

# Interpolating Fuel Cycle Behavior for Scenario Analysis Codes: Thermal IMF Recycle

Tracy Radel

Kara Beharry

*University of Wisconsin – Madison*

Taek K. Kim

*Argonne National Laboratory*





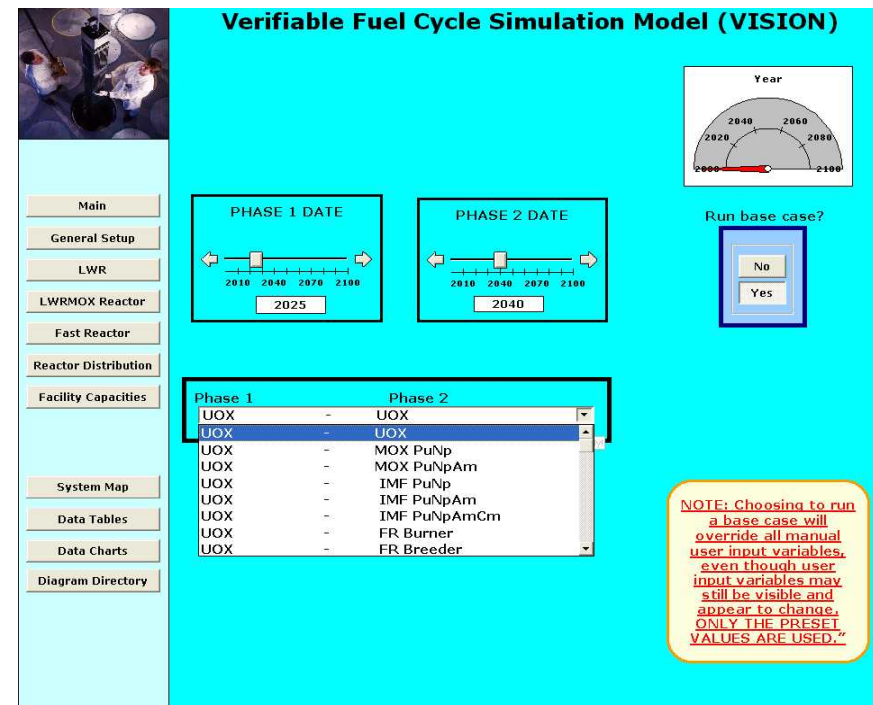
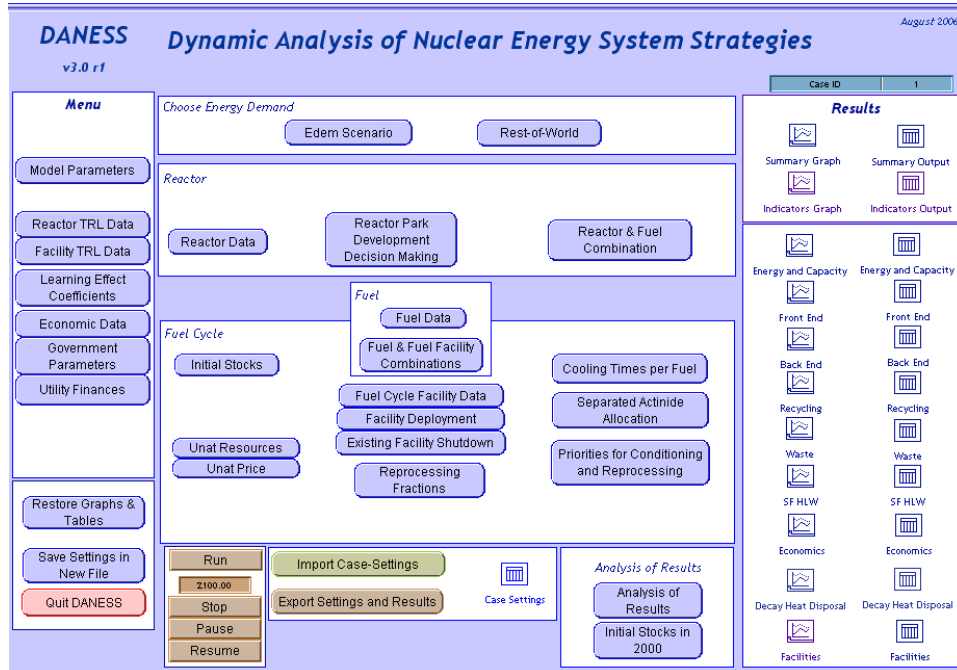
# Outline



- Scenario Analysis Codes
- Methodology
- Inert Matrix Fuel (IMF) Material Description
- Assembly Geometry
- IMF Cycle Description
- Initial Interpolation Scheme
- Revised Interpolations Scheme
- Conclusions

# Scenario Analysis Codes

- DANESS (ANL)
- VISION (INL)
- Model mass flows throughout entire fuel cycle
- Use STELLA and Powersim software

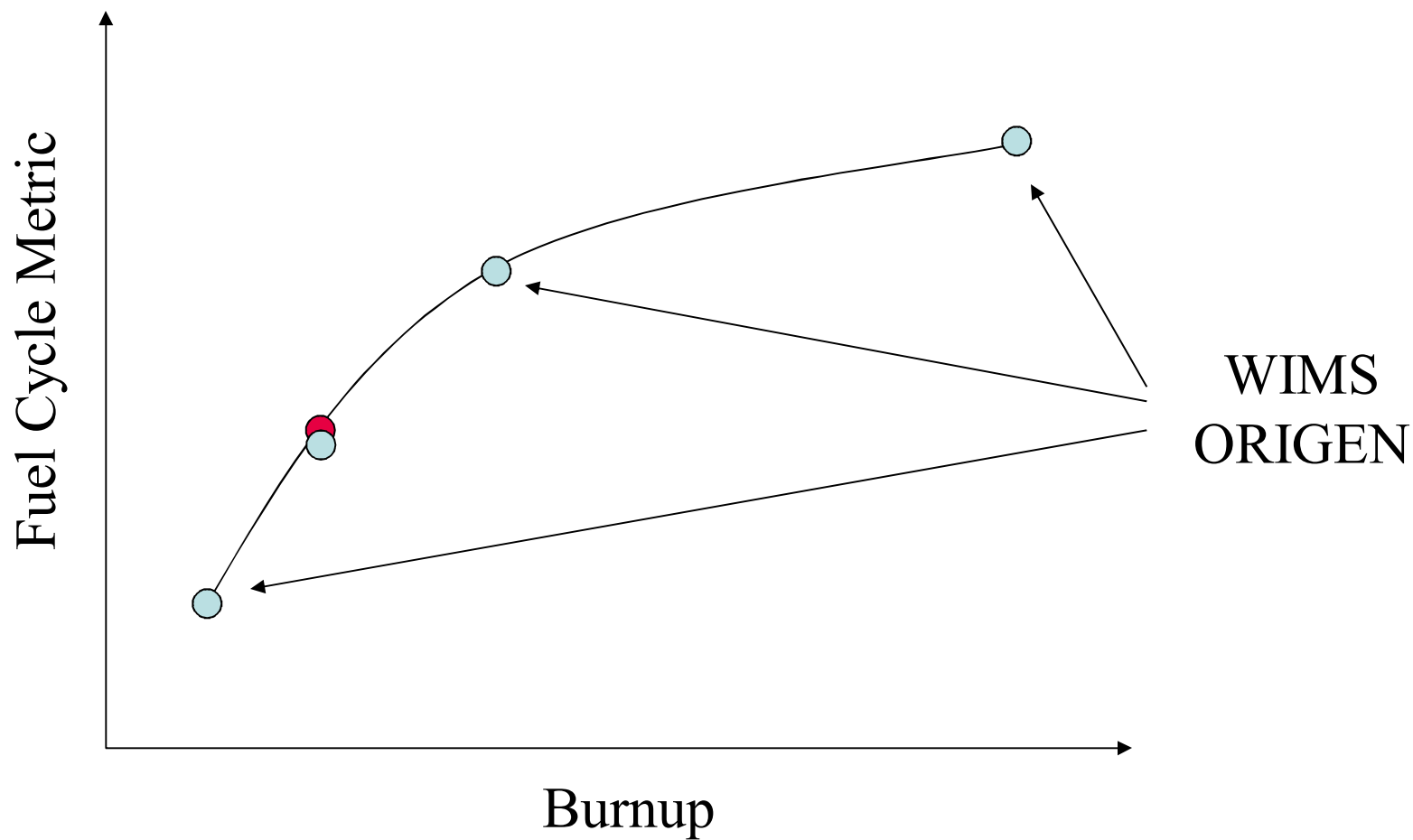




# Scenario Analysis Codes (Cont.)



- Accurate mass flows require knowledge of
  - $\text{TRUO}_2$  Mass Fraction
  - Charge Mass
  - Discharge Mass
  - Blending Ratios
  - Charge Vector
  - Discharge Vector
- Performance requirements rule out detailed physics modules
- Therefore, require simplified interpolation equations





# Inert Matrix Fuel (IMF)



- LWR-IMF system is one of nuclear fuel cycles considered in scenario study under AFCI project of US-DOE
- Pu, Np, and Am extracted from spent nuclear fuel
- Placed in an inert matrix material which is:
  - Neutron transparent
  - Chemically stable
  - Radiation damage resistant
  - Economically reasonable
- Zirconia stabilized by yttrium oxide and combined with spinel to compensate for zirconia matrix's low conductivity

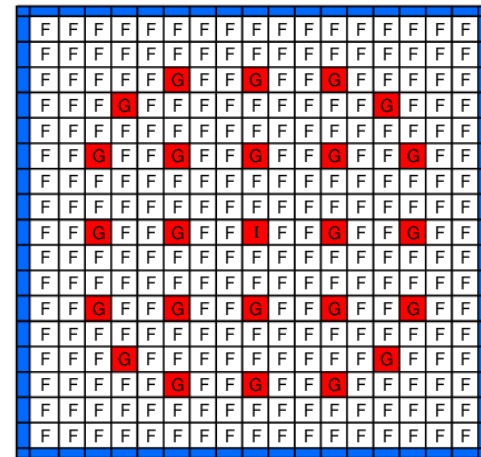
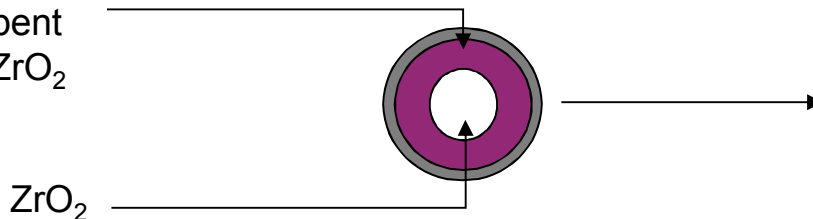


# IMF Fuel Assembly Geometry



- Homogeneous Fuel Assemblies
  - IMF fuel pins are located at all fuel pins positions of 17x17 typical PWR assembly
  - IMF fuel pins are made by blending TRUs of previous IMF cycle and LWR spent fuel
  - Blending ratio and  $\text{TRUO}_2$  mass fraction are key parameters to maintain a desired cycle length

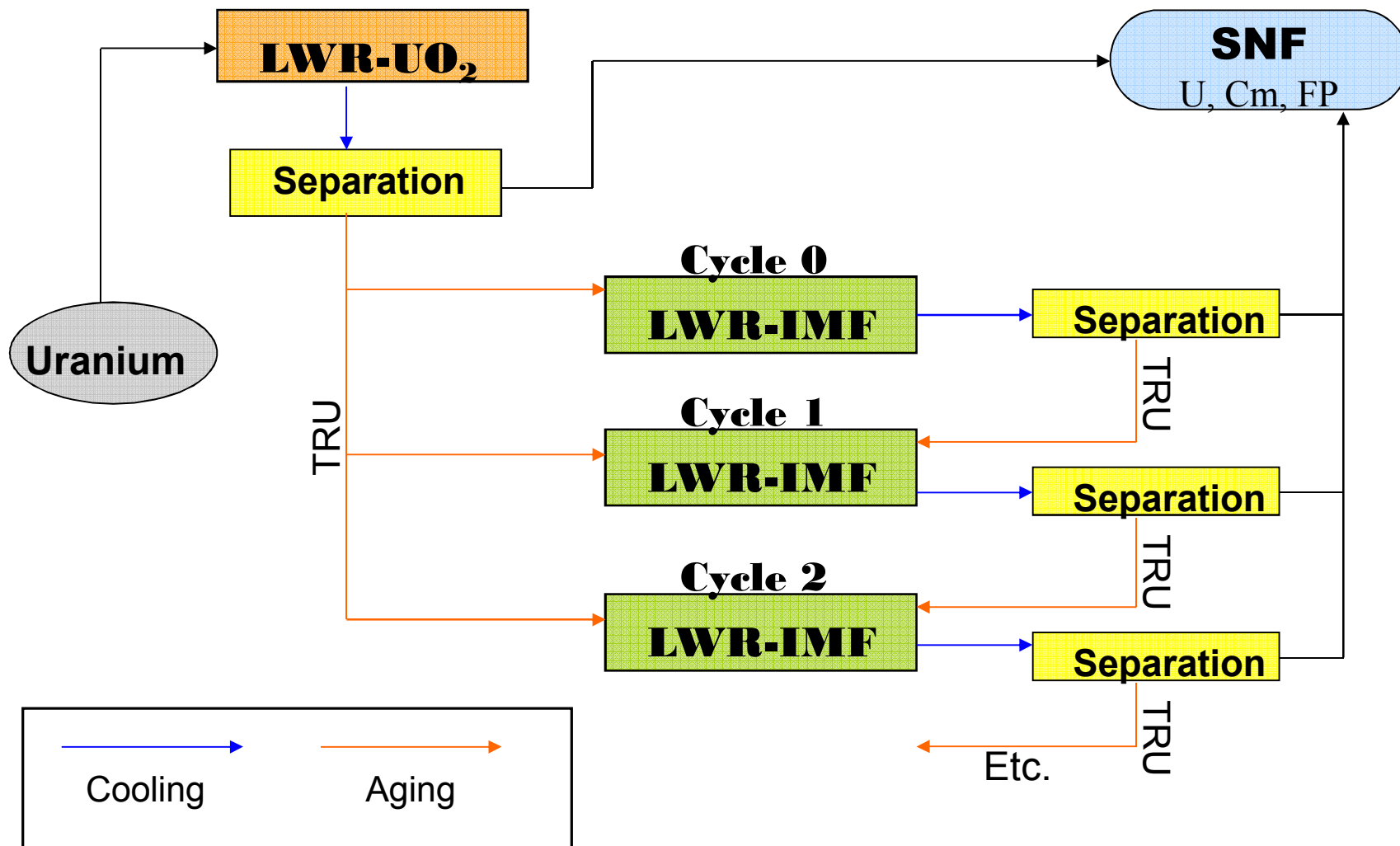
TRU of previous  
IMF cycle, TRU  
of LWR spent  
fuel, and  $\text{ZrO}_2$



**Homogeneous IMF Assembly**



# IMF Cycle







# IMF Calculations



- $\text{TRUO}_2$  mass fraction
- Radius of zirconium center

$$r_{\text{Zr}} = -0.0019 * MF_{\text{TRUO}} + 0.2453$$

- 95% theoretical density

$$\rho_{\text{IMF}_{95}} = \frac{0.95 * V_{ff} \rho_{\text{HMO}} \rho_{\text{ZrO}}}{[MF_{\text{TRUO}} * \rho_{\text{ZrO}} + (1 - MF_{\text{TRUO}}) * \rho_{\text{HMO}}] (1 + D_f)}$$

- Charge mass

$$M_{\text{charge}} = MF_{\text{TRUO}} MF_{\text{TRU}} V_A * \rho_{\text{IMF}_{95}}$$

- Discharge mass
- Blending ratio
- Critical burnup

$$B_c = \frac{n+1}{2n} B_d$$



# Initial IMF Fuel Cycle Interpolation Scheme

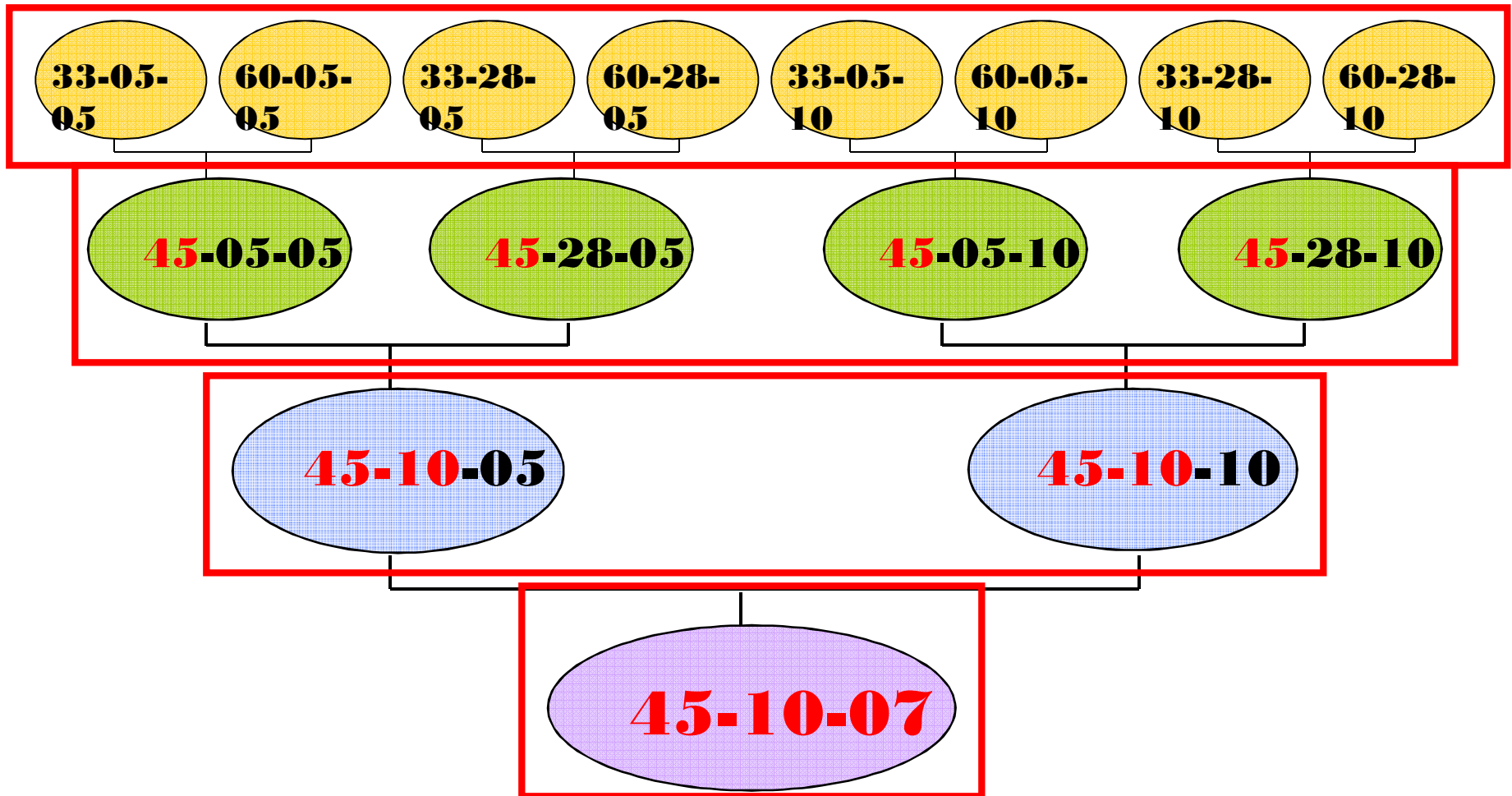


- Three independent variables
  - Burnup of  $\text{UO}_2$  fuel
    - 33, 60, and 100 GWd/t
  - Cooling time after initial  $\text{UO}_2$  cycle
    - 5, 28, and 50 years
  - Cooling time between IMF cycles
    - 5 and 10 years
- Five cycles for each data set
- Two test cases
  - 45 GWd/t burnup, 10 year initial cooling, 7 year IMF cooling
  - 75 GWd/t burnup, 45 year initial cooling, 7 year IMF cooling



## Interpolation Scheme (Cont.)

Example: QUANTILE REGRESSION with  $\lambda_1 = 0$  and  $\lambda_2 = 0.001$  (Cholobing, 2015)  $\rightarrow$  COVE cooling





# Initial Homogeneous Assembly Results



- TRUO<sub>2</sub> Mass Fractions

45 GWd/t burnup, 10 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	Error (%)
Cycle 0	24.21	23.75	1.95
Cycle 1	35.40	37.21	4.86
Cycle 2	44.01	46.70	5.75
Cycle 3	50.48	53.45	5.55
Cycle 4	55.46	58.50	5.19

75 GWd/t burnup, 45 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	Error (%)
Cycle 0	59.61	59.61	0.01
Cycle 1	63.93	64.28	0.54
Cycle 2	68.72	69.13	0.60
Cycle 3	72.47	72.87	0.55
Cycle 4	75.35	75.78	0.56

- Errors higher with lower cooling time due to Pu-241 decay to Am-241
- 57.8% and 93.6% of 109 predicted values had errors less than 5% for 45 and 75 GWd/t burnup cases, respectively



# Effects of Pu-241



- Pu-241 half-life: 14.4 years
- Decays to Am-241
- Fissile isotope to strong absorber
- Mass change varies greatly
  - 5 years: 78.61% remains
  - 28 years: 26.61% remains
  - 50 years: 9.01% remains
- **Cooling times of approximately 10.6 (60%) and 17.5 (43%) years should be added**



# Revised IMF Fuel Cycle Interpolation Scheme



- Three independent variables
  - Burnup of  $\text{UO}_2$  fuel
    - 33, 60, and 100 GWd/t
  - Cooling time after initial  $\text{UO}_2$  cycle
    - 5, **10**, **17**, 28, and 50 years
  - Cooling time between IMF cycles
    - 5 and 10 years
- Five cycles for each data set
- Two test cases
  - 45 GWd/t burnup, **14** year initial cooling, 7 year IMF cooling
  - 75 GWd/t burnup, **40** year initial cooling, 7 year IMF cooling



# Improved Results



- Case 45-14-07
  - Average error on  $\text{TRUO}_2$  mass fraction: 1.3%
  - Overall
    - Values with less than 5% error: 102 of 109
    - Average error: 1.7%
    - Maximum error: 6.1%
- Case 75-40-07
  - Average error on  $\text{TRUO}_2$  mass fraction: 1.1%
  - Overall
    - Values with less than 5% error: 104 of 109
    - Average error: 1.4%
    - Maximum error: 4.9%



# Conclusions and Recommendations



- Mass flows of LWR-IMF fuel cycles were calculated using WIMS9
- Efficient interpolation schemes were proposed to predict system study variables
- Interpolations can now be used to accurately estimate various values within scenario analysis codes





Produced by University Communications

# Questions?

Special thanks to:

Kara Beharry

Taek Kim

Abdellatif Yacout

Paul Wilson





# Backup slides

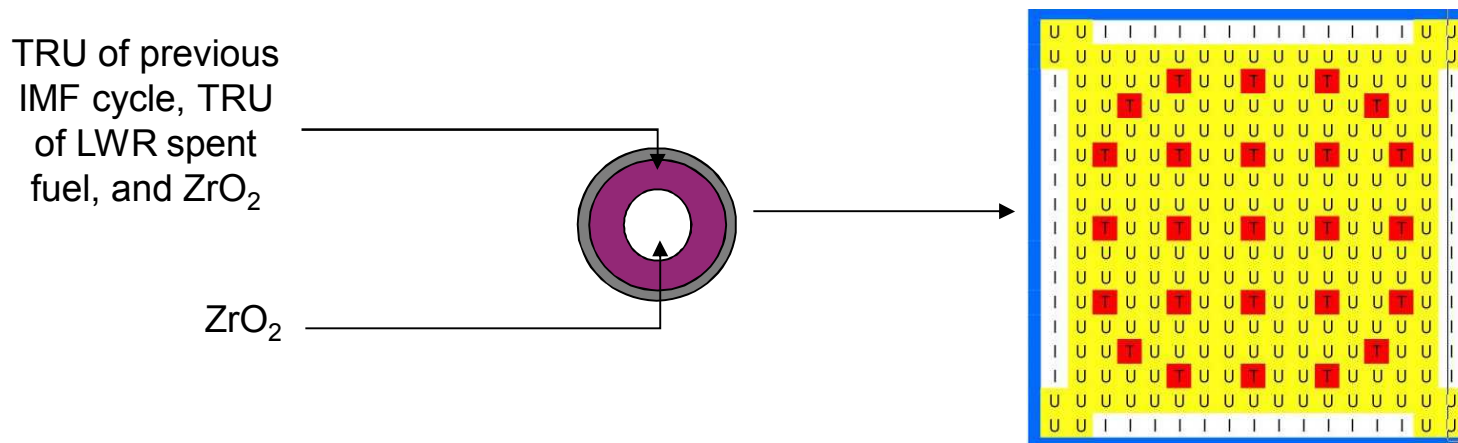




# IMF Fuel Assemblies (Cont.)



- Heterogeneous Fuel Assemblies
  - IMF fuel pins are located at peripheral 52 positions of 17x17 typical PWR assembly
  - IMF fuel pins are made by blending TRUs of previous IMF cycle and LWR spent fuel
  - Cycle length is controlled by Uranium enrichment, blending ratio, and  $\text{TRUO}_2$  mass fraction under power peaking of 1.2



**Heterogeneous IMF Assembly**



# LWR-UO<sub>2</sub> Cycle Interpolations



- Uranium Enrichment
  - 33 GWd/t (2.992% U-235), 60 (5.049%), 100 (8.500%)

	Actual Uranium Enrichment	Predicted Uranium Enrichment (linear)	Errors for Linear Interpolation (%)	Predicted Uranium Enrichment (quadratic)	Errors for Quadratic Interpolation (%)
<b>45 GWd/t</b>	3.872	3.918	1.18	3.879	0.18
<b>75 GWd/t</b>	6.294	6.394	1.56	6.287	0.12

- TRU vector after aging for varied cooling times
  - Three burnups (33, 60, and 100 GWd/t)
  - Five cooling times (5, 14, 28, 40, and 50 years)



# LWR-UO<sub>2</sub> Cycle Interpolations (Cont.)



- Four test cases
  - 45 GWd/t burnup with 10 and 35 year cooling
  - 75 GWd/t burnup with 22 and 45 year cooling
  - Quadratic interpolation for burnup
  - Linear interpolation for cooling time

- Results

Error (%)	Number (out of 56)	Percentage
< 1	38	67.8
1 - 5	16	28.6
> 5	2	3.6

- Larger errors observed with smaller cooling times



# Heterogeneous Results



- TRUO<sub>2</sub> Mass Fractions

45 GWd/t burnup, 10 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	% Error
Cycle 0	20.49	21.47	4.58
Cycle 1	26.18	27.42	4.53
Cycle 2	29.93	31.50	4.99

75 GWd/t burnup, 45 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	% Error
Cycle 0	36.97	37.05	0.22
Cycle 1	38.57	38.00	1.50
Cycle 2	40.76	40.67	0.22

- Uranium Enrichments

45 GWd/t burnup, 10 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	% Error
Cycle 0	4.61	4.60	0.65
Cycle 1	4.75	4.75	0.14
Cycle 2	4.86	4.86	0.12

75 GWd/t burnup, 45 year initial cooling, and 7 year IMF Cooling			
	Predicted	Actual	% Error
Cycle 0	4.93	4.94	0.15
Cycle 1	5.03	5.02	0.28
Cycle 2	5.10	5.11	0.17

- 72.1% and 97.1% of 68 predicted values had errors less than 5% for 45 and 75 GWd/t burnup cases, respectively



# Background



- AFCI (Advanced Fuel Cycle Initiative)
  - Reduce volume and toxicity of nuclear waste
  - Reduce proliferation threat posed by plutonium
  - Reclaim energy contained in spent fuel
- Transuranic Recycling in Commercial Light Water Reactors (LWRs)
  - Manages the inventory of transuranics (TRU) in commercial spent nuclear fuel (CSNF) and impedes further accumulation
  - Helps increase the loading capacity of high-level wastes in the Yucca Mountain repository
  - Capable of utilizing a large capacity of existing nuclear reactor facilities



# Methodology - Codes



- WIMS9
  - 172 group neutron library based on JEF2.2
  - Heavy nuclides and about 100 fission products are explicitly traced in irradiation
  - Calculates physics parameters (eigenvalues and power peaking)
  - Creates one group cross sections for ORIGEN2.1 calculations
- ORIGEN2.1
  - Performs depletion calculation to generate spent fuel composition using one group cross sections generated by WIMS9
  - Simulates cooling, reprocessing, and aging processes





# Trends Observed



- Values increase with burnup, cooling time, and cycle number
  - $\text{TRUO}_2$  mass fraction
  - Charge mass
  - Discharge mass
  - Uranium Enrichment
- Blending ratio decreases between cycle 1 and cycle 2 in high burnup and initial cooling time cases
- Exponential interpolation between cooling times can improve predictions for  $^{241}\text{Pu}$  mass fraction
  - No closed form solution for increasing quantities