

Neutronics computations in support of the CABRI core safety analysis



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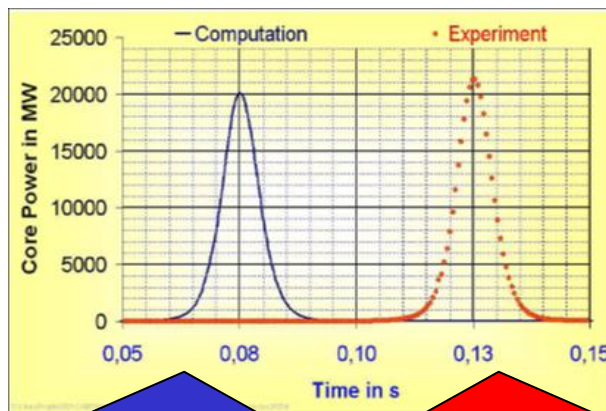
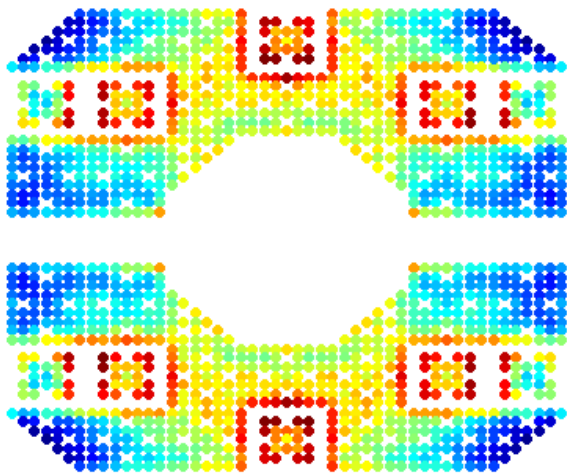


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- ➔ Description
- ➔ Validation
- ➔ Application to safety assessment

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Economic and Safety issues

CABRI is an experimental reactor, operated by the CEA and devoted to safety programs.



■ Historically: Improve the fuel behavior knowledge (RIA&LOCA)

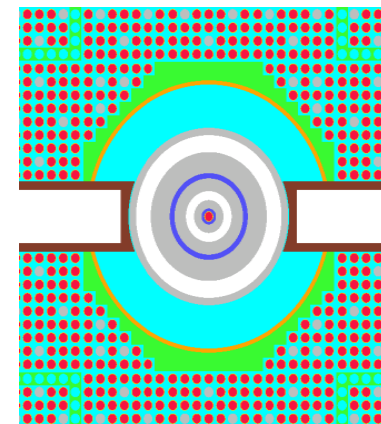
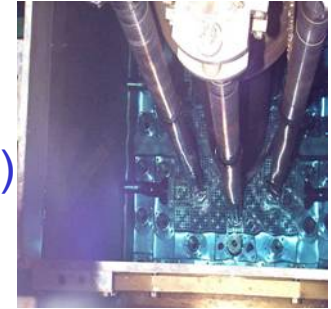
➤ Most of the experiments were dedicated to fast reactors

■ 2003: Installation shut down for renovation

➤ Sodium cooled experimental loop → pressurized water loop:
The main goal is to meet thermal hydraulics parameters identical to LWR.

■ 2008: Safety report for the new configuration

■ 2010: Commissioning tests & start-up

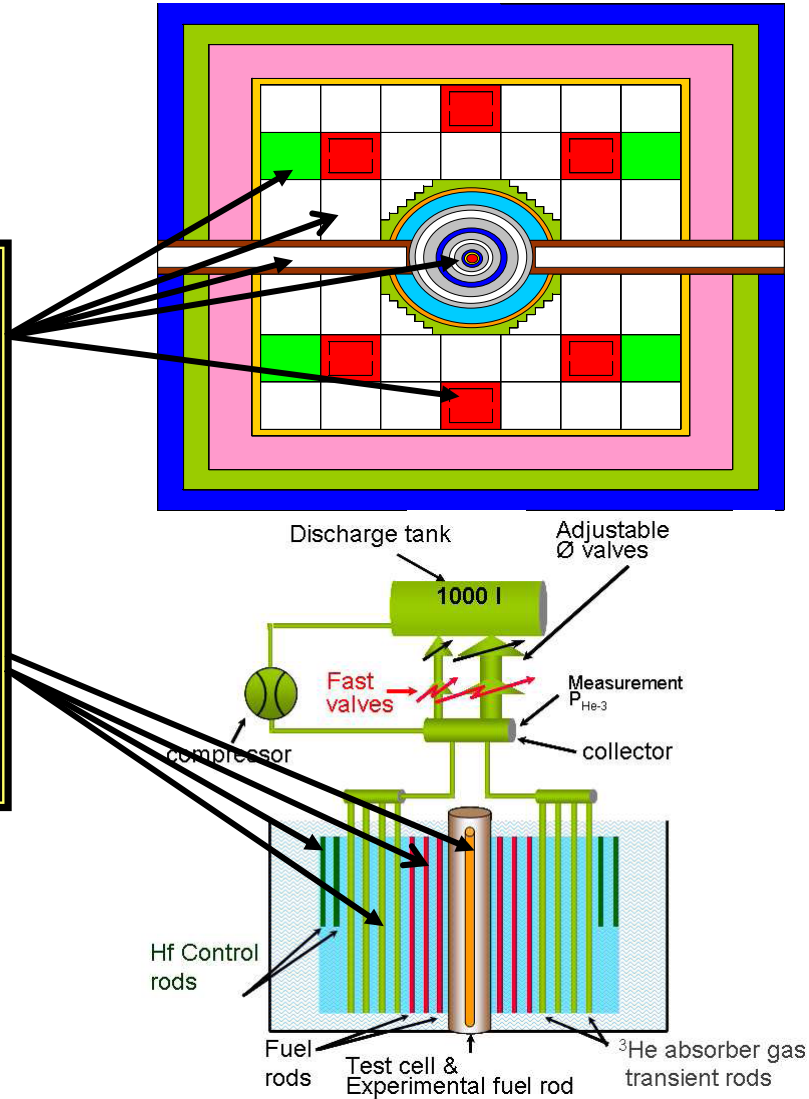


Description of the CABRI core

CABRI is a pool-type reactor.



A vertical core made of 43 rods
 The 30 cm diameter pressurized
 water loop, which circulates at a
 temperature of 250°C, is used to
 cool the core. The core is made
 of rods of 80 cm with a 0.7%
 axial experimental device, the
 experimental device, which can
 be rapidly inserted into the
 discharge tank.



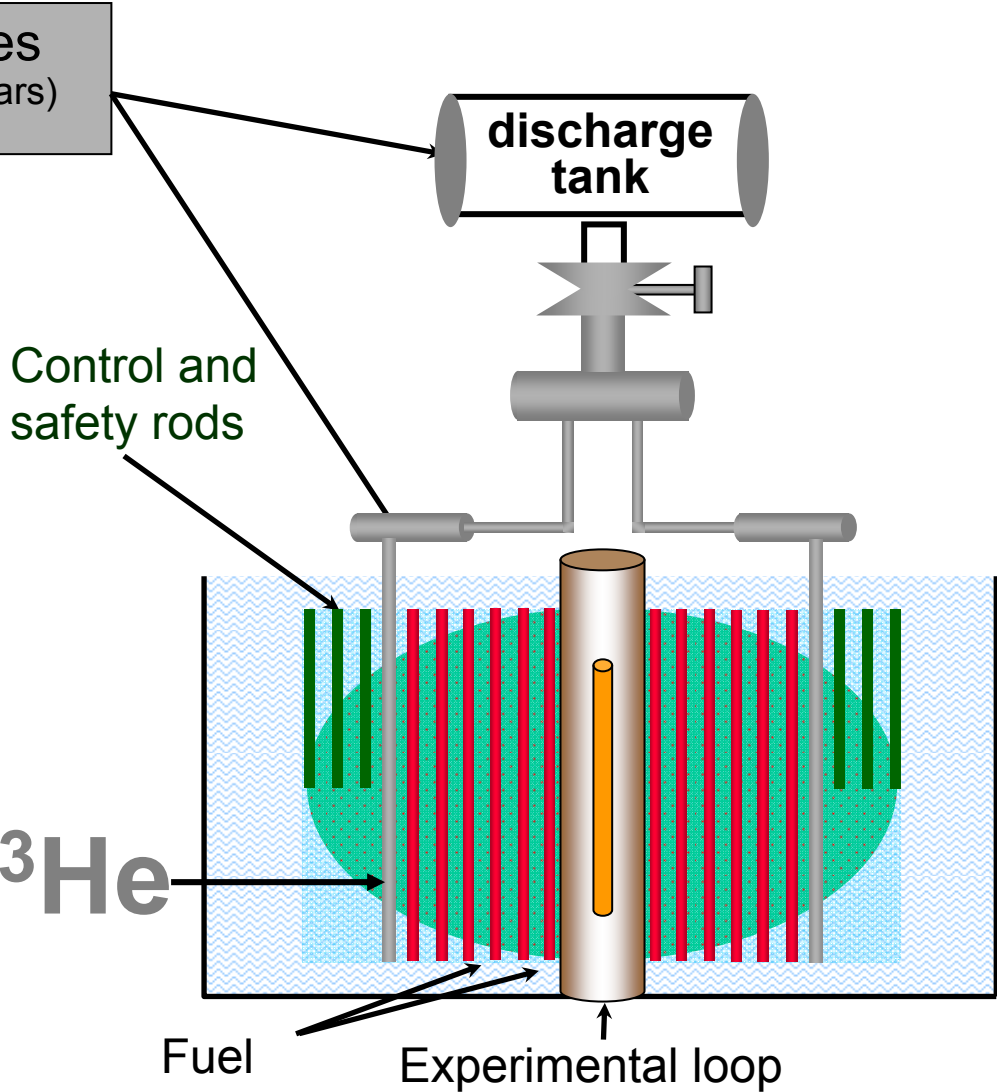
Phenomenology



Pressurized Helium tubes
(maximum allowable pressure is 15 bars)

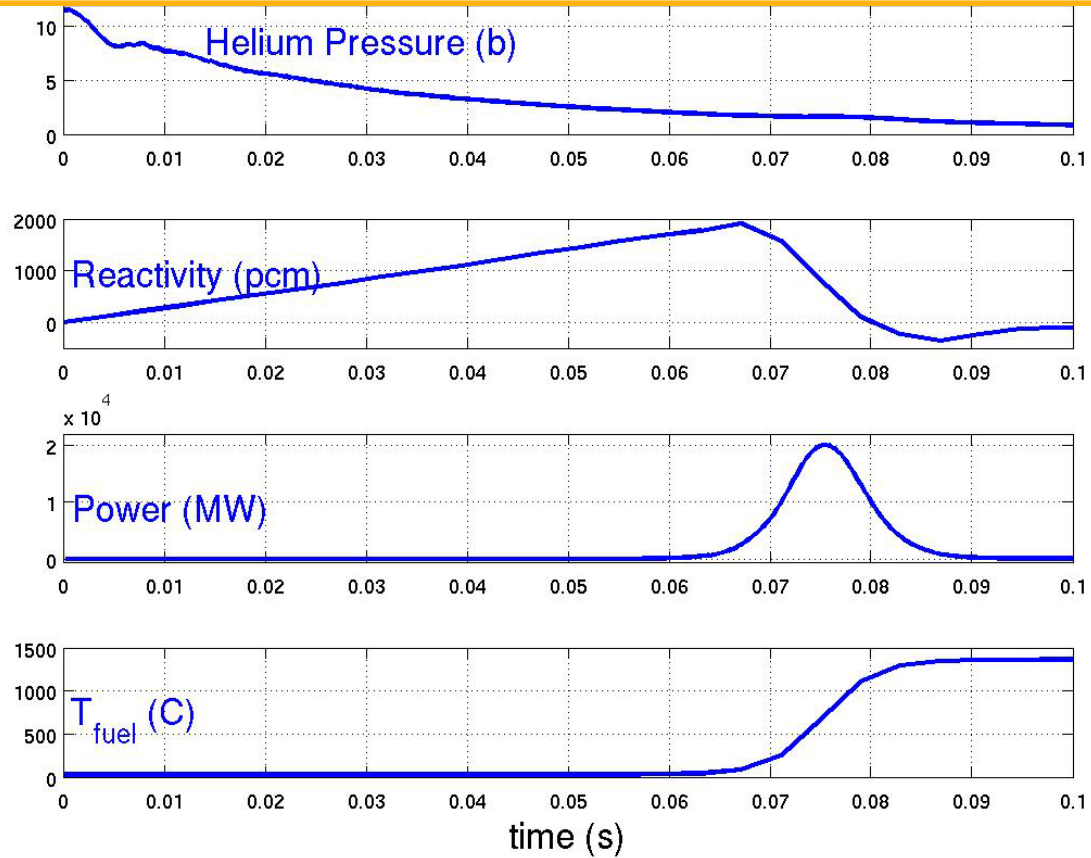
Control rod (Hf) withdrawal
The critical state is obtained by control rod withdrawal to compensate the ^3He antireactivity.
For different helium pressures, TRIPOLI4 predicts the critical position with a good accuracy

Fast ejection of ^3He
after opening of the circuit valves



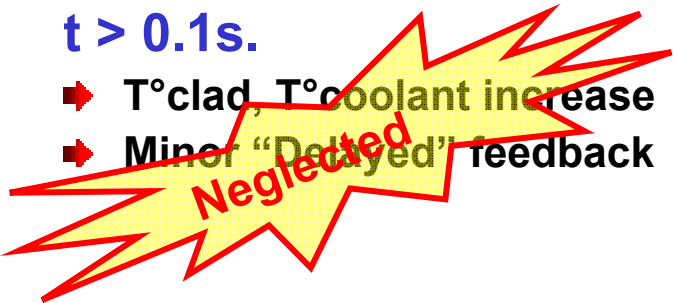
Phenomenology of the experiments

- t < 0.1 s. (adiabatic)**
- ➔ Valves opening
 - ➔ ³He depressurization
 - ➔ ↑ reactivity (3-4 \$)
 - ➔ ↑ Power (0.1 à 20000MW)
 - ➔ ↑ T_{fuel} (> 1500 °C)
 - ➔ Doppler Feedback
 - ➔ New equilibrium state (reactivity = 0 , new fuel T°)



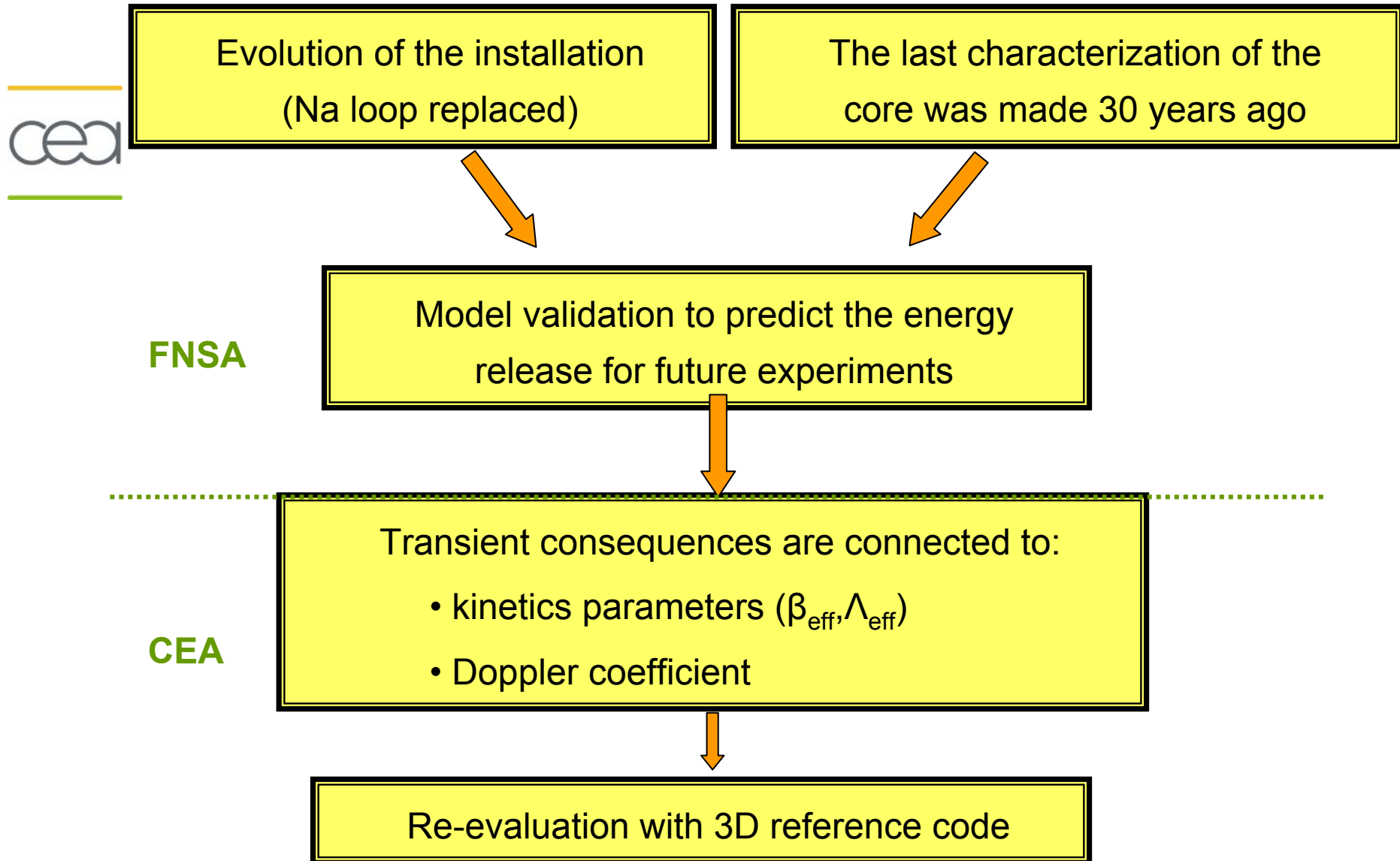
t > 0.1s.

- ➔ T°clad, T°coolant increase
- ➔ Minor "Delayed" feedback



Automatic shutdown ~0.1 – 0.2 s.
by dropping all the control rods
→ End of the transient

What was the problematic in the safety report (2008)?



Neutronics characterizations in CABRI core by MC calculation

Doppler coefficient



Using the reactivity determined at different fuel temperatures by the TRIPOLI4

code: $\Delta\rho^{\text{doppler}}(t) = A_D \cdot \sqrt{\Delta T} \approx 103 \text{ pcm} \cdot K^{-0.5}$

Effective generation time (Λ_{eff})

Using a method based on perturbation theory with MC calculations.

The results are compared with experimental measurements and deterministic computations.

	Expected values	MCNP results
MASURCA (MUSE4 experiment)	$56 \pm 5 \mu\text{s}$	$53 \pm 5 \mu\text{s}$
OSIRIS (deterministic calculation)	$34.7 \pm 2 \mu\text{s}$	$34.0 \pm 1.1 \mu\text{s}$
JHR (deterministic calculation)	$36.0 \pm 2 \mu\text{s}$	$38.7 \pm 1.4 \mu\text{s}$

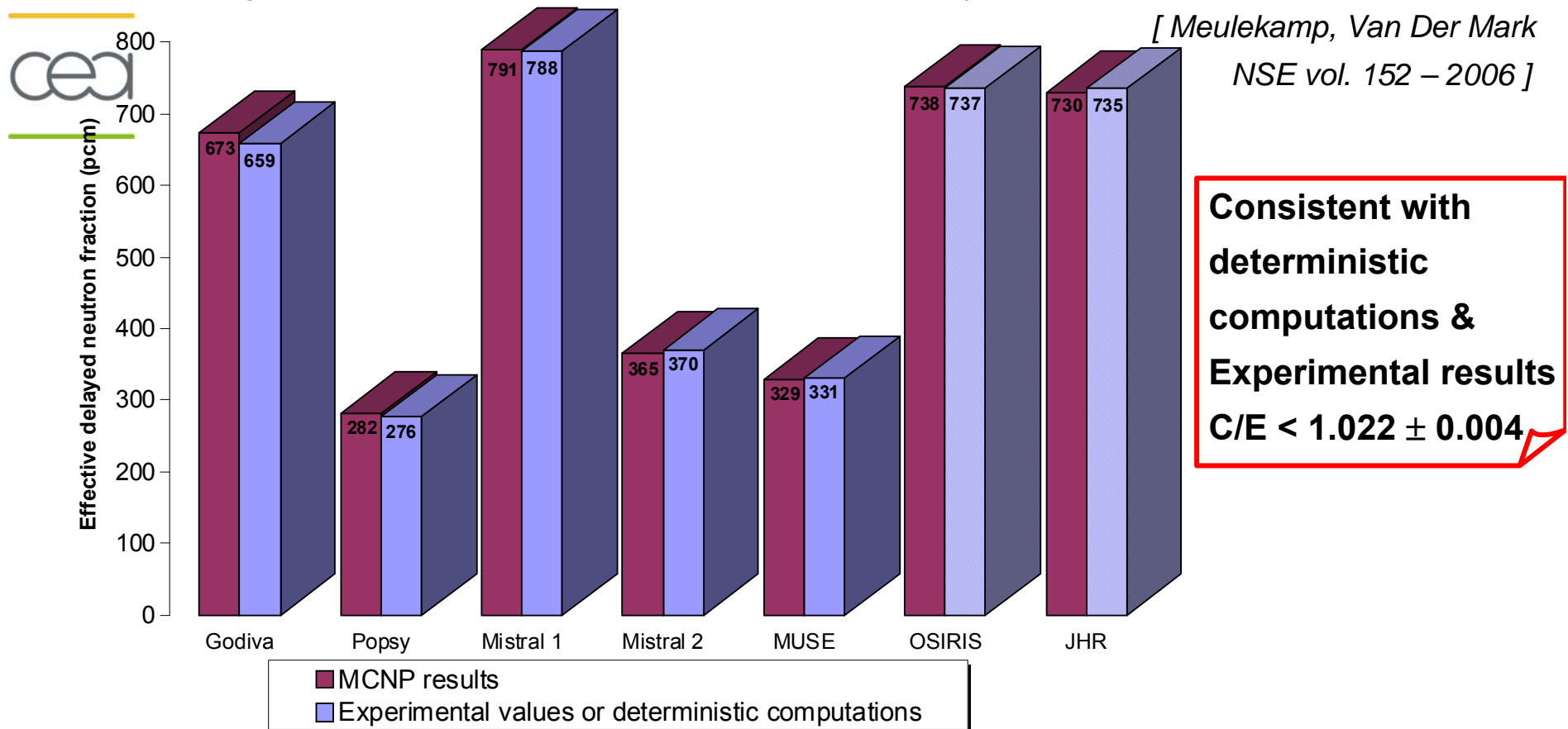
Agreement MCNP results / expected values allows us to trust this method

The application to CABRI core (JEFF-3.1 library) : $\Lambda_{\text{eff}} = 27.7 \pm 2 \mu\text{s}$

Neutronics characterization in CABRI core by MC calculation

Effective delayed neutron fraction (β_{eff})

Using a method based on iterated fission probability and available in a MCNP patch

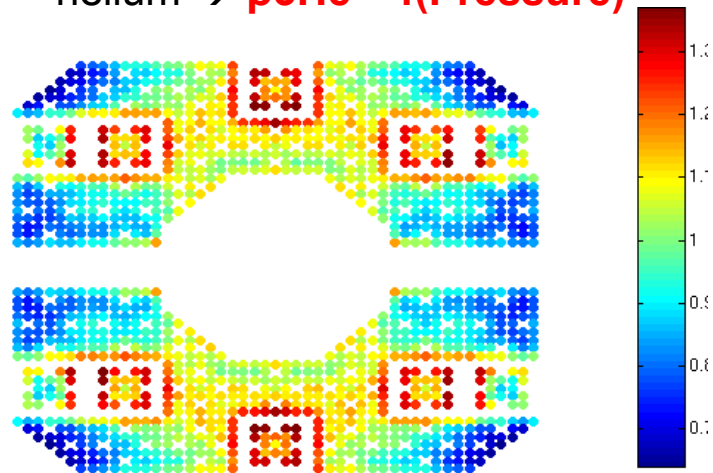


The application to CABRI core (JEFF-3.1 library) : $\beta_{eff} = 758 \pm 2 pcm$

Integral validation : Reactivity calculation

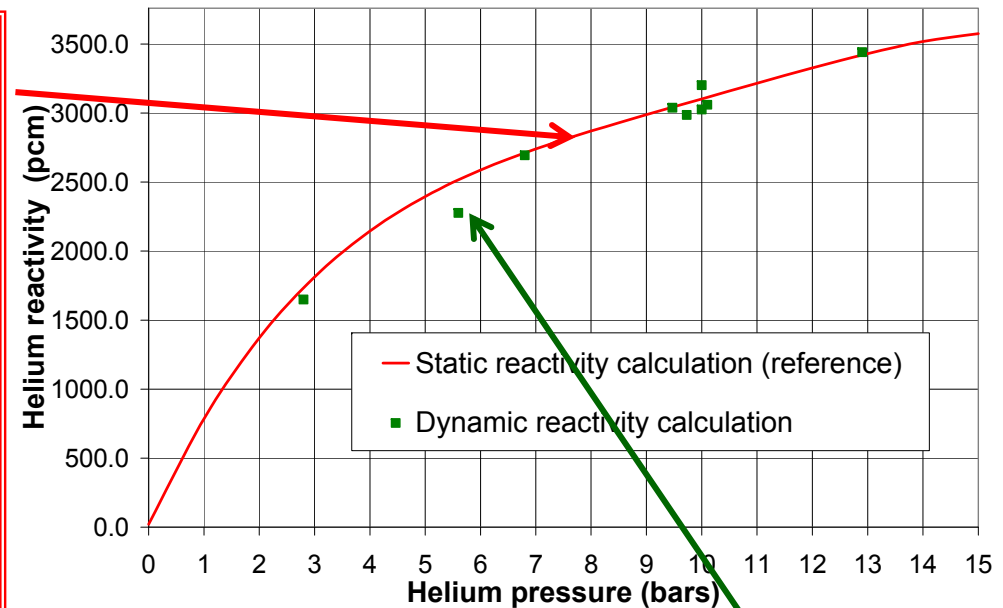
A test is characterized by the injected Helium reactivity

Static calculations with and without helium $\rightarrow \rho^{3\text{He}} = f(\text{Pressure})$



Confident in this result because:

- ✓ TRIPOLI4 is a reference code and helium 3 cross section are well known
- ✓ TRIPOLI4 reactivity calculations are validated (for different helium pressure, the predicted critical position by the code is always consistent with the experimental data)



Dynamic Calculations (using β_{eff} , Λ_{eff} , Doppler)

Experimental power bursts obtained in the past can be interpreted by inverse kinetics. One can determine the injected helium reactivity for different tests (performed at different helium pressure) $\rightarrow \rho^{3\text{He}} = f(\text{Pressure})$

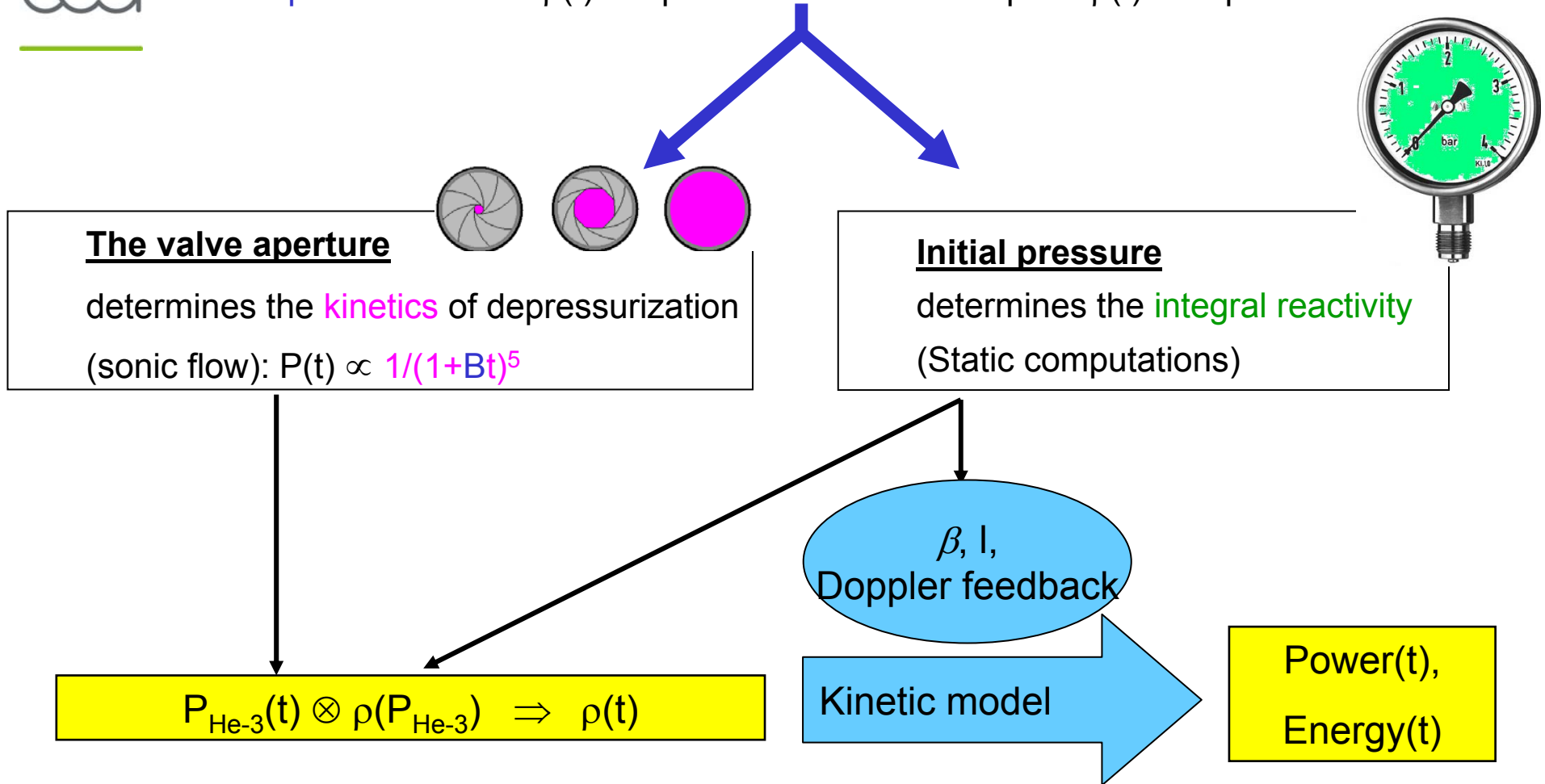
The 2 independent methods are consistent \rightarrow the kinetic model is validated

Predictive methodology definition

- **Problem:** to simulate the pulse power, the model need an input data: Reactivity (ρ).



- **Principle:** To evaluate $\rho(t)$ the phenomena are decoupled: $\rho(t)$ is dependent on



Prospective methodology validation

To assess the accuracy of the predicted energy values, the methodology is applied to the simulation of several tests performed in the past



Test	Energy Measurement (MJ)	Energy Prediction (MJ)	Discrepancy (%)
Test 1	209	195	-6
Test 2	204	197	-3
Test 3	212	199	-6
Test 4	213	203	-4
Test 5	224	227	1
Test 6	172	176	2

Maximum Discrepancy ~ 6%

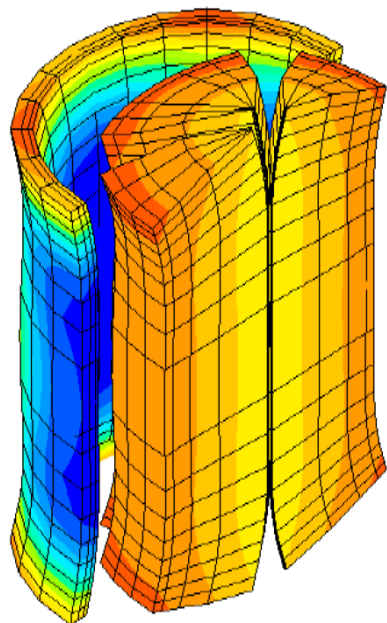
Good agreement between measured and predicted energy better than 6%.

Safety Assessment

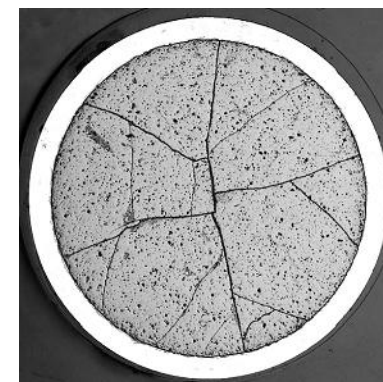
The same predictive approach is used to characterize the energy release during the most severe possible accident



	Peak Power (GW)	Energy release (MJ)
Reference Accident 15 bars-valves at full aperture	38.6	382 +6% → 403



Thermo-mechanical computations
SCANAIR code (provided by IRSN)



- attests that the consequences are acceptable to both fuel and cladding.

Conclusion

Reevaluation

✓ The main neutronics parameters driving a power pulse were re-evaluated by 3D Monte Carlo reference codes.

Validation

✓ The code including thermal and point kinetics models is validated.

Predictive methodology

✓ The comparison between predictions and experimental results shows the model predictions are better than 6%.

Safety assessment

✓ This model is used to assess the power and energy release during the reference accident of the CABRI facility: computations attest the fuel and cladding integrity during the transient.



**Thank you
for your
attention**

ANNEXE 1 : Calcul du β_{eff}

$$\beta_{eff} = \frac{\langle \Phi^* \chi_d \nu_d \Sigma_f \Phi \rangle}{\langle \Phi^* \chi \nu \Sigma_f \Phi \rangle}$$



Dans un calcul Monte-Carlo : Pas de modèle physique, pas d'équation : la notion de flux adjoint Φ^* (ou d'importance neutronique) n'existe pas.

Alternatives :

- On supprime Φ^* \rightarrow on calcule β_0 (\sim environ 10 % inférieur à β_{eff})
- Méthode prompte : approximation $\Phi = \Phi_p$ et $\Phi^* = \Phi_p^* \rightarrow \beta_{eff} \cong 1 - \frac{k_p}{k}$
avec $k_p = k_{eff}$ si $\chi = \chi_p$ et $\nu = \nu_p$.
- Méthode NRG \rightarrow on estime Φ^* par IFP ou plus précisément NFP : MCNP stocke la « longueur d'histoire » afin de restituer la notion d'importance.

Calcul MCNP CABRI – bib. JEFF3.1	
β_0	716 pcm
β_{eff} (méthode prompte)	762 pcm
β_{eff} (méthode NRG)	758 pcm

Calcul du temps de génération (Λ_{eff})

$$\Lambda_{eff} = \frac{\langle \Phi^*, \frac{1}{v} \Phi \rangle}{\langle \Phi^*, F \Phi \rangle}$$



Théorie des perturbations :
$$\Delta\rho = \frac{\langle \Phi_0^*, [\lambda_0 \Delta F - \Delta K] \Phi_1 \rangle}{\langle \Phi_0^*, F_1 \Phi_1 \rangle}$$

Cas particulier : ajout uniforme d'un poison de section efficace microscopique de capture en c/v

- $\Delta K = c/v$: l'absorption se rajoute aux causes de disparitions des neutrons ;
- $\Delta F = 0$: pas de modification de l'opérateur production.

$$\Delta\rho = \rho_c - \rho_0 = -\frac{\langle \Phi_0^*, \frac{c}{v} \Phi_c \rangle}{\langle \Phi_0^*, F_0 \Phi_c \rangle}$$

$$\left(\frac{\partial \rho}{\partial c} \right)_{c=0} = \lim_{c \rightarrow 0} \frac{\rho_c - \rho_0}{c} = \lim_{c \rightarrow 0} \frac{1}{c} \frac{\langle \Phi_0^*, \frac{c}{v} \Phi_c \rangle}{\langle \Phi_0^*, F_0 \Phi_c \rangle} = \frac{\langle \Phi_0^*, \frac{1}{v} \Phi_0 \rangle}{\langle \Phi_0^*, F_0 \Phi_0 \rangle}$$

$$\Lambda_{eff} = -\frac{\Delta\rho}{c}$$

Compromis : c doit tendre vers 0 tout en permettant une précision statistique sur $\Delta\rho$ suffisante. « Verboomen et al. » préconisent $\Delta\rho \sim 300$ pcm

Rq : le temps de génération est accessible avec T4 (conforme à MCNP)

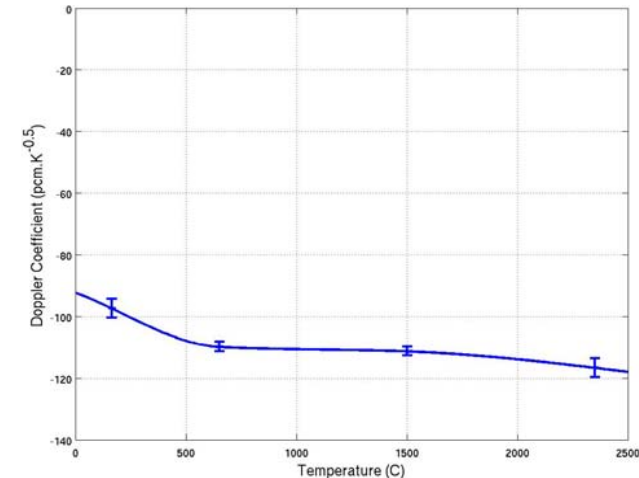
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