

PWR 3D-Core Simulation using different ATF fuels

M. Muñoz et al. (Grupo INGENIA 2019-2020)

MASTER IN INDUSTRIAL ENGINEERING

E.T.S. de Ingenieros Industriales

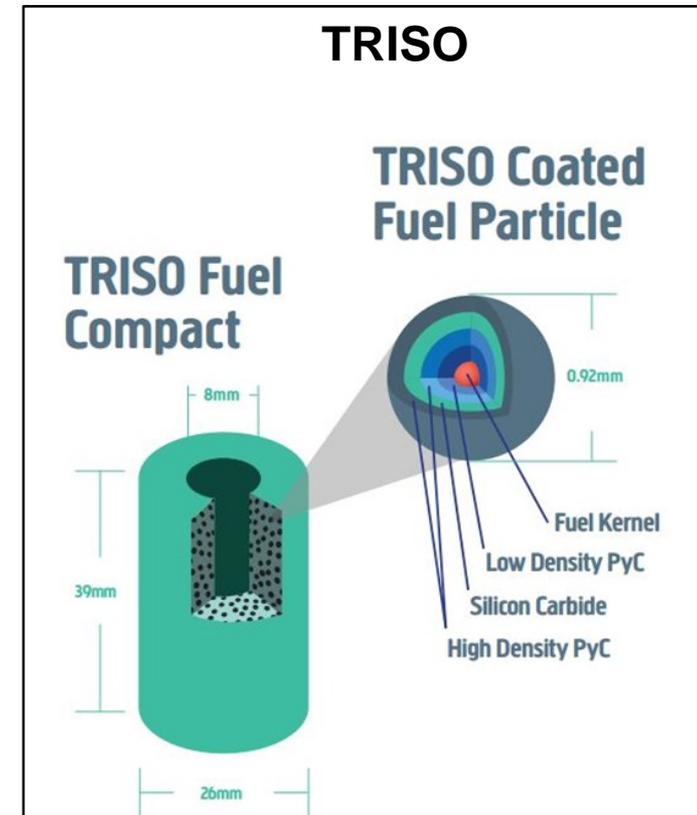
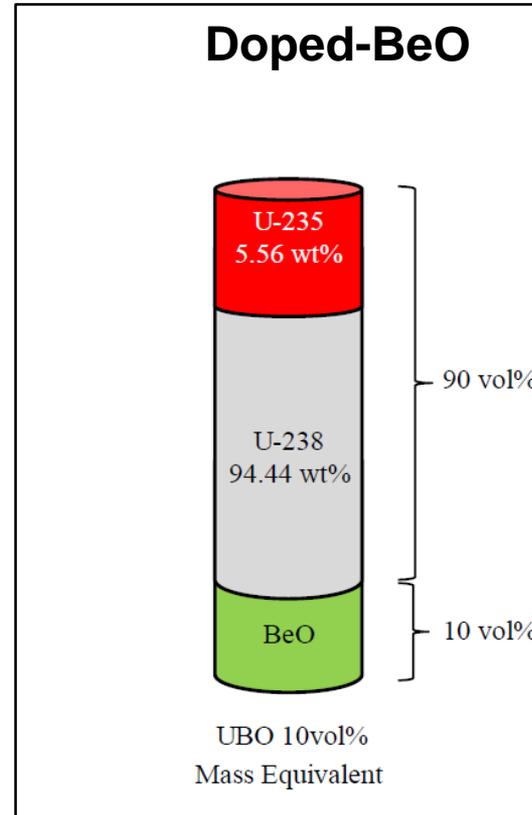
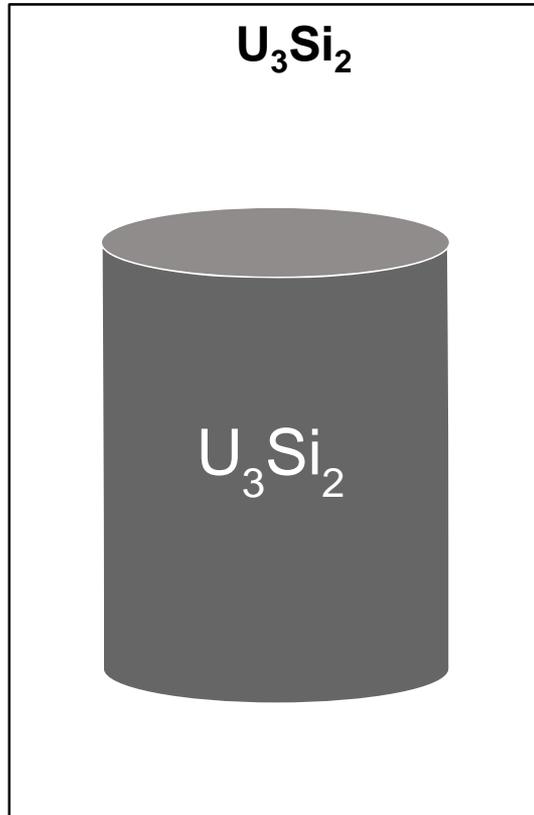
Universidad Politécnica de Madrid (UPM), Madrid, Spain

O. Cabellos

Universidad Politécnica de Madrid (UPM), Madrid, Spain

E-mail: oscar.cabellos@upm.es

□ ATF Fuels

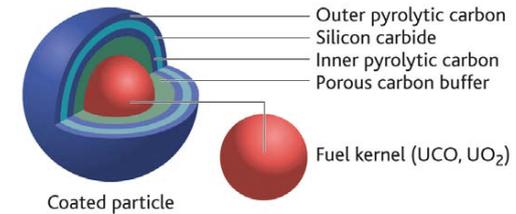


Ref. Neutronbytes, "TRISO Fuel Drives Global Development of Advanced Reactors" (2017)

Ref. J. R. Smith, "Enhanced Thermal Conductivity UO₂-BeO Nuclear Fuel: Neutronic performance studies and economic analyses, MSc (2012)

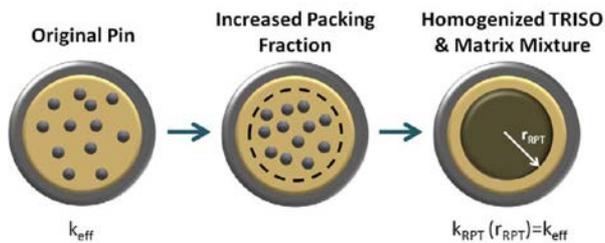
Multi-coated TRISO fuel particles

- Double-heterogeneity arises from the distribution of TRISO particles within the graphite matrix
- The packing or volume factor



Methodology Reactivity-equivalent Physical Transformation (RPT) Pin Cell

The RPT model creates a single-heterogeneous problem out of double-heterogeneous one



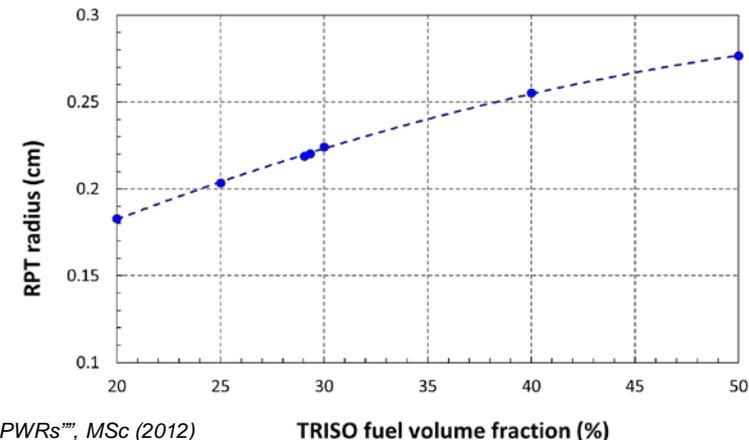
RPT method of fuel zone homogenization

- TRISO particles are condensed in the center of the fuel zone
- Then, CSi and TRISOs are homogenized into a single mixture

Calculation of RPT Radius.

- SCALE/KENO PWR Pin-cell with TRISO UO₂/wo-VF%
- WIMSD5 equivalent pin-cell (densities,....)

Figure 1. RPT radius TRISO/Csi-UO₂-20.0%wt



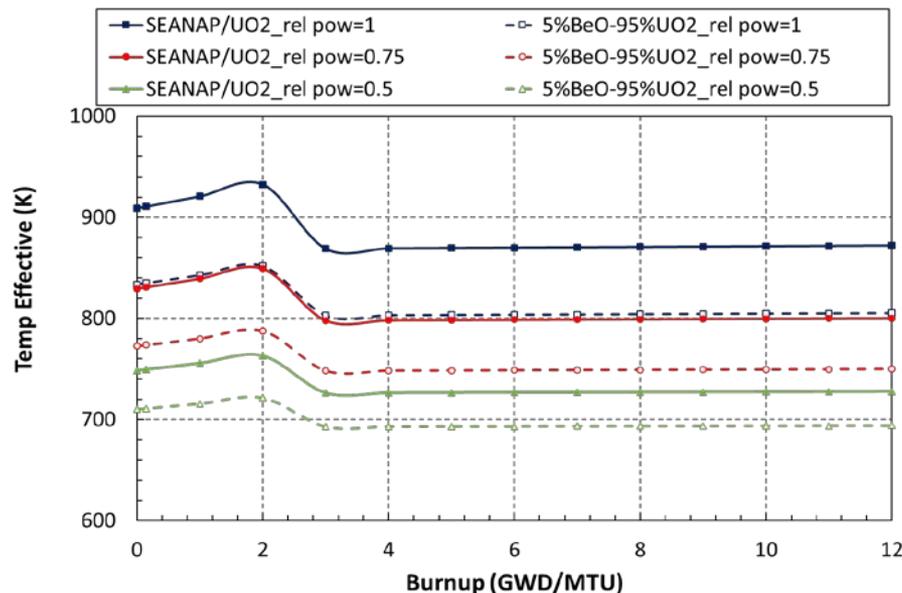
Ref. Steven T. Lynch, "Reactivity-Equivalent Physical Transformation Model for Pin Cell Arrays", MSc (2010)

Ref. C.A. Gentry, "An investigation of the use of Ceramic Microencapsulated Fuel for Transuranic Waste Recycling in PWRs", MSc (2012)

The “Effective” fuel-temperature

- Assumption in Lattice code: the temperature profile across a fuel pin is spatially flat
- Then, we need an “effective Doppler temperature”. This value can be calculated using different approaches such that the Doppler feedback is properly accounted for
- For instance, the “NEA” effective fuel temperature: $T_{NEA}=0.7 \cdot T_S+0.3 \cdot T_C$
- $T_{eff}=f(\text{fuel, relative power, burnup})$
 - T_S = temperature pellet surface
 - T_C = temperature pin center-line

Figure 2. Teff as a function of burnup



Adjusted Teff value for Doped-BeO is based on temperature profiles of this work:

Ref. J. R. Smith, “Enhanced Thermal Conductivity UO₂-BeO Nuclear Fuel: Neutronic performance studies and economic analyses, MSC (2012)

□ Modelisation of ATF Fuels: %wt enrich. for “equivalent cycle”

□ TRISO ATFs

- 84 FCM (CSi-UO₂-5%wo-80%VF) fuel rods loaded on the periphery of the fuel assembly
- 180 remaining fuel of enriched UO₂ fuel rods

□ U₃Si₂ and Doped (UO₂-BeO) ATFs

- 264 fuel rods

Figure 3. CORAIL type-fuel assembly

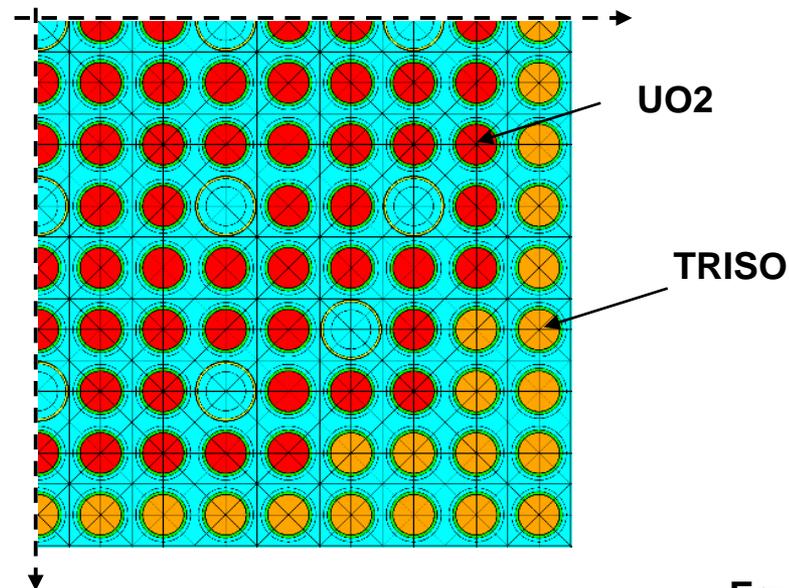
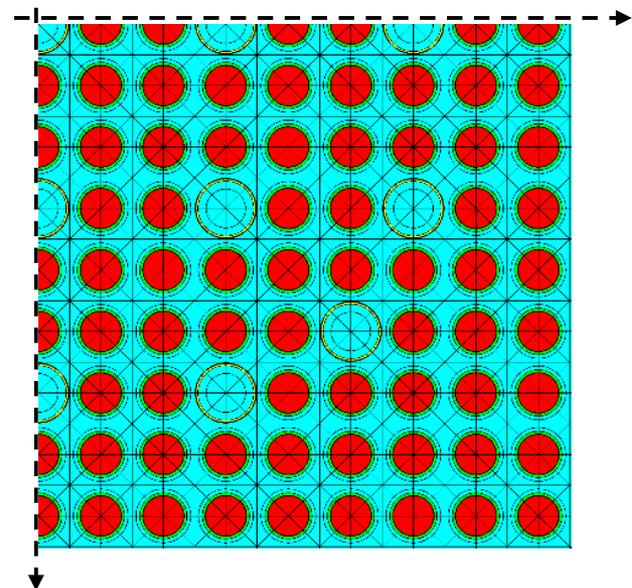


Figure 4. MIX type-fuel.

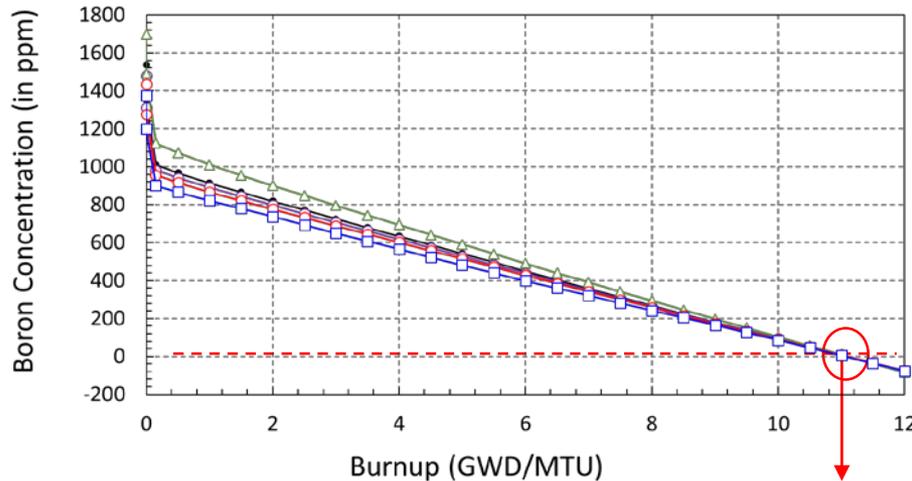
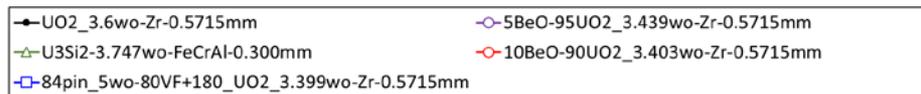


**Equivalent
Enrichment ?**

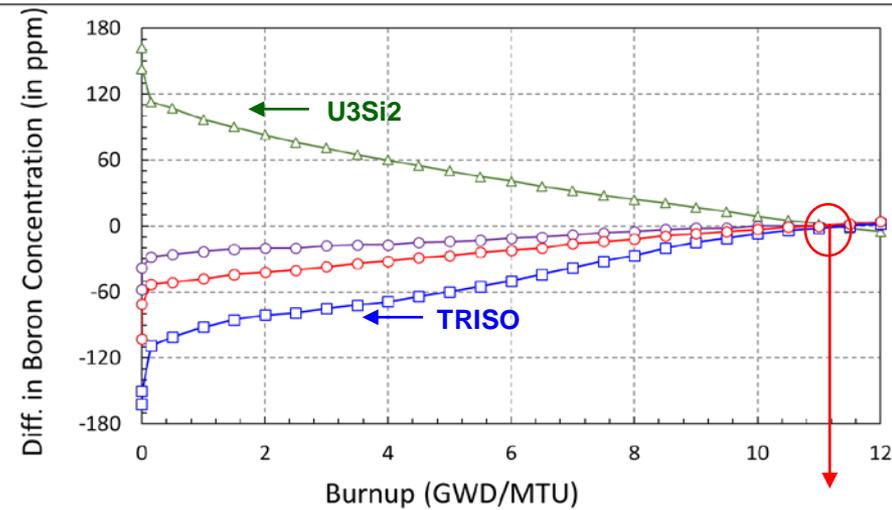
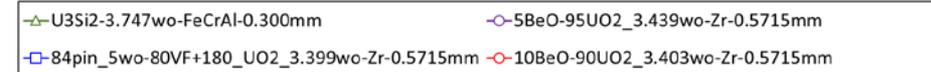
□ Core average burnup analysis with “equivalent cycle length”

Table 1. Enrichment for each technology

	UO ₂ -Zr	U ₃ Si ₂ -FeCrAl	TRISO-Zr	5% BeO + 95% UO ₂ -Zr	10% BeO + 90% UO ₂ -Zr
Enrichment (% wo)	3.60	3.75	3.40	3.44	3.40



Cycle length ~ 11.08 GWD/MTU



Cycle length ~ 11.08 GWD/MTU

□ Axial Offset

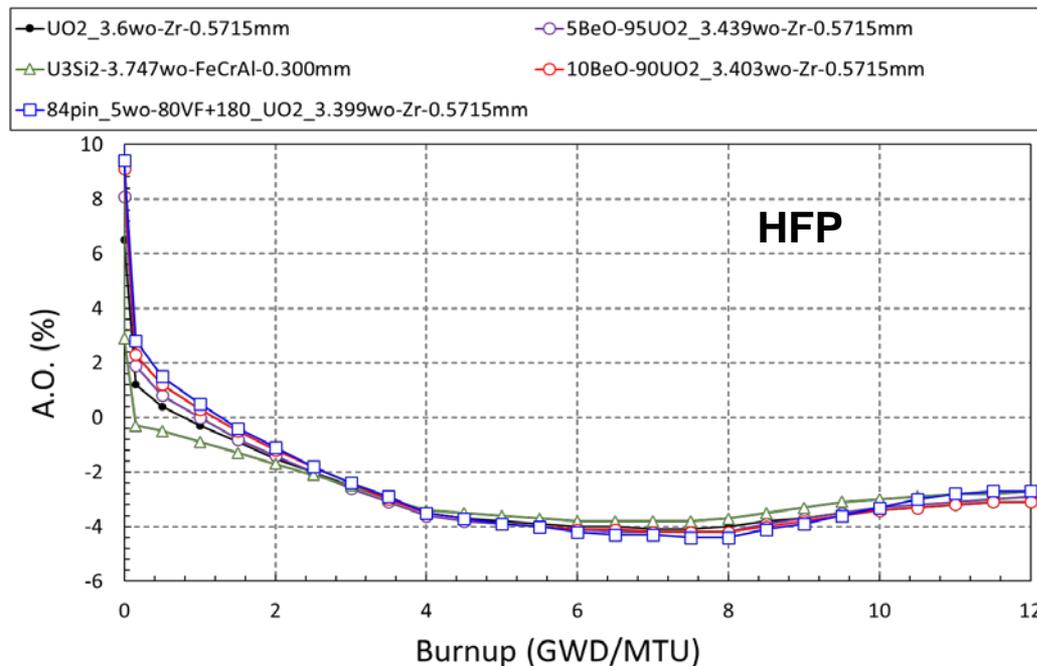
$$\text{Axial Offset (\%)} = \frac{P_T - P_S}{P_T + P_S} * 100$$

P_T is the power fraction in the top-half of the core

P_S is the power fraction in the bottom-half of the core

Table 2. A.O.(%) at BOC/HZP

Case	A.O. (%) BOC/HZP
UO ₂ /3.60wo – Zr/0.5715mm	33.9
U ₃ Si ₂ /3.747wo – FeCrAl-300μm	29.2
95%UO ₂ /3.439wo+5%BeO – Zr/0.5715mm	36.2
90%UO ₂ /3.403wo+10%BeO – Zr/0.5715mm	37.8
TRISO – Zr/0.5715mm 84FCM: 5%wo-80%VF 180UO ₂ : 3.4wo	40.0



□ Heat flux hot channel factor (F_Q)

$$F_Q = \frac{\text{max. local LHGR}}{\text{average LHGR}} = \max[Q(z)]$$

LHGR is Linear Heat Generation Rate (W/cm)

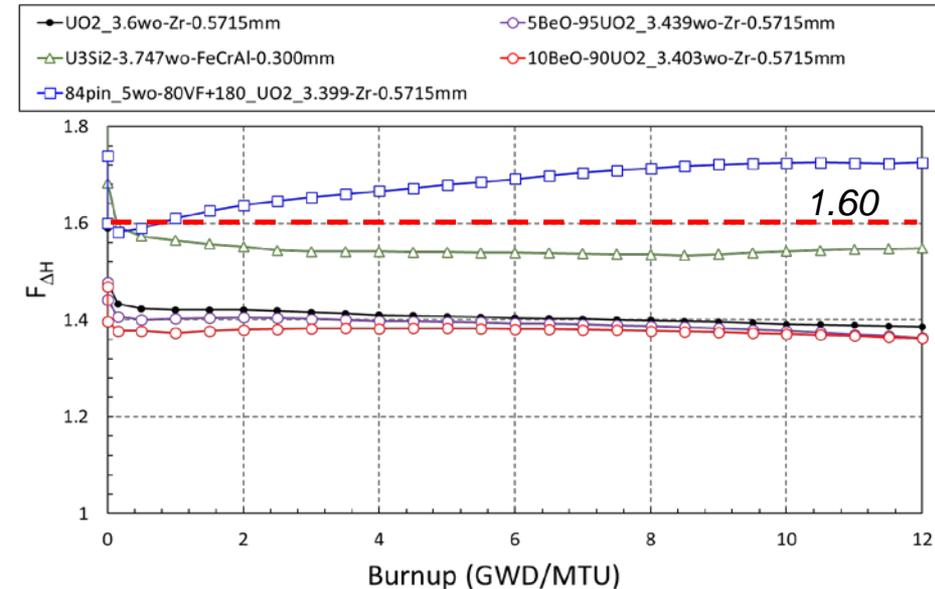
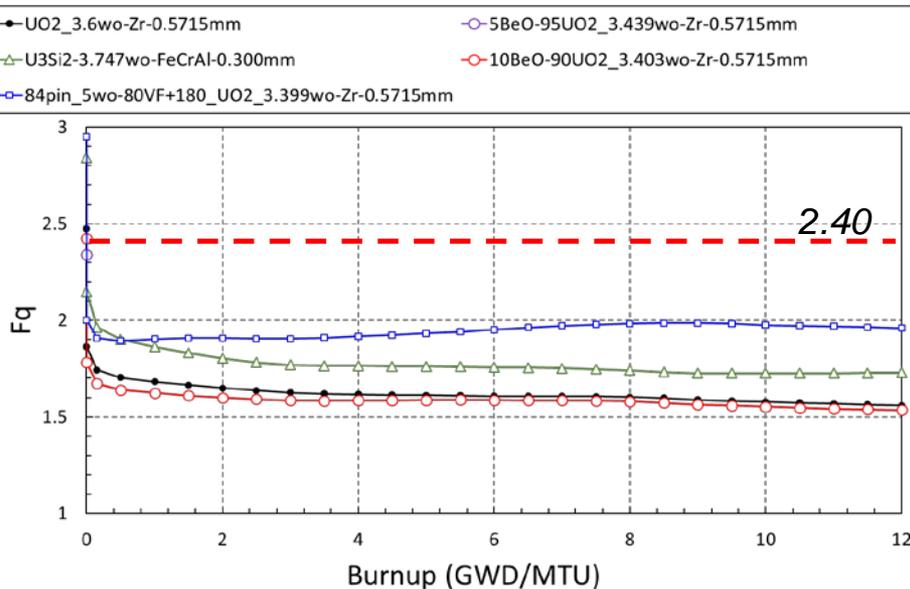
$Q(z)$ is the maximum linear power at elevation z

Design Limit: $F_Q \cdot P < F_Q^{\text{max}} = 2.40$

□ Enthalpy Peaking factor ($F_{\Delta H}$)

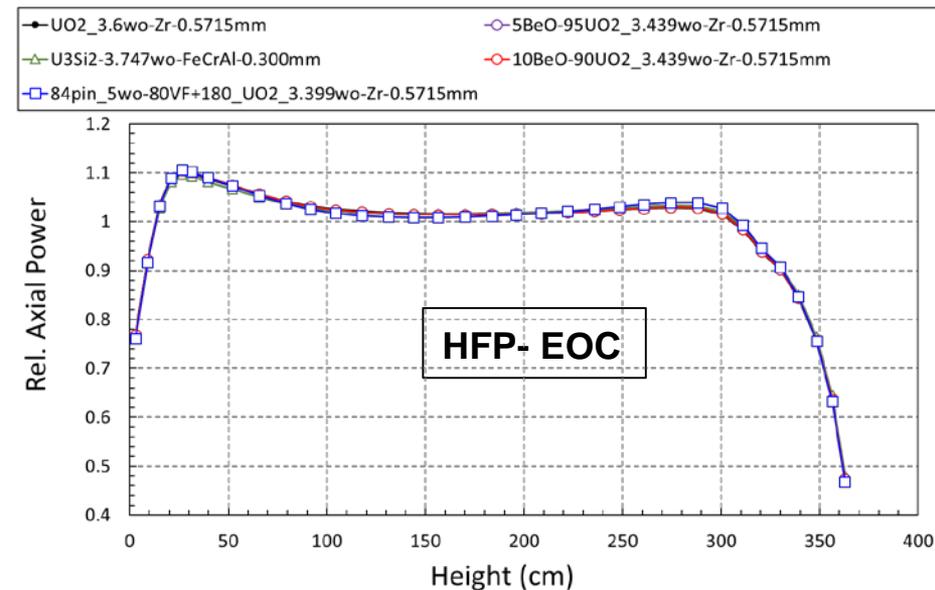
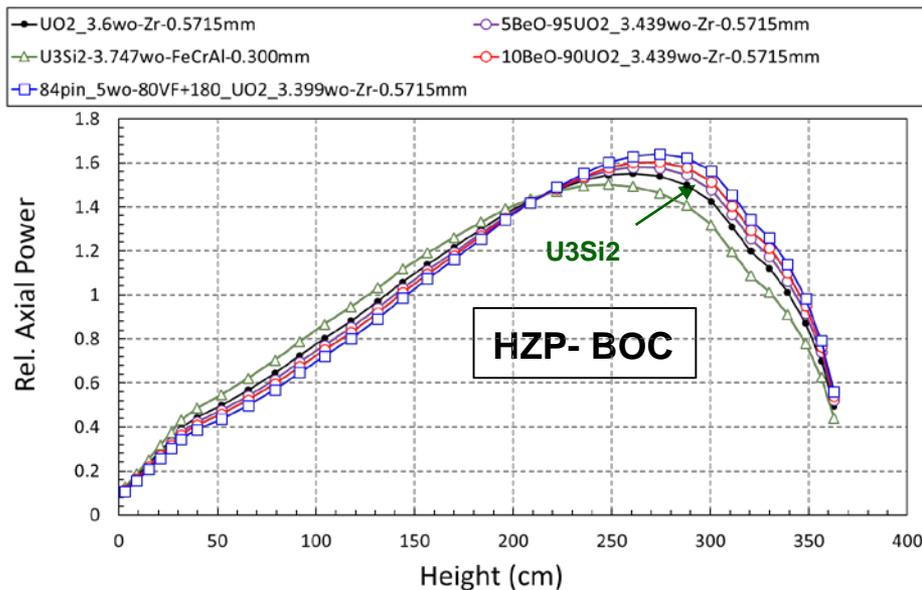
$$F_{\Delta H} = \frac{\text{max. channel enthalpy rise}}{\text{core average enthalpy rise}} = \max\left[\frac{\Delta H}{\text{avg } \Delta H}\right]$$

Design Limit: $F_{\Delta H} < F_{\Delta H}^{\text{max}} \approx 1.60 \cdot [1 + 0.3(1-P)]$



□ Axial power distribution: HZP/HFP – BOC/EOC

- Axial power distribution change with cycle depletion
- At BOC/HZP is peaked at the bottom of the core
 - $\text{UO}_2/\text{Zr} \Rightarrow \text{A.O.}(\%) = 33.9\%$
 - $\text{U}_3\text{Si}_2 \Rightarrow \text{A.O.}(\%) = 29.2\%$
 - TRISO $\Rightarrow \text{A.O.}(\%) = 40.0\%$
- At EOC, the axial power becomes flattened



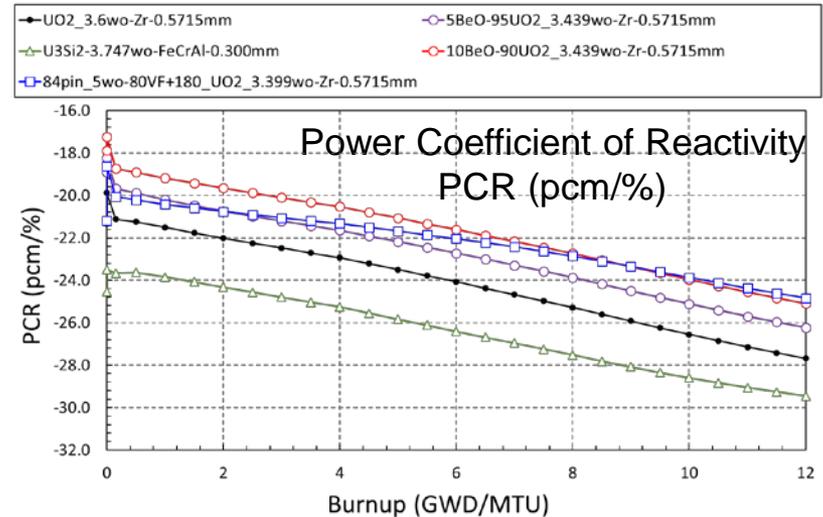
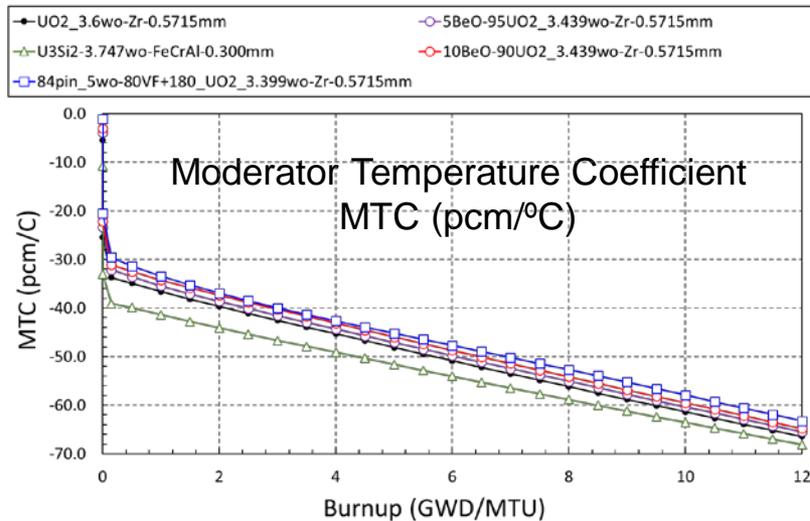


VALBOR = Boron worth coefficient
 MTC = Moderator temperature coefficient
 DOP = Power coefficient of reactivity (only Doppler)

FTC = Fuel temperature coefficient
 PCR = Power coefficient of reactivity

Table 3. Reactivity coefficients at BOC/HZP

Reactivity Coefficients at BOC/HZP	Boron Worth (pcm/ppm)	MTC (pcm/°C)	FTC (pcm/°C)	CISO (pcm/°C)	DOP (pcm/%)	PCR (pcm/%)
UO ₂ /3.60wo – Zr/0.5715mm	-8.3	-5.5	-5.1	-10.5	-16.0	-21.3
U ₃ Si ₂ /3.747wo – FeCrAl-300μm	-7.3	-10.8	-5.1	-15.9	-15.4	-24.6
95%UO ₂ /3.439wo+5%BeO – Zr/0.5715mm	-8.6	-3.9	-5.1	-9.1	-14.4	-18.9
90%UO ₂ /3.403wo+10%BeO – Zr/0.5715mm	-8.8	-3.0	-5.3	-8.3	-13.8	-17.9
TRISO – Zr/0.5715mm 84FCM: 5%wo-80%VF 180UO ₂ : 3.4wo	-9.1	-1.0	-5.7	-6.8	-17.6	-21.2





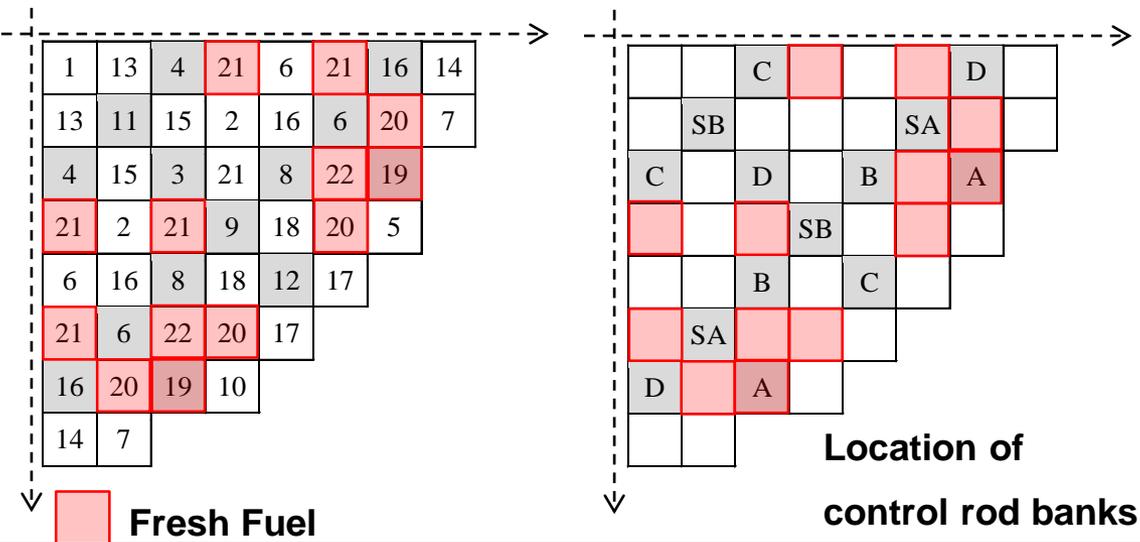
Bank Worths

- Note: the worth of control bank rods are calculated at HZP/BOC using Rod Swap and boron dilution technique.

Table 4. Bank Worth (in Boron ppm) at HZP/BOC

1/4 CORE

1/4 CORE



Control rod bank worth (in ppm)	UO2 3.6wo	U3Si2 3.747wo	95%UO2_3.439wo +5%BeO	90%UO2_3.403wo +10%BeO	TRISO 84FCM: 5%wo-80%VF 180UO2: 3.4wo
	Zr-0.5715mm	FeCrAl-300mm	Zr-0.5715mm	Zr-0.5715mm	Zr-0.5715mm
BANK D-in	114	118	111	109	106
BANK C-in	87	80	96	102	107
BANK B-in	135	143	127	120	110
BANK A-in	90	106	71	62	51
BANK SB-in	86	78	98	106	113
BANK SA-in	120	138	102	92	80
BANK D+C-in	221	220	228	231	232
BANK D+C+B-in	400	403	403	402	396
BANK D+C+B+A-in	542	565	517	502	481
BANK D+C+B+A+SA-in	664	674	657	653	637
ARI (all rods in)	884	910	865	854	829

□ Core average burnup analysis with “equivalent cycle length ”

➤ Impact nuclear data: JEFF-3.3, JEFF-3.1, JEF-2.2, ENDF/B-VII.1 and ENDF/B-VIII.0

Table 5. Value of wt% for the ATF to maintain an “equivalent cycle” length as UO₂/3.6wo fuel

Nuclear Data Evaluations	U3Si2	95%UO ₂ +5%BeO	90%UO ₂ +10%BeO	TRISO 84rods 5%wo-80%VF
	FeCrAl-300μm	Zr-0.5715mm	Zr-0.5715mm	Zr-0.5715mm
JEFF-3.3	3.75	3.44	3.40	3.40
JEFF-3.1	3.75	3.43	3.39	3.38
JEF-2.2	3.76	3.47	3.47	3.39
ENDF/B-VII.1	3.76	3.44	3.40	3.41
ENDF/B-VIII.0	3.76	3.44	3.40	3.40



□ **Conclusions**

➤ **ATFs may get simulated in SEANAP**

- an updated Teff correlation will be needed

➤ **TRISO must be specially adapted to get simulated**

- an optimal layout must be found to make it suitable in our simulation

➤ **General speaking good performance for doped fuels, uranium silicide and TRISO**

➤ **Assessing the impact of nuclear data**

- ND Evaluations: JEFF-3.3, JEFF-3.1, JEF-2.2, ENDF/B-VII.1 and ENDF/B-VIII.0
- low differences are found searching the %wt for the “equivalent cycle”
- slightly differences were found in JEF-2.2 due to Be nuclear data?
- other core parameters to be investigated: power distributions, etc...
- impact of new nuclear data: new Thermal Scattering Laws for graphite in TRISO?