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Current Status of Neutronics Code System in Japan

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Contents

- Review of developing neutronics code systems in Japan
- MARBLE development project in JAEA
 - Objectives
 - Software Architecture
 - Burn-up analysis system for fast reactors
 - Database of ZPPR experiment and analysis

Review of neutronics codes in Japan

- **No standardized software platform** for neutronics code system development
- Development projects in Japan
 - Presented in the latest annual meeting of the Atomic Energy Society of Japan (AESJ)
 - **AEGIS/SCOPE2** by NFI, NEL and Nagoya Univ.
 - **NGM** by MHI
 - **MOSRA** by JAEA-Tokai (former JAERI)
 - ... more ...

The AEGIS/SCOPE2 System

- Next-generation ICFM System by
 - Nuclear Engineering, Ltd. (NEL)
 - Nuclear Fuel Industries, Ltd. (NFI)
 - In cooperation with Nagoya University

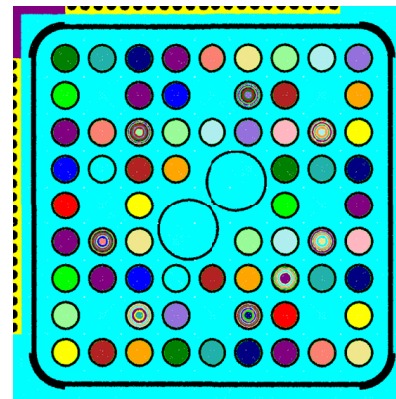
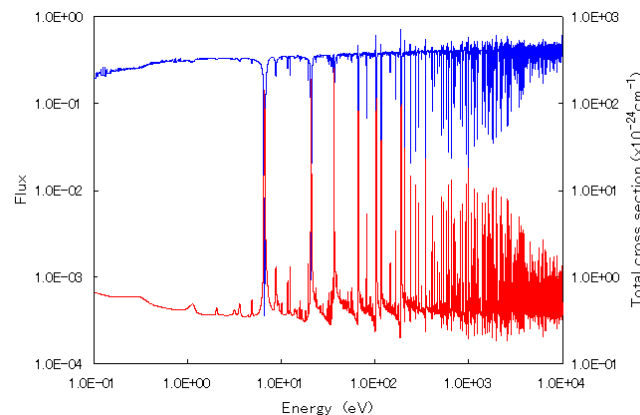
- Precise Calculation Models
 - Ultra-fine group spectrum calc. for XS
 - 3-D pin-by-pin calculation (SP3-nodal)

- Practical Computation Time
 - Acceleration and parallel calculation

M. Tatsumi, et al., "Verification of AEGIS/SCOPE2, a Next-Generation In-Core Fuel Management System," PHYSOR 2006, (2006)

Overview of AEGIS

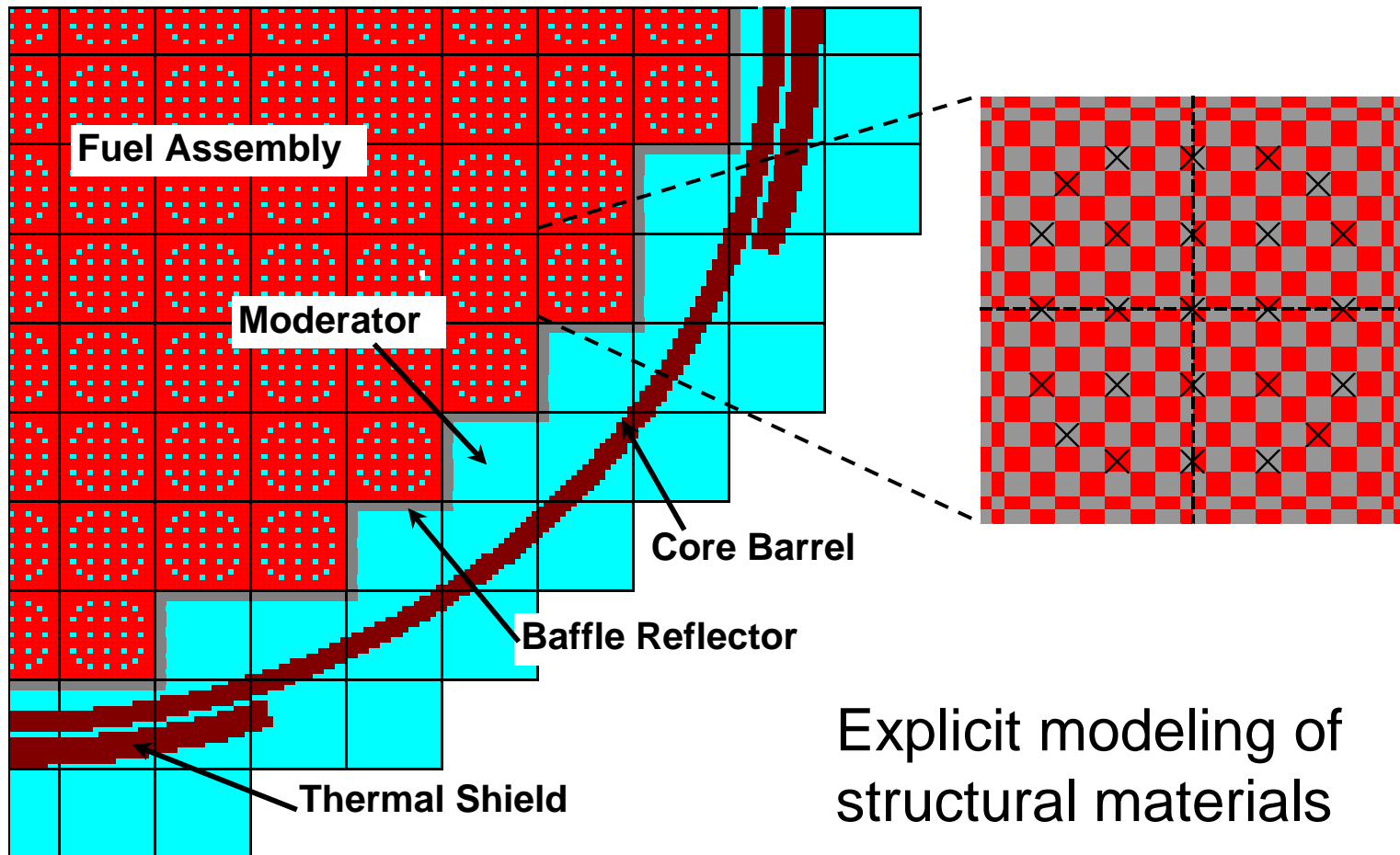
- XS generation
 - Dedicated XS library (AELIB)
 - » H.M 43, F.P. 193, Other 194
 - 32000 gr. UFG spectrum calc.
- Transport calculation
 - 172-group MOC calculation
 - Flexible geometry handling
- Burnup calculation
 - Krylov subspace method
 - Input arbitrary burnup chains



Overview of SCOPE2

- Advanced Core Calculation Code
 - 3-D fine-mesh, pin-by-pin geometry
 - SP3 nodal-transport method
 - 9-group, SPH corrected cross sections
 - Macroscopic/microscopic depletion models
 - Acceleration methods (GCMR, etc)
 - Fully parallelized using MPI ($\alpha > 99.7\%$)
 - » Identical solutions, independent processor configuration

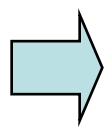
3-D, Pin-by-Pin Fine-Mesh Geometry (340x340x26 meshes for 4-loop WH type plant)



Explicit modeling of structural materials

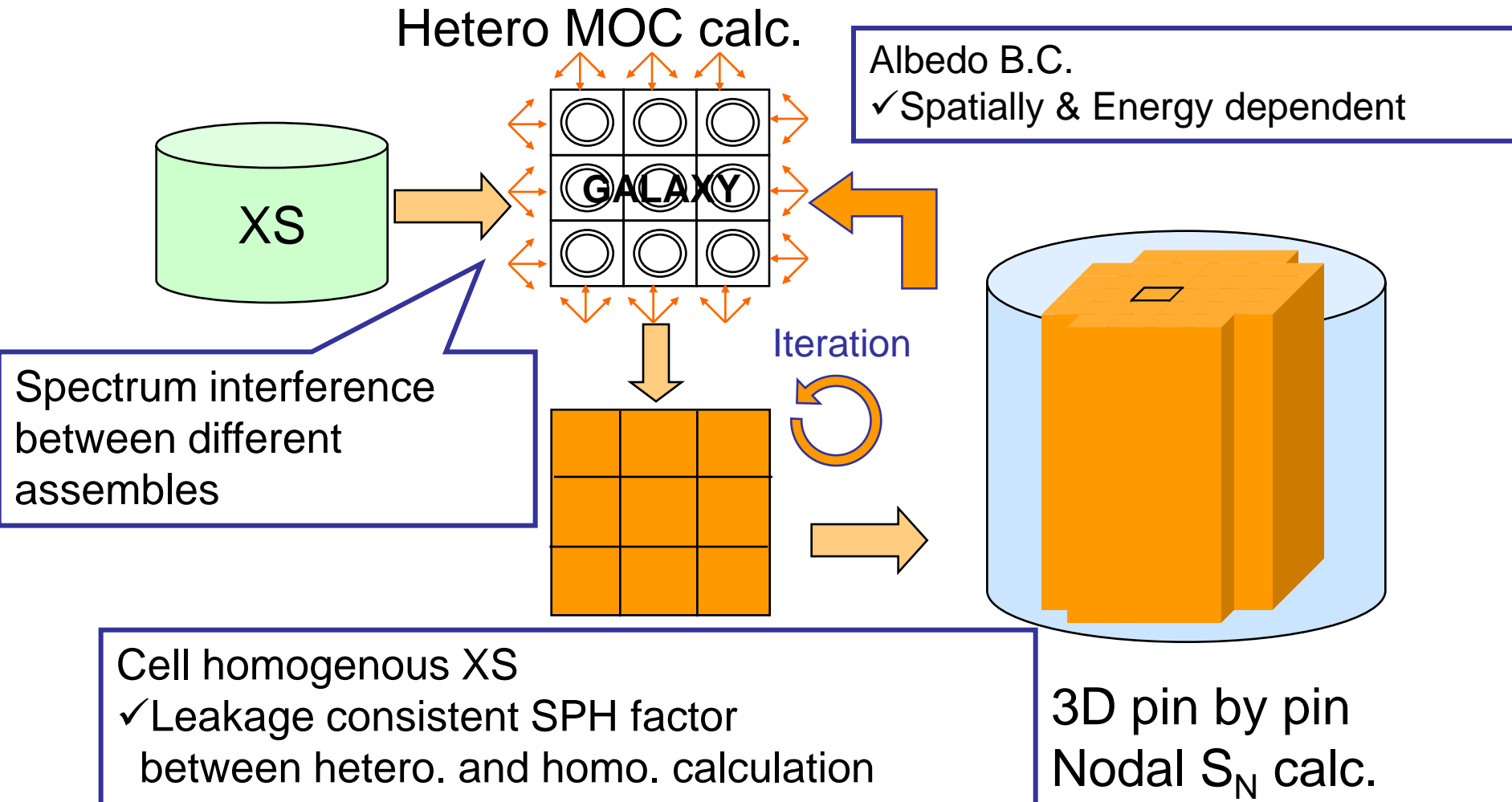
NGM: Overview

- MHI is developing the next generation method (NGM).
 - High accuracy of pin power
 - Complex geometry such as assembly gap and irregular fuel rod arrangement
- Direct 3D core calculation
 - Pin-by-pin, Multi-group, Transport
- Our NGM can treat **cell-heterogeneity**.
 - “On-flight” method



Coupling 2D method of characteristics (MOC) with 3D pin-by-pin S_N

NGM: On Flight Method



MOSRA: Background & Objectives

Development of SRAC

1978 --

:


2007 SRAC2006
Release

Completion of SRAC

- **Obsolete base design**
- **Constraint to extend (e.g. I/O compatibility)**
- **Complexity of the system**

MOSRA (**MO**dular code **S**ystem for **R**eactor **A**nalyses)

Combination, addition and deletion of modules

- 
- **Upgrading of analysis methods**
 - **Diversifying into various types of reactors**
 - **Reducing costs of maintenance and development**

MOSRA: System Components

Library processing
module (**LibMakers**)

Cell calculation module
(**M-SRAC**, J-MOC)

Core calculation module
(**M-Citation**, **Partisn+**,
M-Light)

Thermal-hydraulics module
(M-Hydro)

Fuel temperature
calculation module
(M-Fondu)

Feed-back cross section
calculation module
(M-Macron)

Reaction rate/reactivity
editing module
(**More-MOSRA**)

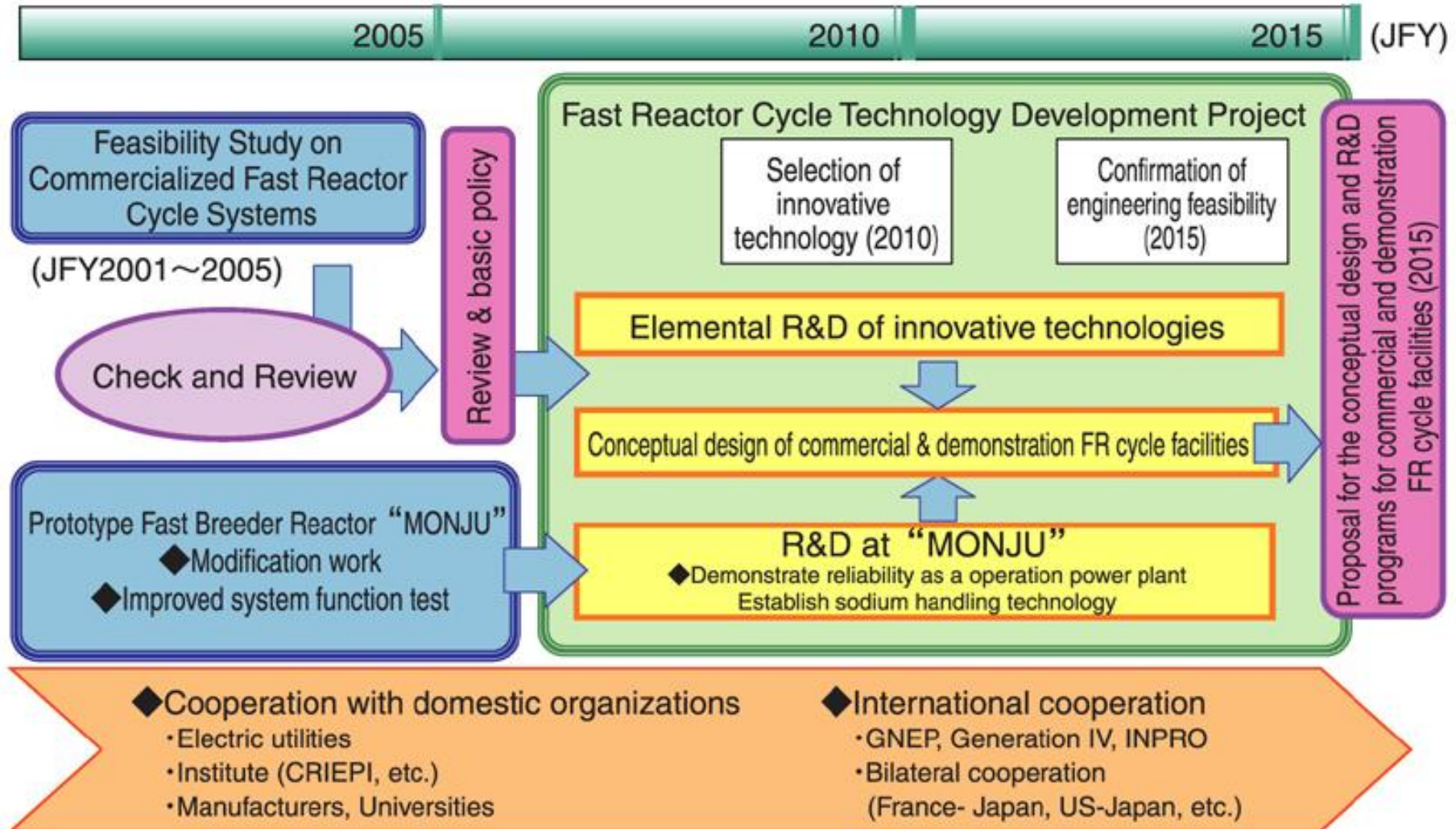
Data interface module
(PonPon)

Coupled
neutronics/thermal-
hydraulics core
burn-up calculation
frame (**M-FCBT1**)

MARBLE by JAEA-Oarai

- Background
- Objectives
- Software architecture
- Burn-up analysis system for fast reactors
- Database of ZPPR experiment and analysis

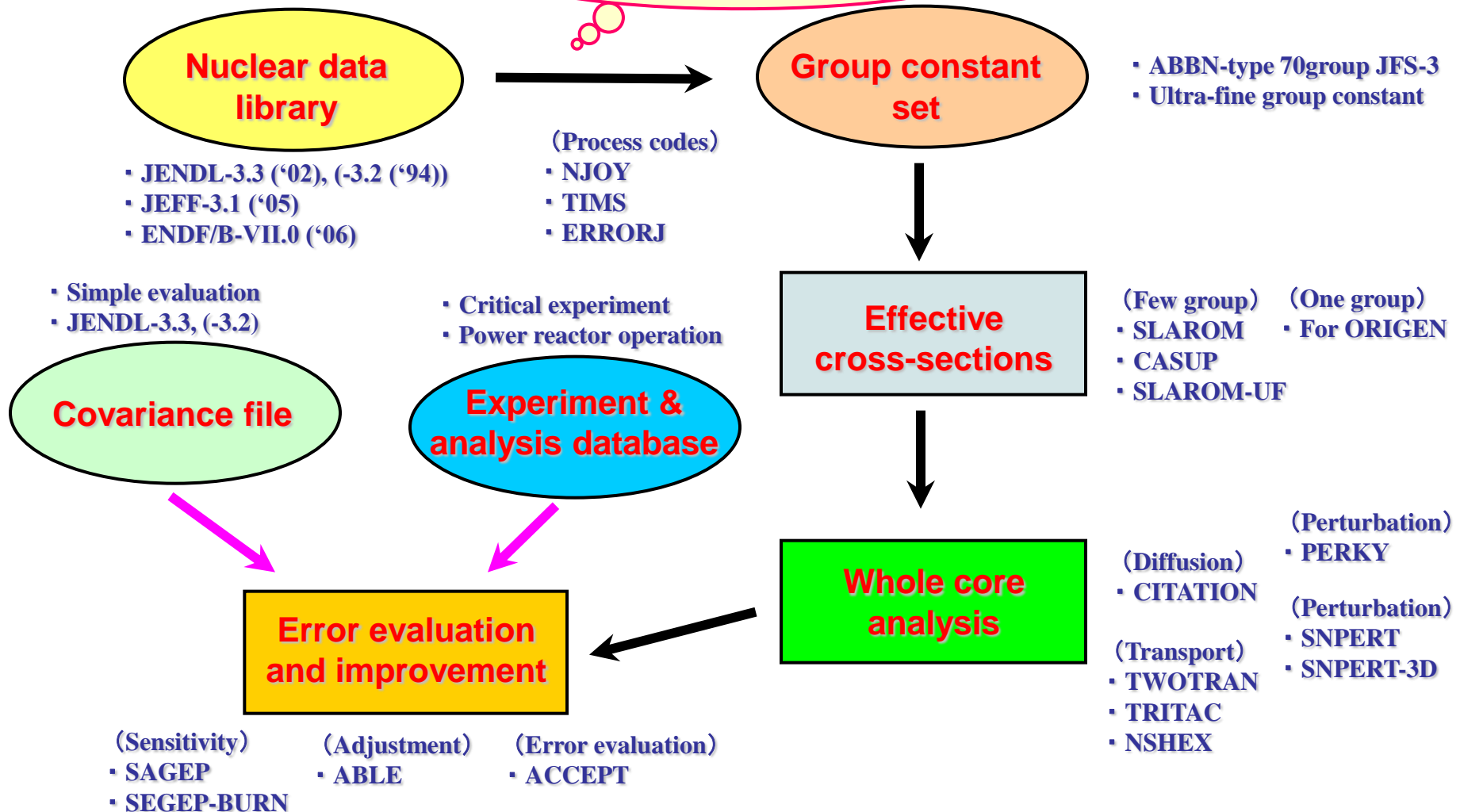
FaCT Project in JAEA



Fast Reactor Cycle Technology Development Project (FaCT)

Flow of Nuclear Analysis for FR Cores

Ultra-fine energy, 3-D, Transport, Error evaluation



Core Experimental Data



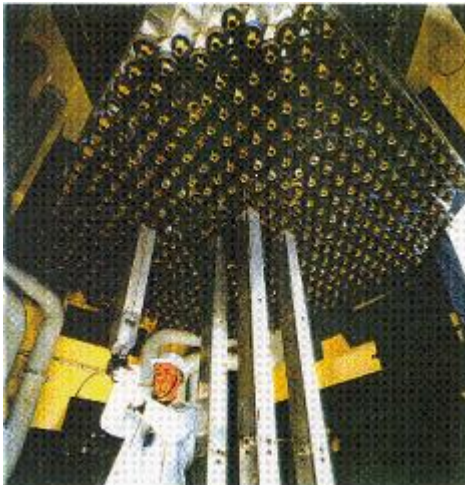
BFS-2
(Russia)



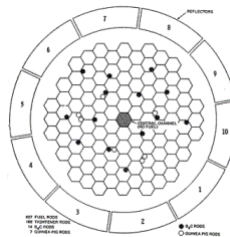
JOYO (Japan)



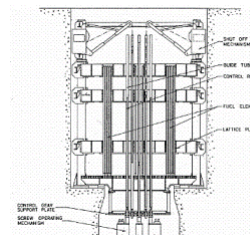
ZPPR (USA)



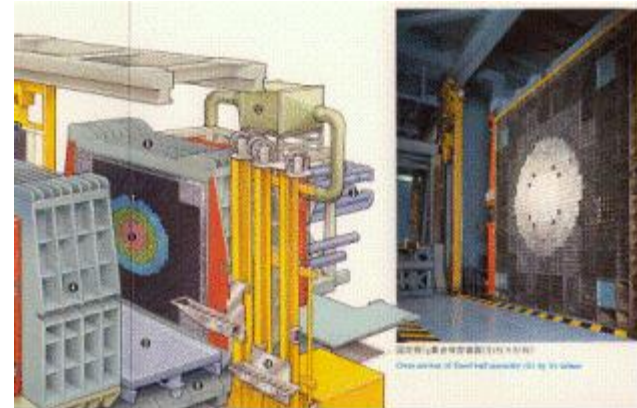
MASURCA (France)



SEFOR (USA)



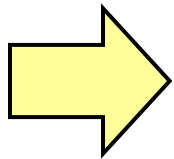
ZEBRA (U.K.)



FCA (Japan)

Problems in Conventional System

- Troublesome and/or too-many inputs
- Complicated input cards
- Difficulties in data conversion among codes



- **Inefficient processes in analysis**
- **Poor quality management**

Aim for improved efficiency and quality with rational automation mechanism.

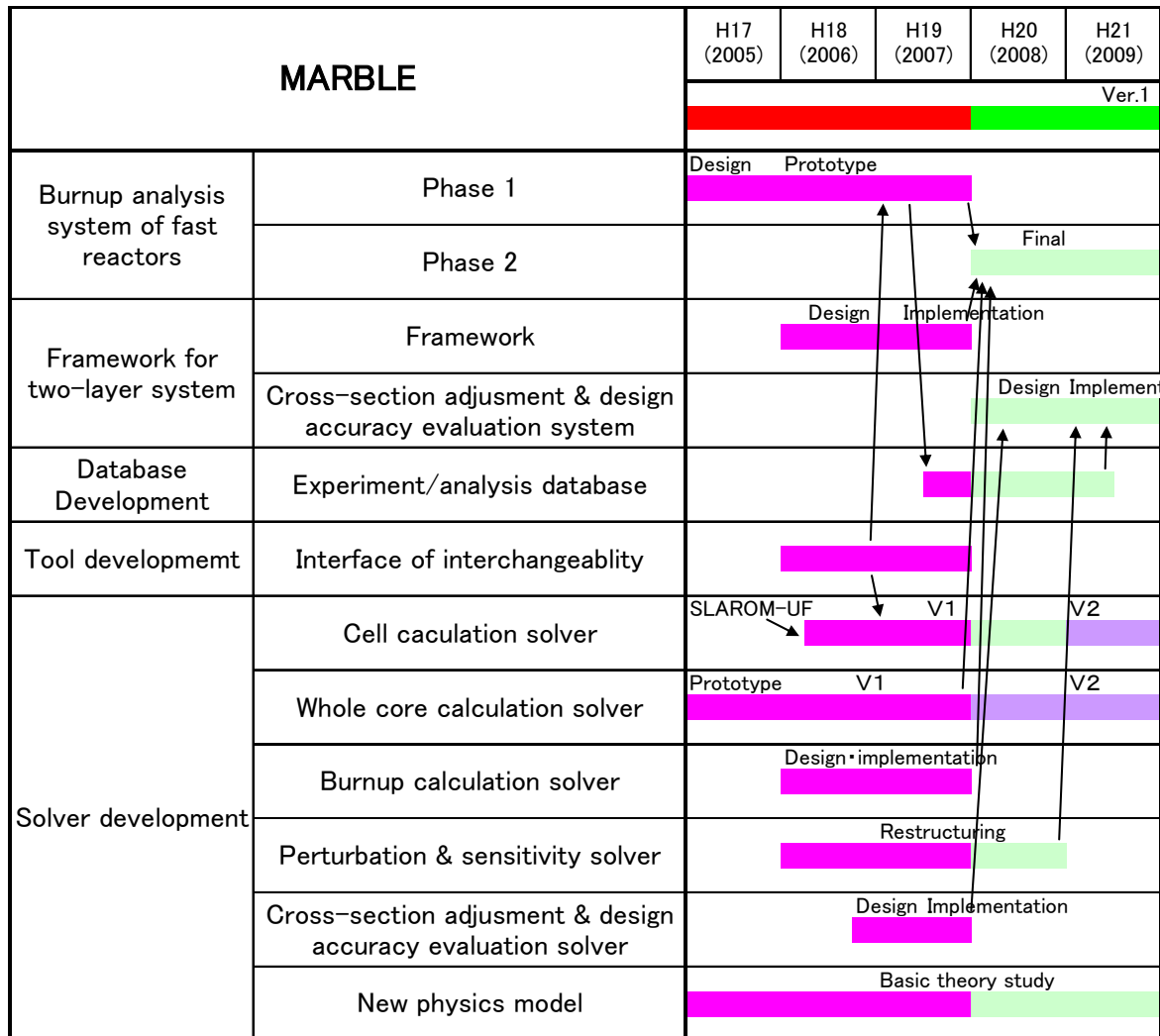
MARBLE Project (1/2)

New reactor physics analysis code system MARBLE (Phase 1: JFY2006—JFY2009)

- To improve the **quality** of **analysis** and **software**
 - **Reliability**, **reusability**, **maintainability**, and **extendibility**
- **Power reactor analysis** with fuel depletion
- **Common data model** for major calculation routes

MARBLE Project (2/2)

Developing Schedule



- **Resources**
 - ~5 man / yr
 - 90,000 euros / yr
- **Scope**
 - **Until JFY2009**
 - Replacement of the conventional code system for fast reactors
 - **After JFY2010**
 - Extension of functionality
 - Application to other types of reactors



**End of JFY2009
MARBLE Ver. 1**

Goals for MARBLE Ver.1 (1/2)

- All major existing calculation routes on new software architecture
 - Replacement of JOINT-FR and SAGEP-FR
 - Two-layer system with object-oriented languages Python and C++
 - Based on common data models
 - both for critical assemblies and power reactors
- Extension of calculation routes
 - Burn-up analysis for fast reactors

Goals for MARBLE Ver.1 (2/2)

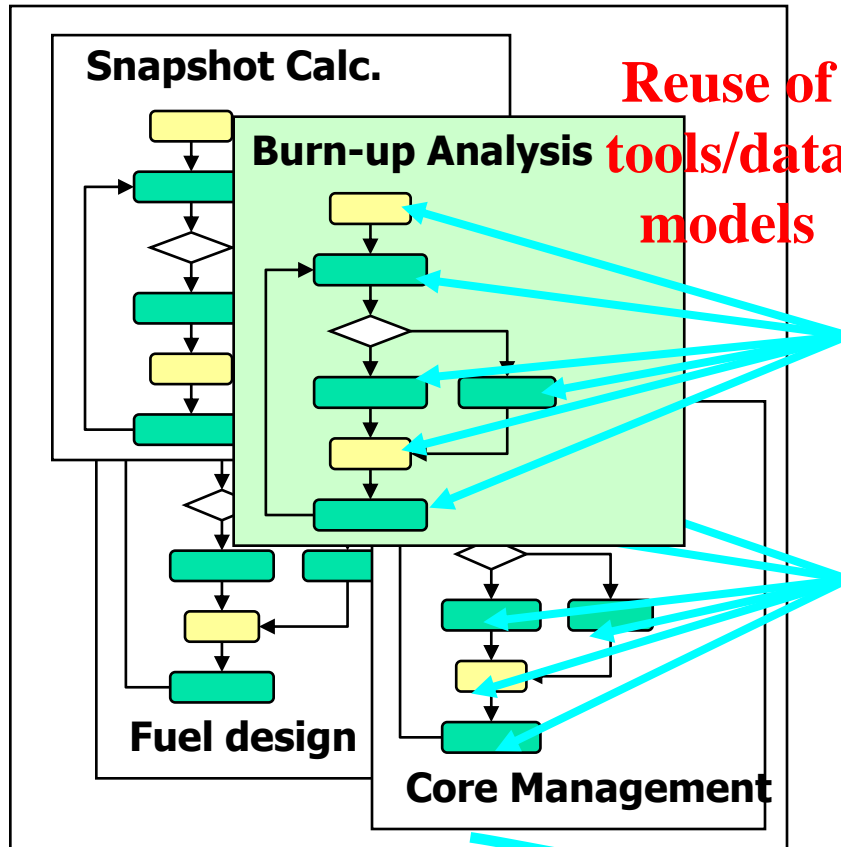
- Improvement of **operating efficiency**
 - Management of **database for fast reactor** core nuclear design (~2000 parameters)
 - **Semi-automatic analysis** with the database

Target time order for major calculation routes

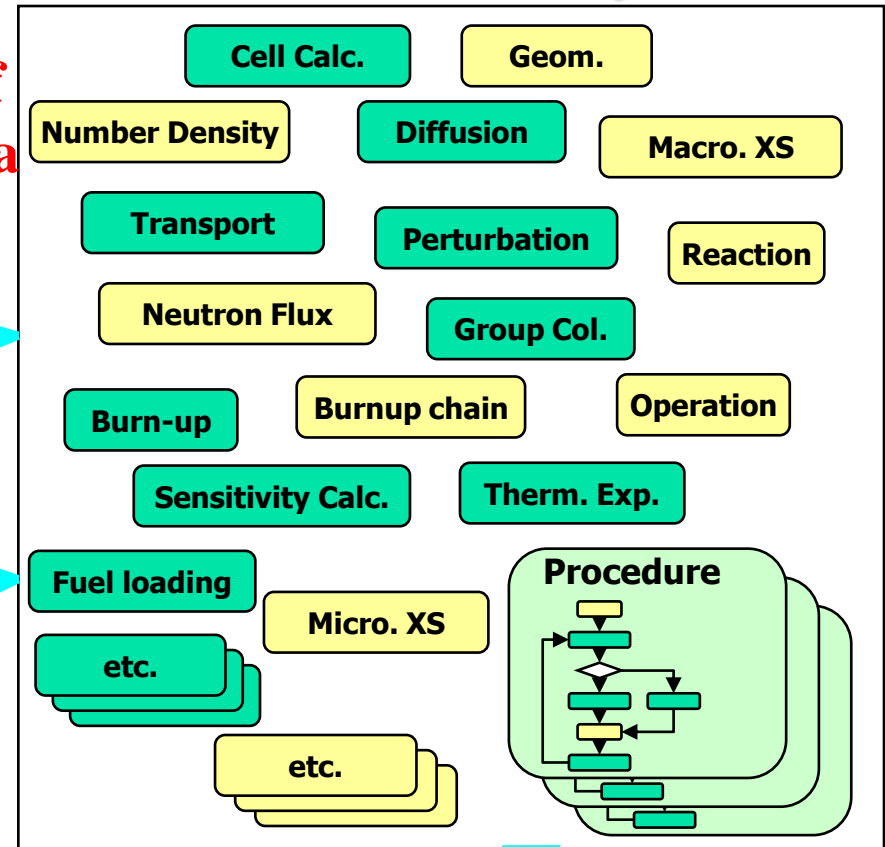
Calculation route	Model	Calculation example	Target time order
Critical assembly	Detail	Recalculation DB for XS adjustment (~200 parameters)	week
Power reactor	Detail	“JOYO” tracking calculation (~30 cycle)	day
Design analysis	Simple	A major parameter for a new fast reactor	minute

MARBLE Architecture

Control layer



Calculation layer



Reuse of
tools/data
models

- MARBLE consists of **common tools/data models**
- New sub-system (e.g. burn-up analysis) also consists of **common tools/data models**

MARBLE Modules

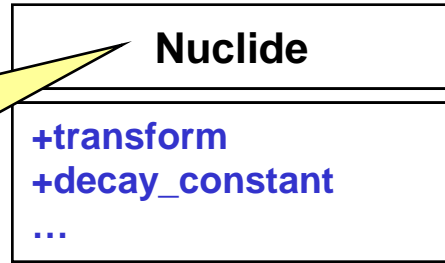
Common Data Models

- Model
 - Quantity
 - Mass, Length, Substance, ...
 - Material
 - Nuclide, Material, ReactionType, EffectiveCrossSection, ...
 - Coordinates
 - Xyz, Triz, Hexz, Zone, ...
 - Geometry
 - Point, Line, Circle, Hex, ...
 - Mesh
 - MeshConverter, ...
 - Pattern
 - Assembly, LoadingPattern, ...
 - Info
 - CoreInfo, CrossSectionInfo, ...
 - Solver
 - CitationSolver, TritacSolver, ...
- Orpheus (burn-up analysis)
 - Input
 - CalcCondition, InputInterface
- Code
 - Adjust
 - XSAdjustment, BiasMethod, Parameter, Covariance, ...
 - Burn
 - Burnup, BurnupChain, ...
 - CBG
 - PLOS (diffusion)
 - SNT (transport)
- Utils
 - ConventionalCodeUtils
 - SLAROM-UF
 - JOINT-FR
 - SAGEP-FR
 - Python
 - FortranUtils

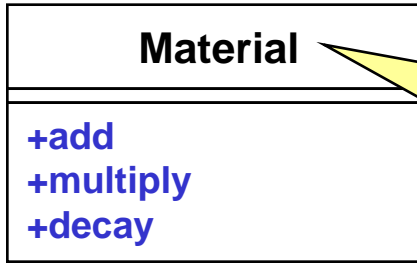
Designed to be **re-usable** and identical to **neutronic concepts**

Re-use of Data Models

Nuclide such as Pu-239, U-238, Na-23



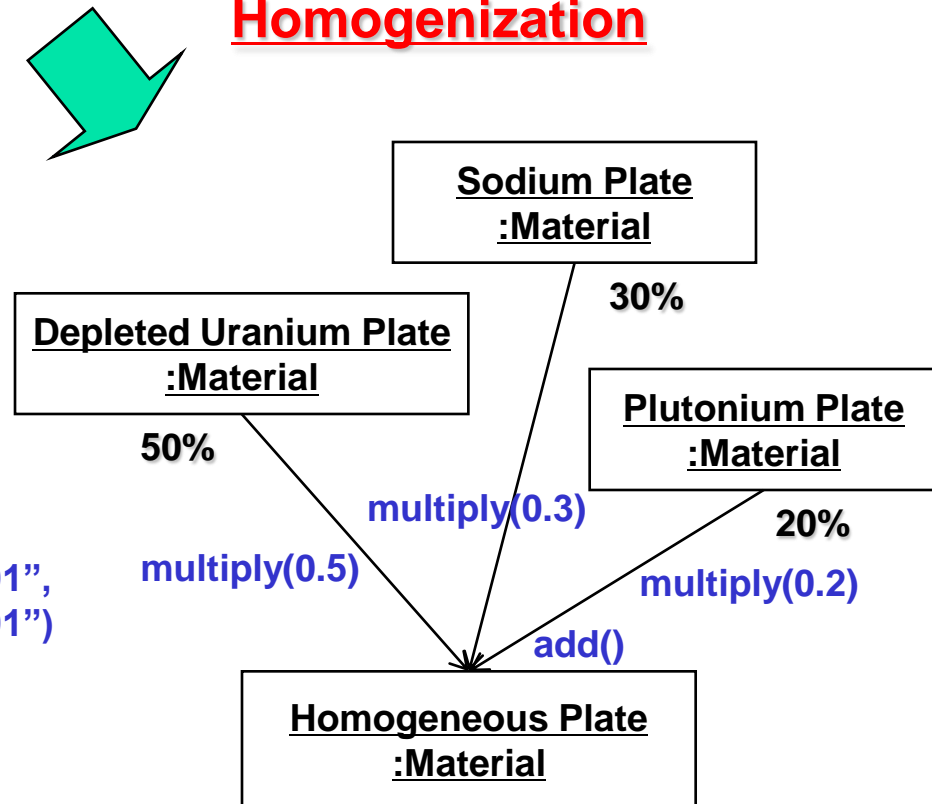
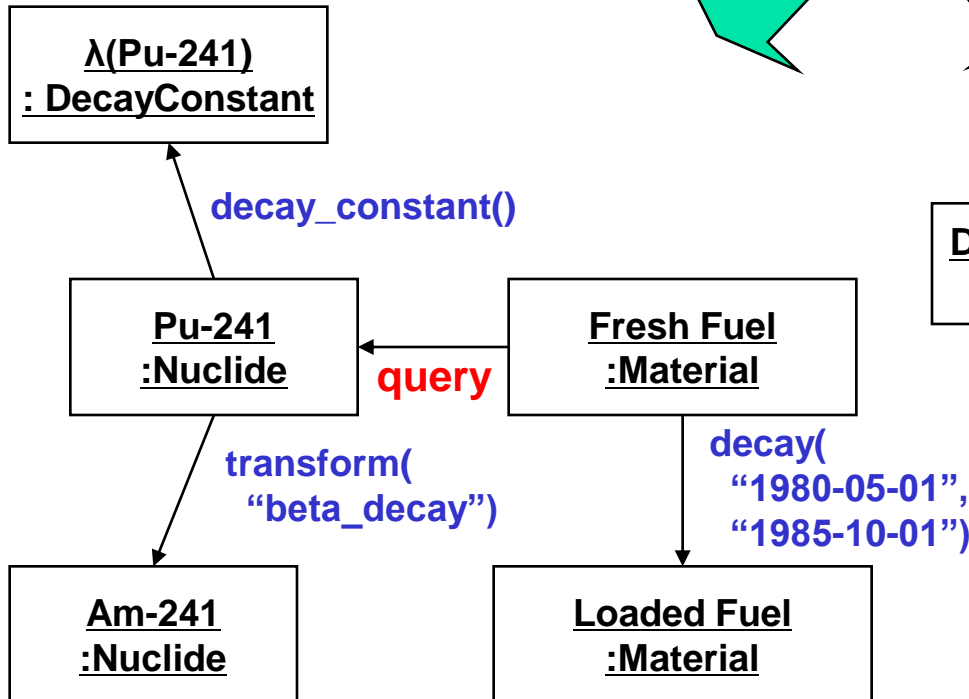
Material



Homogeneous material defined by nuclide densities

Decay correction

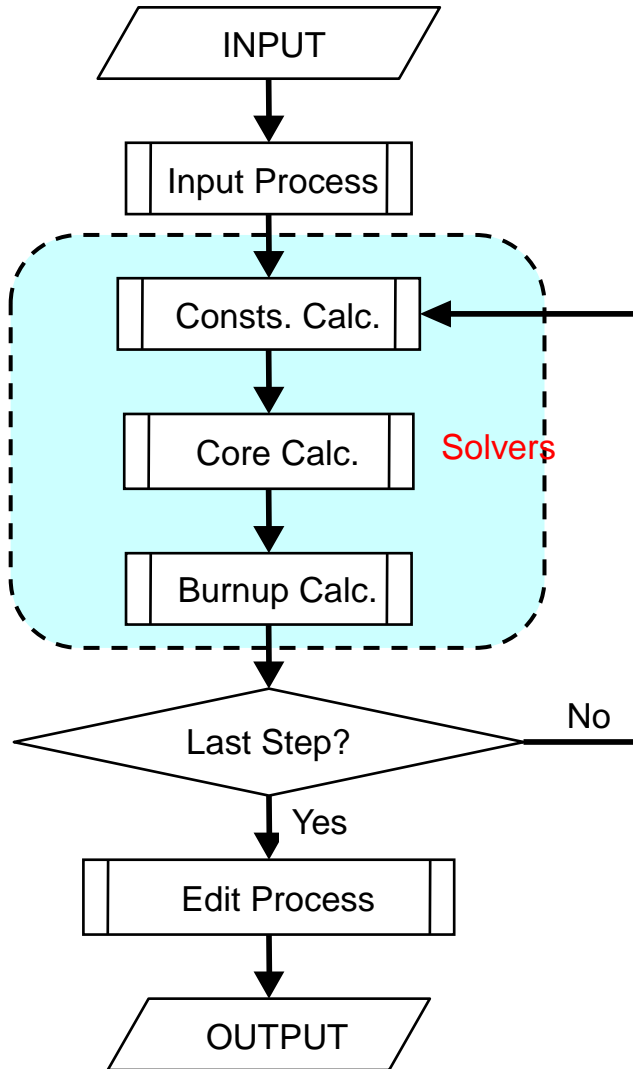
Homogenization



Requirements for Burn-up Analysis

- Adopt **arbitrary calculation codes** : lattice code, diffusion/transport codes...
- Treat “**meta inputs**” such as geometry/material composition/fuel loading pattern...
- Realize **automatic input/output conversion** among various codes
- **Switch codes** for analysis routes with ease

Burn-up Analysis System Overview



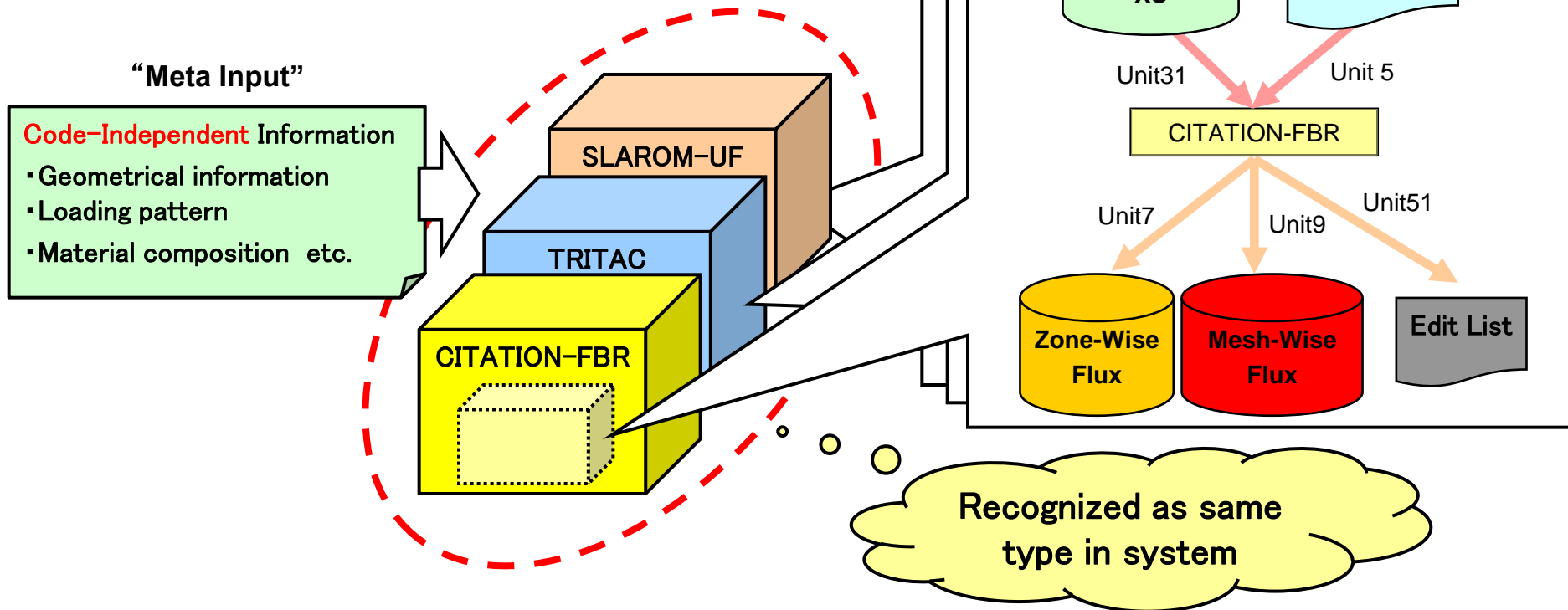
- **Input Process**
 - Input read, data preparation
- **Constants Calculation**
 - Calc. of microscopic cross sections
- **Core Calculation**
 - Calc. of neutron flux/reaction rates
- **Burnup Calculation**
 - Number densities
- **Edit Process**

**Design and Implementation of
Solver Management Mechanism**

Solver Management (1/3)

Concept:

Unified treatment for all codes



Control calculation codes with "meta input"

Solver Management (2/3)

Material Composition

```
material:
- name: fuel1
  composition:
  - ["U-234", 1.0e-20]
  - ["U-235", 0.00196132]
  - ["U-236", 7.76703e-10]
  ...
```

Fuel Loading Pattern

```
fuels:
- label: PFD066
  type: fuel_pfd1

load_fuels:
- { address: 1A1, label: PFD066 }
- { address: 1C1, label: PFD067 }
- { address: 1E1, label: PFD068 }
- { address: 2B1, label: PFB010 }
- { address: 2E2, label: PFC010 }
...
```

Calc. Condition

```
calc_system:
  coordinates:
  core : xyz
  burnup: xyz

solver:
  core:
  name: tritac
  burnup:
  name: burnup

step:
- period: 13.905
  power : 97.394

- period: 13.905
  power : 97.394
...
```



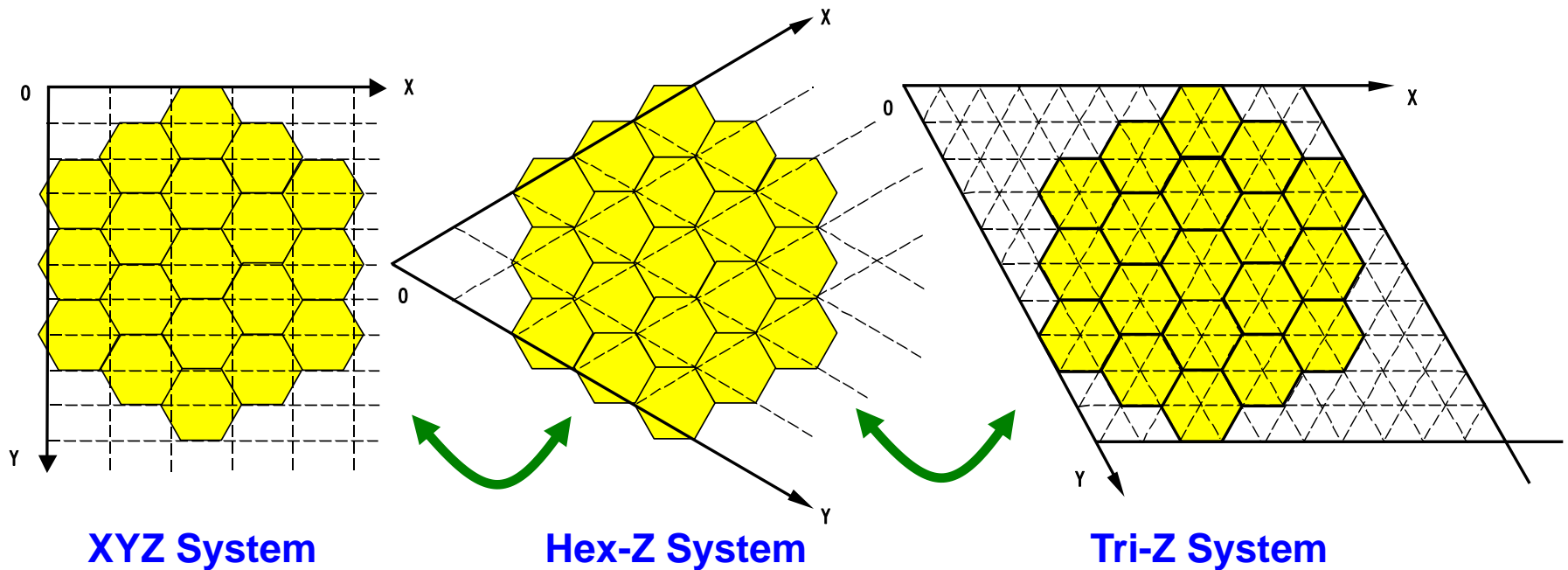
TRITAC: XYZ System

```
JOYO MK-2 Burn-UP 100.00MW x 27.81day 2781.00MWD 0cy 13.9050day x 2
0
1 1 0 0 0 0 1 1 0
21 42 96 70 555 8
21 42 18
0 0 0 0 0 0
1.00000E-04 1.00000E-04 1.00000E-05 1 10 99999.00000 0.00000
10 1.00000E-02 1.00000E+00 900
1 1 1 1 1 1 1 1 1 1 1 1
1 1 1 1 1 1 1 1 1 E
1 1 1 1 1 1 1 1 1 1 1 1
1 1 1 1 1 1 1 1 1 1 1 1
1 1 1 1 1 1 1 1 1 1 1 1
1 1 1 1 1 1 E
5 17 10 6 6 1 3 2 1 4 2 10
1 3 3 12 2 8 E
0.0000 7.0855 14.1710 21.2565 28.3420 35.4275
42.5130 49.5985 56.6840 63.7695 70.8550 77.9405
85.0260 92.1115 99.1970 106.2825 113.3680 120.4535
127.5390 134.6245 141.7100 148.7955 E
0.0000 4.0908 8.1816 12.2724 16.3633 20.4541
24.5449 28.6357 32.7265 36.8173 40.9082 44.9990
49.0898 53.1806 57.2714 61.3622 65.4531 69.5439
...
```

Minimum change to switch the calculation code!

Solver Management (3/3)

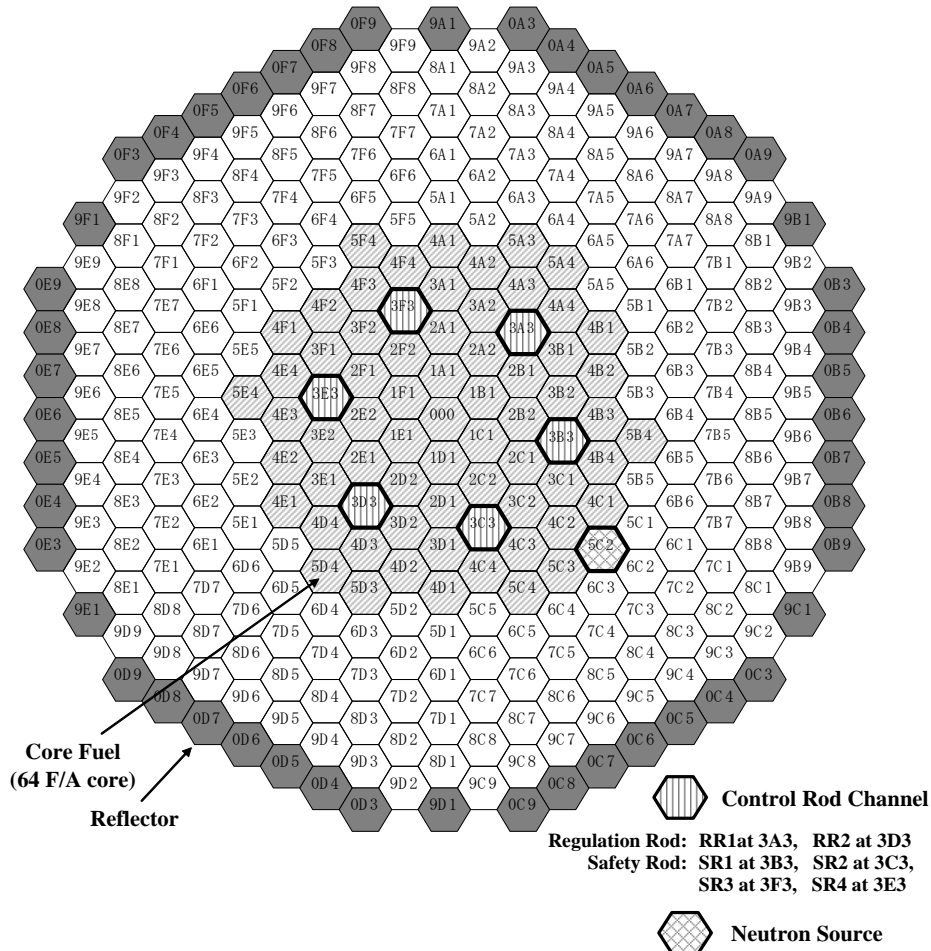
- Automatic construction of computational mesh system with lattice geometry information and fuel loading pattern
- Interoperability between different mesh system
 - Ex. Tri-Z for core calculation and Hex-Z for burnup calculation



No need to worry about mesh conversion!

Verification (1/2)

■ Burn-up reactivity coefficient of JOYO MK-I



Item	Specification
Core fuel	U-Pu mixed oxide
Pu weight fraction (w/o)	17.7
Fissile Pu ratio (w/o)	80.4
Blanket fuel	Depleted-U oxide
Coolant	Sodium
Absorber	Boron carbide
Reactor thermal power (MW)	50
Core fuel height (mm)	600
Equivalent core diameter (mm)	~720
Axial blanket thickness (mm)	400
Radial blanket thickness (mm)	~300

Verification (2/2)

Calculation results of the burn-up reactivity coefficient measured in JOYO MK-I 50MW power-up test

		JENDL	
		3.2	3.3
Base calculation [a]		-0.797	-0.801
Correction factor [c]	Heterogeneous cell	0.998	
	Ultra-fine group	0.999	
	Transport & mesh	1.012	
	CR position [b]	1.036	
Corrected calculation		-0.835	-0.831
Experiment [a]		-0.80	
C/E		1.038	1.041

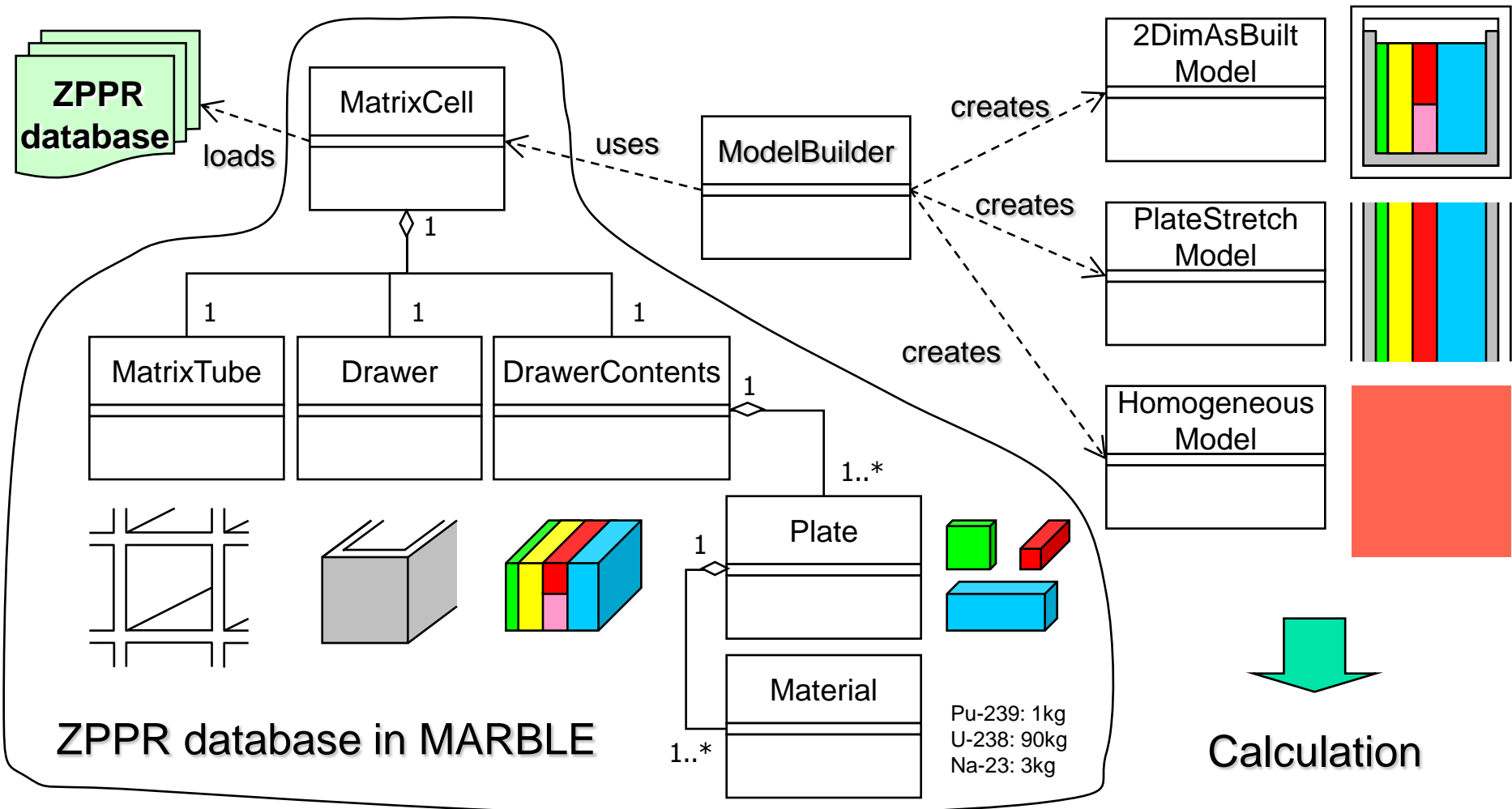
[a] burn-up reactivity coefficient (unit: 10^{-4} dk/kk'/EFPD)

[b] control rod interaction effects evaluated by JOINT-FR (conventional code system)

[c] Calculated with JENDL-3.2

Database Management

- Keeping all as-built data and supplying the data to users
- Systematic creation of calculation models



Development Methodology

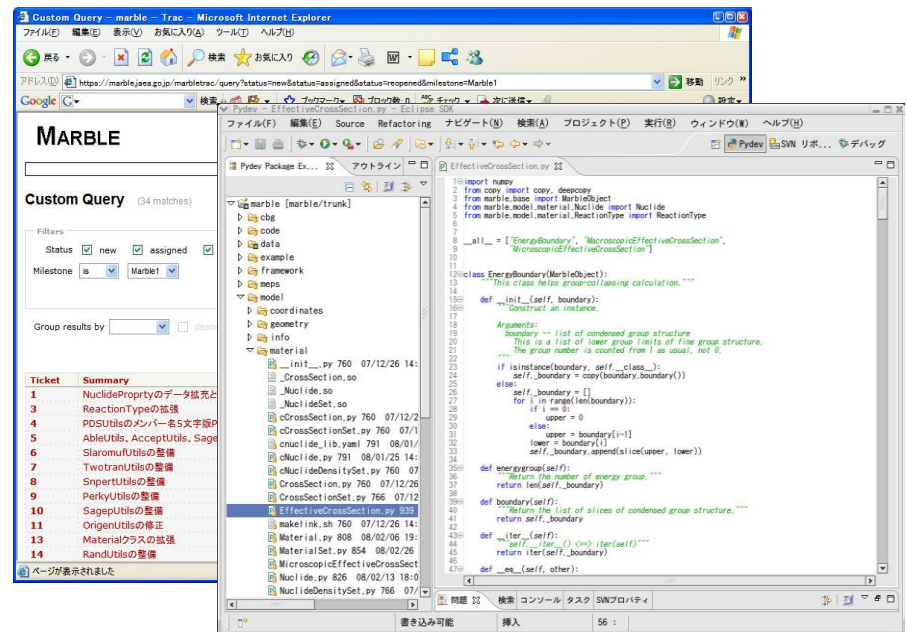
- Object-Oriented Analysis and Design

- Modeling in UML (Unified Modeling Language)
- Agile Development Scheme
- Iterative and incremental development
- Test-first programming
- Active refactoring, ...etc.

- Rapid Implementation with Python Scripting Language

- Development Tools

- Subversion(version control), Trac (bug tracking), Eclipse (IDE) etc.



Concluding Remarks

- Review of recent developing projects in Japan
 - A lot of efforts for next generation code systems
 - AEGIS/SCOPE2 by NFI, NEL and Nagoya Univ.
 - NGM by MHI
 - MOSRA by JAEA
- Development of MARBLE by JAEA
 - MARBLE Ver.1 for fast reactors (end of JFY2009)
 - Replacement of the conventional code system JOINT-FR and SAGEP-FR
 - Better burn-up analysis system for nuclear design
 - Improving database management for experiment and analysis