

Comparison of Core Design Parameters for BANDI-60 Using UO₂ and U-Mo Fuels

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1. Introduction

The small modular reactor (SMR) has several advantages such as power supply to remote locations, seawater desalination, and marine propulsion [1]. Recently, the BANDI-60 [2] SMR of 200 MWth developed by KEPCO E&C is designed as a core with 52 fuel assemblies (FAs) using UO₂ fuel enriched to 4.95% and the Pyrex burnable absorber (BA) for excess reactivity control [3]. In the core, five types of FAs in the same number of BA rods of 24 with different concentrations of Pyrex are loaded to have core characteristics of its cycle length of about 4.9 years and the maximum excess reactivity of around 3800 pcm [4].

In this paper, the core design parameters of Bandi-60 with changing the UO₂ fuels into U-Mo are compared with the existing ones. The main core design target is to extend the cycle length by using the U-Mo fuel. Since the U-Mo fuel has a higher uranium density than UO₂ fuel, U-Mo fuel can make the cycle length longer with the same amount of uranium enrichment [5, 6]. For the comparison of core design parameters, the core burnup calculation is performed by the Monte Carlo particle transport analysis code, McCARD [7]. For the two cores, their neutronics parameters such as the effective multiplication factor (k_{eff}), the power peak factor, and the temperature coefficient are compared.

2. Core Design of SMR

2.1 Fuel materials

Table I shows a comparison of information of UO₂ and U-Mo fuels. U-Mo fuel has a higher thermal conductivity than UO₂ fuel, which leads to a lower maximum fuel temperature, while the melting temperature of U-Mo is lower than UO₂ [5, 6]. Because U-Mo fuel has a high density, using it enables to load more uranium into the core.

Table I: Information for UO₂ and U-Mo fuels

Division	UO ₂ fuel	U-Mo fuel
Thermal Conductivity [W/m/K] (@ 377 °C)	4.78	25.6
Melting Temperature [°C]	2,865	1134
Density [g/cm ³]	Uranium	9.04
	The others	1.22
	Total	10.26
Initial uranium loading in BANDI-60 [kg]	11,894	22,497
Uranium enrichment	4.95 wt%	

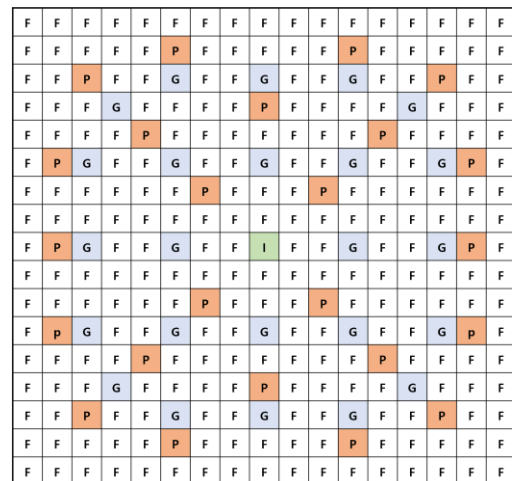
2.2 Core design parameters

Table II shows the design parameters of SMR [3, 4]. The thermal power of the core is 200 MW. FA is based on the Westinghouse 17×17 FA. 52 FAs are loaded in the core. The average linear power density is 7.28 kW/m, which is lower than that of commercial PWR and there is no soluble boron in the coolant.

Table II: The design parameters of SMR

Parameters	Value
Reactor type	PWR
Thermal power	200 MW
Average Linear power density	7.28 kW/m
Coolant & Moderator	Light water
Coolant Average Temp.	580.65 K
Number of FAs	52
Active core height	200 cm
FA pitch / Pin pitch	21.50 cm / 1.26 cm
FA type	Westinghouse 17×17
Uranium enrichment	4.95 wt%
Fuel Material	UO ₂ or U-Mo
BA material	Pyrex

The core is designed using FA loaded with Pyrex. As shown in Figure 1, BA rods of 24 are loaded instead of fuel pins. The neutron absorption capacity of BA is proportional to the weight percent (w/o) of B₂O₃ contained in Pyrex [3]. Figure 2 shows the cross-sectional view of the fuel pellet and BA.



F: Fuel Pin (240) / P: Pyrex (24) / I: Instrument Tube (1) / G: Guide Tube (24)

Fig. 1. Fuel assembly configuration

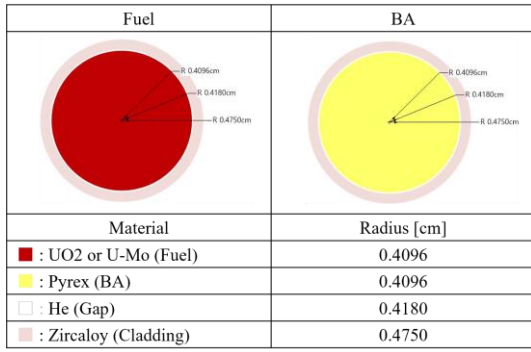


Fig. 2. The cross-sectional view of the fuel and BA.

2.3 Loading Pattern of the Core

Table III shows the information of FA types by the concentration of BA. FAs using a high concentration of B₂O₃ are placed in the center of the core and FAs using a low concentration of B₂O₃ are placed in the periphery to make power distribution smooth in the radial direction [3, 4]. Figure 3 presents the loading pattern of SMR. In the soluble boron-free SMR, the core needs a large amount of the reactivity control mechanism instead of soluble boron, and the reactivity change should be controlled by using the control rods. As shown in Figure 4, there are forty (40) Control Element Assemblies (CEAs) in the core for the reactivity control [2]. The location of CEAs is determined to control for normal operation and provide sufficient control rod worth to overcome the reactivity feedback caused by the core state change [3, 4].

Table III: FA types by the concentration of BA

B ₂ O ₃ w/o in Pyrex	Number of fuel pins	Number of BA pins	Number of FA
5	240	24	8
10	240	24	12
25	240	24	16
35	240	24	12
40	240	24	4
Total	12,480	1,248	52

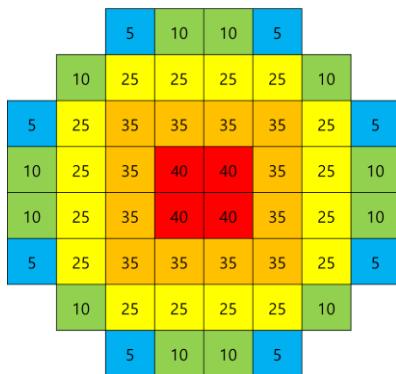
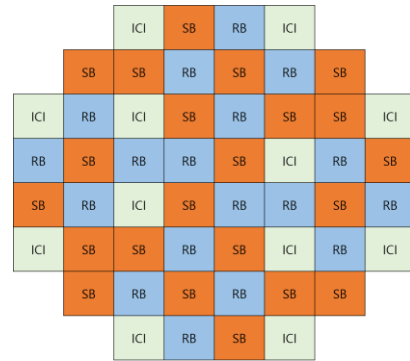


Fig. 3. The loading pattern of SMR



* RB: Regulating Bank (18), SB: Shutdown Bank (22), ICI: In-Core Instrumentation (12)

Fig. 4. The location of the CEAs

3. Numerical Results

The McCARD burnup calculation is conducted with 50,000 or 100,000 histories per cycle on 150 inactive and 300 active cycles using the continuous-energy cross section libraries produced from ENDF/B-VII.1. Table IV shows the McCARD burnup calculation options.

Table IV: The McCARD burnup calculation options

Division	UO ₂ core	U-Mo core
Neutron histories	100,000	50,000
Active/inactive cycle	300 / 150	300 / 150
Fuel Avg. Temp.	700 K	700 K
Moderator Avg. Temp.	580.65 K	580.65 K
Moderator density	0.7105 g/cm ³	0.7105 g/cm ³

3.1 The effective multiplication factor (k_{eff})

Figure 5 shows k_{eff} vs. the effective full power day (EFPD) behavior of UO₂ and U-Mo cores. The calculated maximum cycle lengths for UO₂ and U-Mo cores are $1,784 \pm 6$ and $2,905 \pm 10$ days, respectively. The excess reactivity for the U-Mo core is calculated as $7,080 \pm 15$ pcm, which is $3,291 \pm 15$ pcm more than the UO₂ core. Table V shows the difference in results about EFPD, burnup, and the excess reactivity.

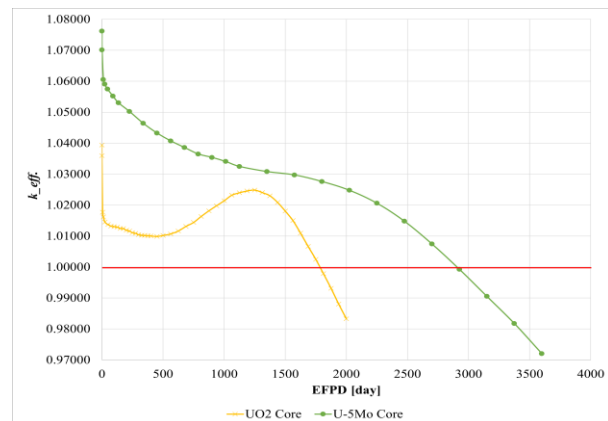


Fig. 5. k_{eff} vs. EFPD behavior of UO₂ and U-Mo cores

Table IV: The calculation results of UO₂ and U-Mo cores

Division	UO ₂ Core	U-Mo Core
Max. EFPD [day]	1,784 ± 6	2,905 ± 10
Max. burnup [MWD/kgU]	29.99 ± 0.09	25.82 ± 0.09
Excess reactivity [pcm]	3,790 ± 10	7,080 ± 18
BOC k_{eff}	1.03939 ± 0.00012	1.07620 ± 0.00017

3.2 The power peaking factor

The assembly power peaking factor (Fr) for both cores are compared in Figure 6. Both graphs decrease from BOC to the middle of the cycle (MOC) and reach their highest value at the end of the cycle (EOC). The maximum Fr values for UO₂ and U-Mo cores are 1.355 ± 0.002 and 1.328 ± 0.003 , respectively.

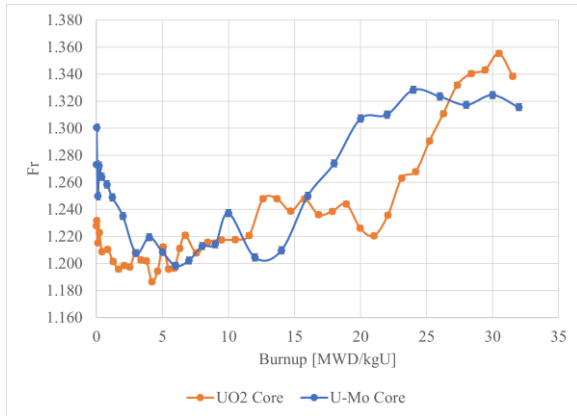


Fig. 6. Fr vs. burnup behavior of UO₂ and U-Mo cores

Figure 7 shows the pin power peaking factor (Fq) vs. burnup of UO₂ and U-Mo cores. Both cores show similar Fq behavior over the cycle, with the U-Mo core maximum peaking factors being slightly higher within the error bars indicating 1 σ statistical uncertainty in calculated results. The maximum Fq values for UO₂ and U-Mo cores are 2.33 ± 0.10 and 2.57 ± 0.17 , respectively.

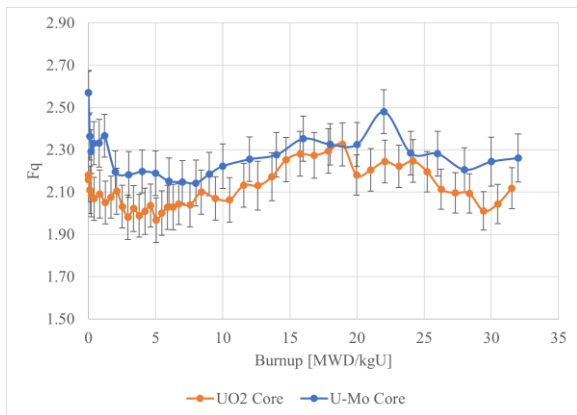


Fig. 7. Fq vs. burnup behavior of UO₂ and U-Mo cores

3.3 The temperature coefficient

Table IV shows the McCARD burnup calculation options for the temperature coefficients of both cores. The fuel temperature coefficients (FTCs) for UO₂ and U-Mo cores are compared in Figure 8. The U-Mo core FTCs are more negative than the UO₂ core FTCs over the cycle.

Table V: The calculation options for FTC & MTC

Division	UO ₂ Fuel		U-5Mo Fuel	
	FTC	MTC	FTC	MTC
Delta fuel temp. [K]	200	0	50	0
Delta moderator temp. [K]	0	7.5	0	7.5
Fuel avg. temp. [K]	900	700	650	700
Moderator avg. temp. [K]	580.65	573.15	580.65	573.15
Moderator density [g/cm ³]	0.7105	0.7270	0.7105	0.7270
Neutron histories	100K	50K	50K	25K
Active cycle	300	300	300	300
Inactive cycle	150	150	150	150

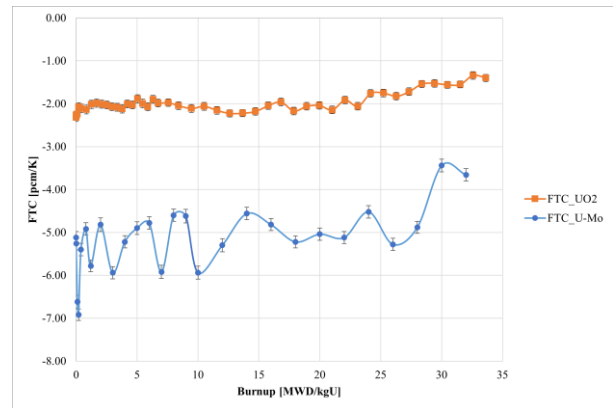


Fig. 8. FTC vs. burnup of UO₂ and U-Mo cores

Figure 9 shows a comparison of the moderator temperature coefficient (MTC) for both cores. In BOC, the UO₂ core MTC is more negative than the U-Mo core MTC, and vice versa in EOC.

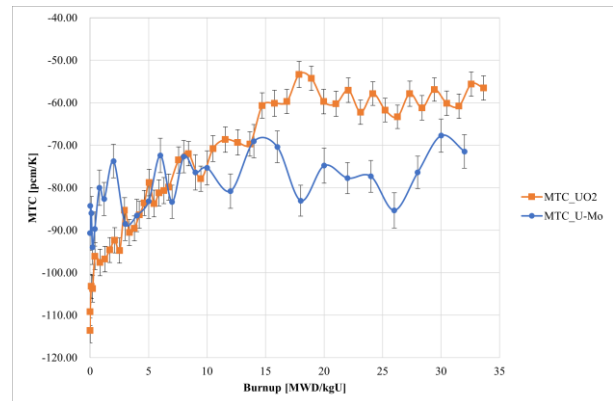


Fig. 9 MTC vs. burnup of UO₂ and U-Mo cores

4. Conclusions

In this paper, the core design parameters of the Bandi-60 SMR using UO_2 and U-Mo fuels enriched to 4.95% are compared. The main design target is to extend the cycle length of the core by using U-Mo fuel. This study presents that the cycle length of the U-Mo core is $2,905 \pm 10$ days, which is about 1.63 times that of the UO_2 core. Because the U-Mo fuel is a higher density than the UO_2 fuel, the core contains a large amount of uranium, which can extend the cycle length. The maximum excess reactivity for the U-Mo core is $7,080 \pm 15$ pcm, which is about 3,290 pcm higher than that of the UO_2 core. The maximum F_q values for the UO_2 and U-Mo cores are 2.33 ± 0.10 and 2.57 ± 0.17 , respectively. The U-Mo core has more negative FTC than the UO_2 core.

As for future work, the core design adopts a newly proposed BA concept for Gd_2O_3 (gadolinia), the cylindrically inserted and mechanically separated burnable absorber (CIMBA) [8].

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