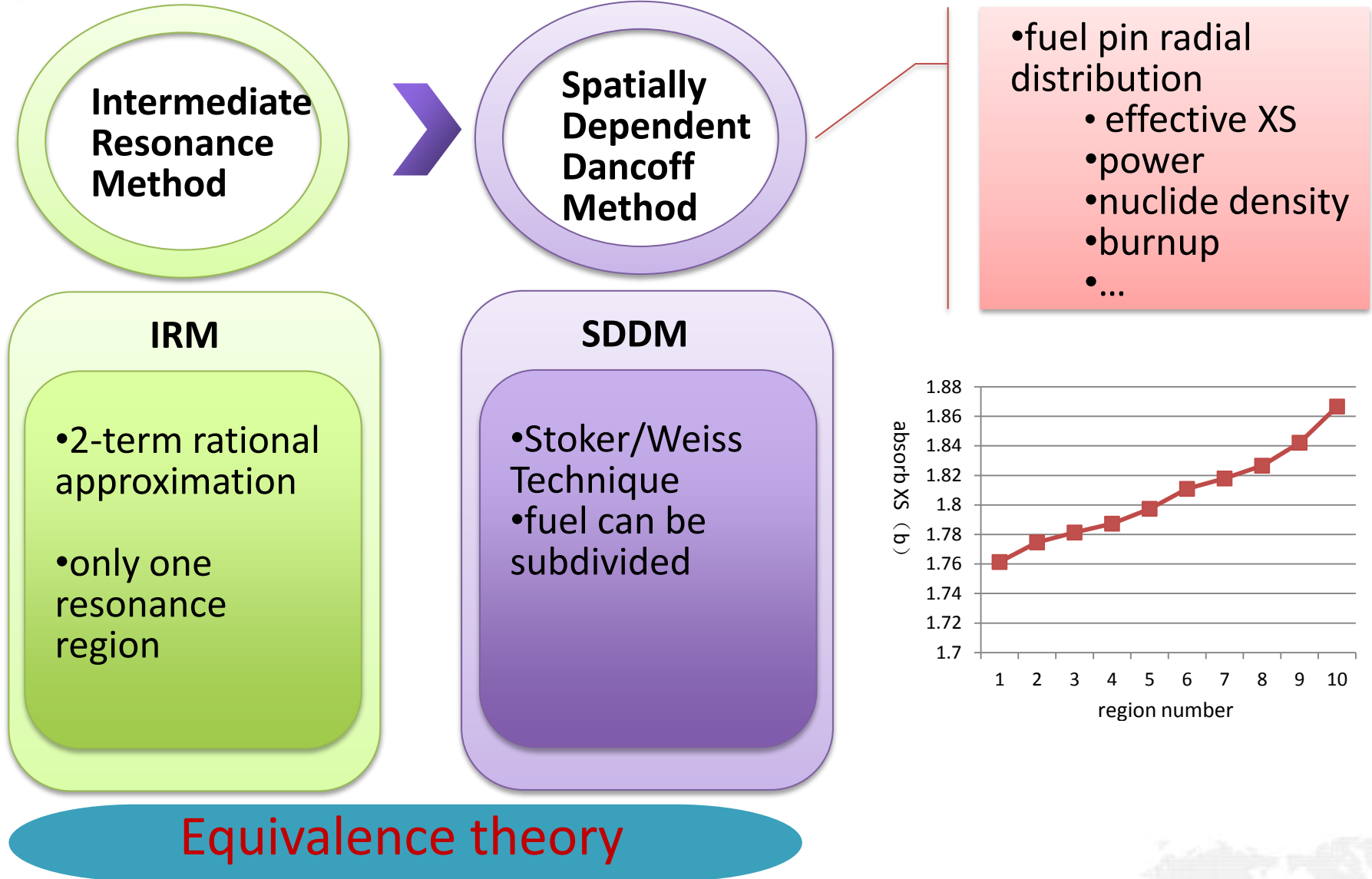
IN

- Linear heating rate
- moderator temperature
- flow rate
- material conductance

Steady State
Heat Equation

OUT

- Clad, fuel effective temperature interpolation table
- Depend on the burn up



Predictor-Corrector Method

Burn-up Equation Solving:

- The matrix exponential method
- Effective for stiff equation
- Bigger burn-up Step

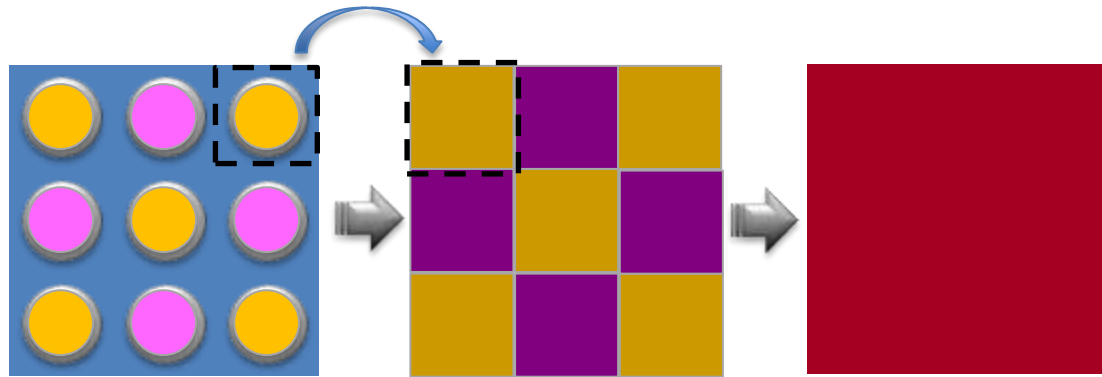


Gadolinium Depletion

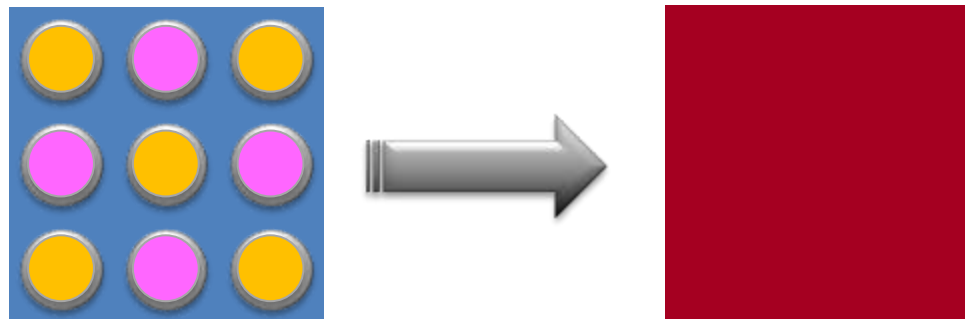
- Projected PC method
- Log Linear Rate Method



◆ Traditional two-step scheme: CP + SN



◆ One-step scheme : MOC



The transport equation can be written in the following discretised form:

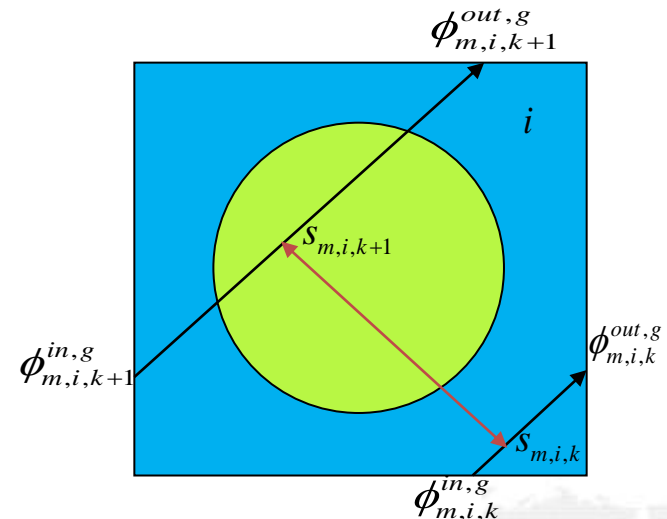
$$\frac{d\phi_{m,i}^g}{ds_m} + \Sigma_{tr,i}^g \phi_{m,i}^g = Q_{m,i}^g \quad \text{Eq (1)}$$

The outgoing angular flux along the path k in i mesh can be given by:

$$\phi_{m,i,k}^{out,g} = \phi_{m,i,k}^{in,g} \exp(-\Sigma_{tr,i}^g s_{m,i,k} / \sin \theta_m) + \frac{Q_{m,i}^g}{\Sigma_{tr,i}^g} [1 - \exp(-\Sigma_{tr,i}^g s_{m,i,k} / \sin \theta_m)] \quad \text{Eq (2)}$$

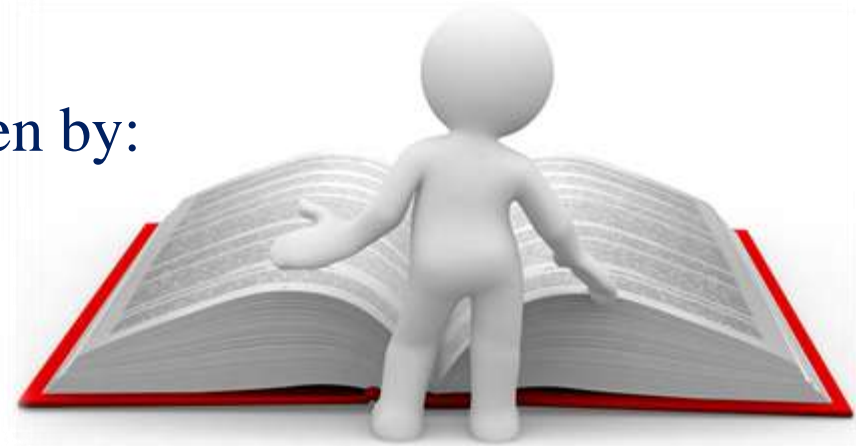
The average angular flux can be obtained by integrating the Eq (2) along the path k:

$$\bar{\phi}_{m,i,k}^g = \frac{Q_{m,i}^g}{\Sigma_{tr,i}^g} + \frac{\phi_{m,i,k}^{in,g} - \phi_{m,i,k}^{out,g}}{\Sigma_{tr,i}^g s_{m,i,k}} \quad \text{Eq (3)}$$



The average mesh angular flux is given by:

$$\bar{\phi}_{m,i}^g = \frac{\sum_k \bar{\phi}_{m,i,k}^g s_{m,i,k} \delta A_m}{\sum_k s_{m,i,k} \delta A_m} \quad \text{Eq (4)}$$



By summing up all the average angular fluxes within region i, the region average scalar flux is obtained as follows:

$$\bar{\phi}_i^g = \sum_m \bar{\phi}_{i,m}^g w_m \quad \text{Eq (5)}$$

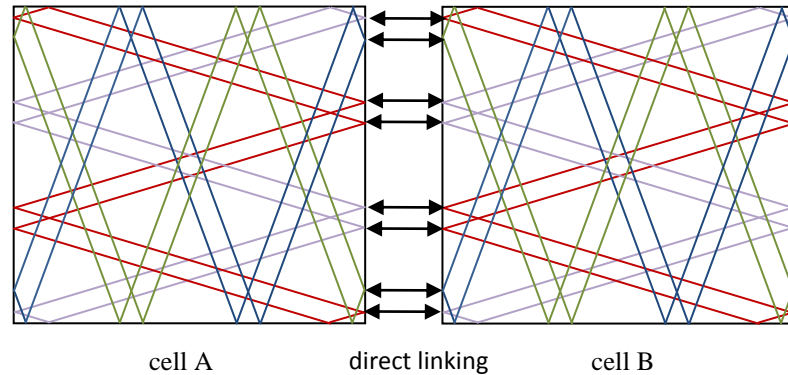


Modular Ray Tracing

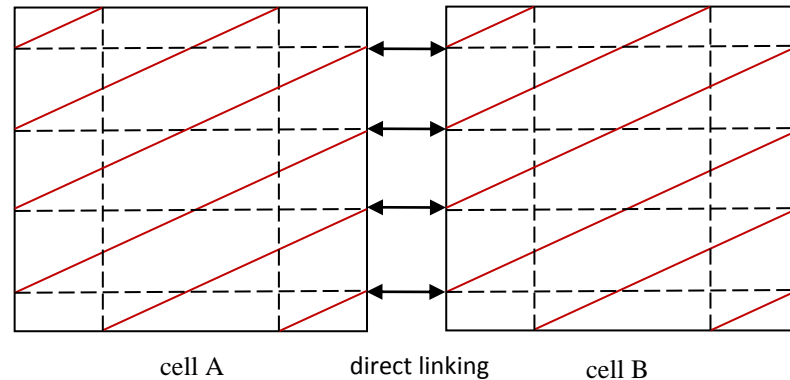
- MOC ray tracing process often takes large memory storage to save the neutron path information.
- Fortunately, in PWR core, many assembly have the same types and the number of types is not very large.
- In order to make the ray tracing be modular, each type of cell has its own neutron path information , and the path of every cell is associated with other's path

Two modular ray tracing in LATC

DNPL technique
(Direct Neutron Path Linking)



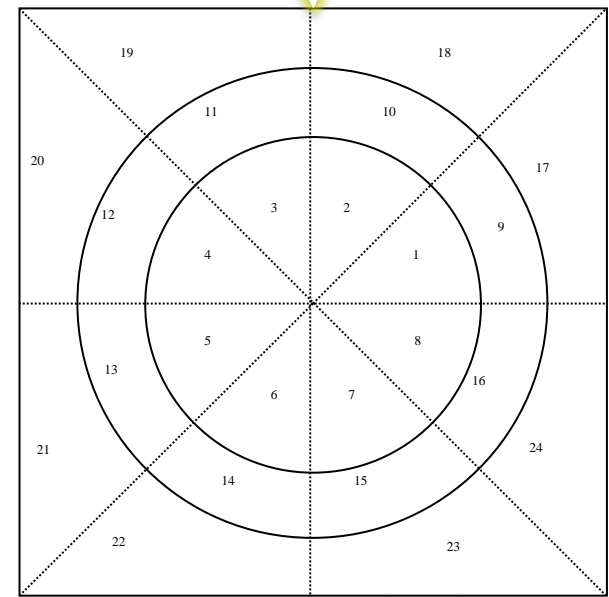
CMRT technique
(Cell Modular Ray Tracing)



Both of the two schemes can make sure that each neutron flight path of a cell is linked to a corresponding path of the neighboring cell in order to perform a continuous characteristics calculation along the path through the cell interface.

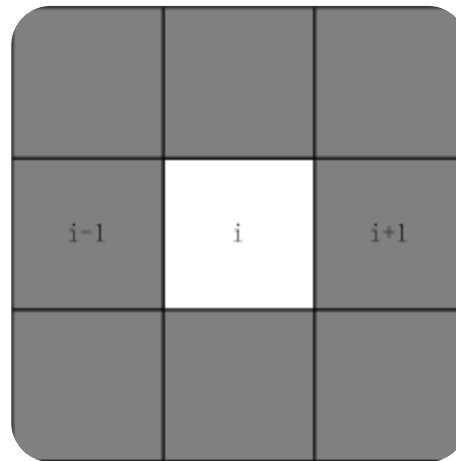
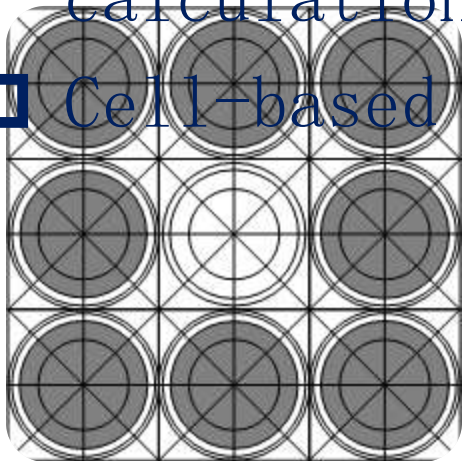
- Each cell in the problem needs to be broken into numbers of sub-regions.
- In the every sub-region, the source and flux are existed and flat.
- In this way ,it lends itself to an easy calculation of true areas for each mesh.

subdivided into octants areas.

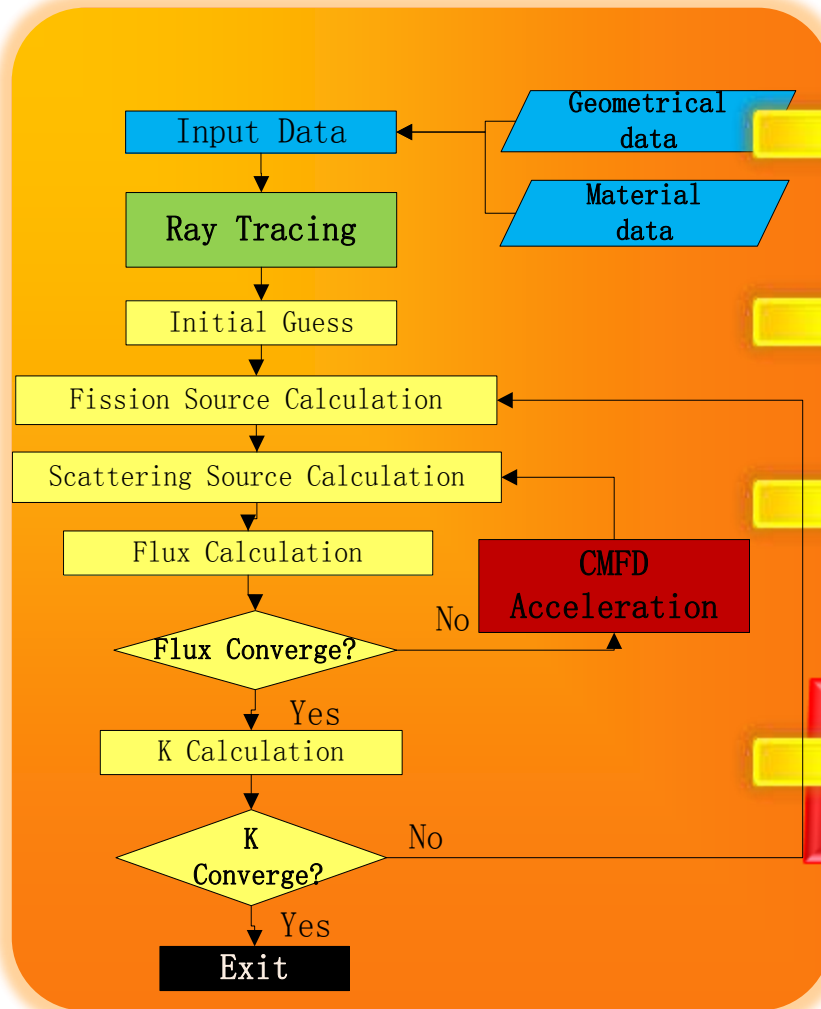


- ❑ The Coarse Mesh Finite Difference acceleration solver is implemented in LATC.
- ❑ It has been successful in reducing the computational burden for MOC transport calculation

❑ Cell-based CMFD



Transport Calculation Flow



the cross section comes from resonance calculation

generates all the necessary data related to neutron flight paths.

inner and outer iteration to calculate the flux and the eigenvalue.

the CMFD accelerate the inner iteration.



- V&V are part of the development process and with the quality assurance (QA) process.
- V&V are means by which the LATC is checked, and by which its performance is demonstrated and assured to be a correct interpretation of the requirements.



LATC Versus WIMSD code



UME-LW-AECL-aecl_um
A uranium metal critical experimental
benchmark of AECL Laboratory.

Pitch (cm)	5.359
Lattice geometry	Square
Water/fuel volume ratio	2.305
Moderator	H2O
Fuel material	U-met (0.714 wt% U-235)
Fuel Density (g/cm ³)	18.95
Radius of fuel rods (cm)	1.6255
Clad material	Al
Clad Density (g/cm ³)	2.6999
Outer radius of clad (cm)	1.7475
Thickness of clad (cm)	0.102
Temperature (all components) (K)	293.0
Experimental buckling B ₂ (cm ⁻²)	-0.00193 ± 0.00008

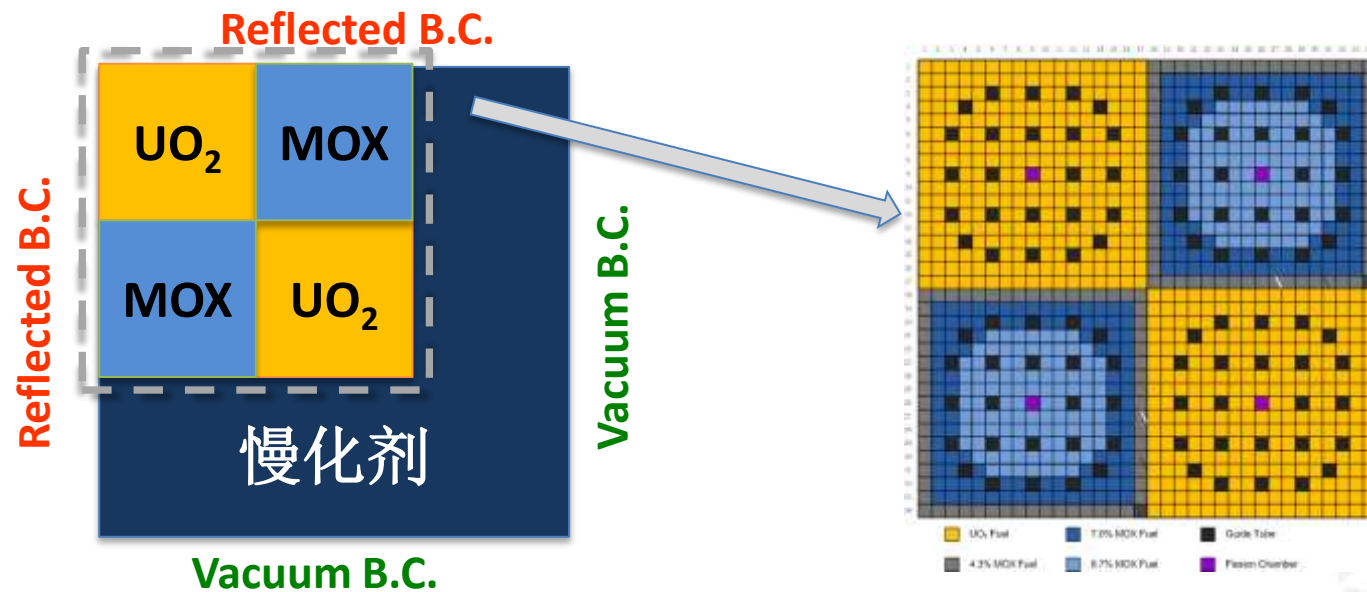
eigenvalue result

- Under the condition of
using the same iaea library.

Code	eigenvalue	Error
WIMSD-5B	1.00259	----
LATC	1.00323	64pcm

Comparison against C5G7 Benchmark

- ❑ The C5G7 benchmark is proposed by OECD/NEA for deterministic Transport calculation without spatial homogenization.
- ❑ The benchmark geometry is the sixteen assembly(quarter core symmetry) C5 MOX fuel assembly problem.
- ❑ Each fuel assembly is made up of a 17×17 lattice of square pin cells.



configurations of C5G7

The result of C5G7 benchmark

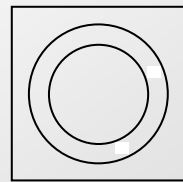
	Reference MCNP	LATC	error
Eigenvalue	1.18655	1.18744	89 pcm
Max pin power	2.498	2.502	0.16%
Min pin power	0.232	0.233	0.42%
inner UO ₂	492.8	493.4	0.12%
MOX assembly	211.7	211.5	0.09%
outer UO ₂	139.8	139.5	0.21%

CMFD acceleration effect :

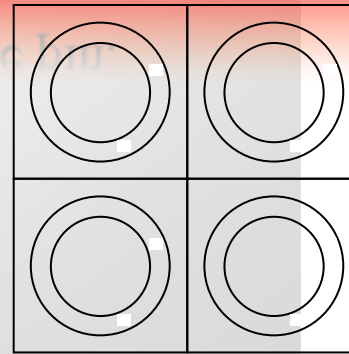
	LATC(CMFD)	LATC(no CMFD)	error	Accelerate rate
Eigenvalue	1.18744	1.18741	3 pcm	— —
Calculation time	693 s	22665 s	— —	32.7

Validation of Ray Tracing Technique

The ray tracing routine can be validated by comparing the results of different kinds of layout with the same pin.



single pin



2X2 pin by pin

- ❑ The pin by pin calculation results are fully coincided with the single pin calculation result.
- ❑ It proves that the angular fluxes at the pins interfaces are correctly transferred between pins.

Summary & outlook

01

LATC, the characteristics transport theory code, has been developing by **SNPSDC**.

02

Two ray tracing cyclical technique

- ◆ **DNPL** technique
- ◆ **CMRT** technique.

03

Effective CMFD acceleration technique is used in LATC.

04

Comparison against some codes.
Much more V&V work need to do in the next phase.



**Thank you for your
attention!**

