

Establishing a Methodology for Performing a Multiphysics Run-In Analysis of the GPBR200

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Ryan Hunter Stewart, Paolo Balestra, Jack M Cavaluzzi





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Ryan Hunter Stewart, Paolo Balestra, Jack M Cavaluzzi

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Idaho National Laboratory Idaho Falls, Idaho 83415

http://www.inl.gov

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Ryan Stewart, Jack Cavaluzzi, Paolo Balestra

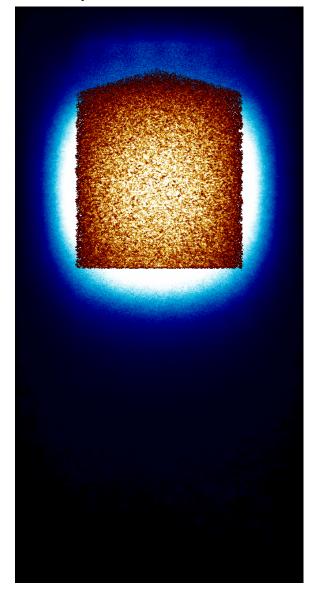
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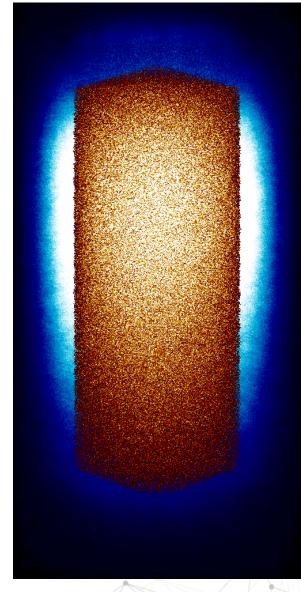
Introduction

- The running-in phase of a pebble-bed reactor (PBR) is a complex time-dependent problem
 - Involves the use of multiple fuel types, graphite pebbles and a ramp-up of power
- Modeling this problem using high-fidelity simulation tools allows us to examine multiple physical phenomena that is important to PBR operations
 - quantities of interest: discharge burnup, time to full power, pebble power peaking, etc.
- Understanding the temperature distribution will provide insight into key physics missing in isothermal core analysis

Step 1 – Critical Core

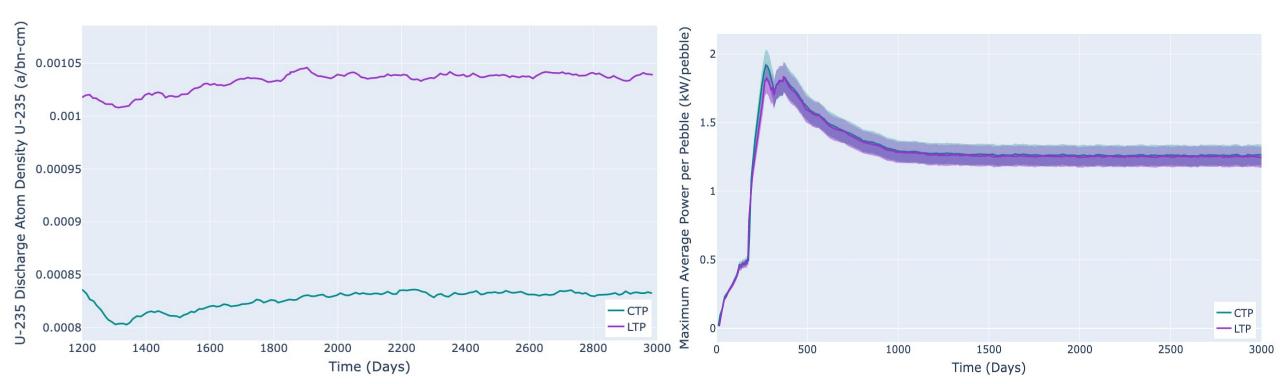


Step N – Equilibrium Core



Justification of Research

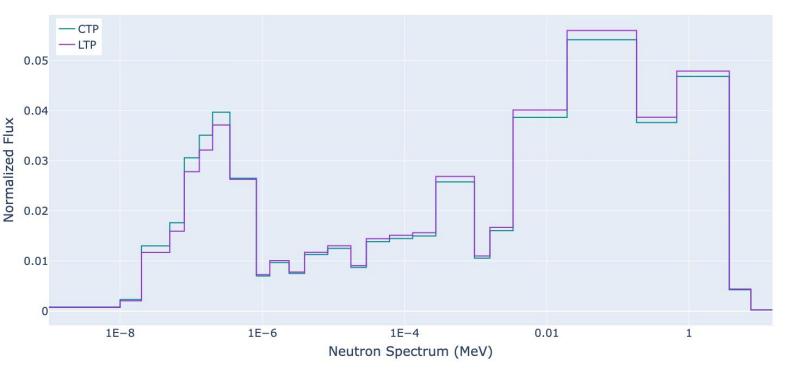
- Previous research showed that temperature can influence core neutronics
 - Isothermal vs linear temperature distribution

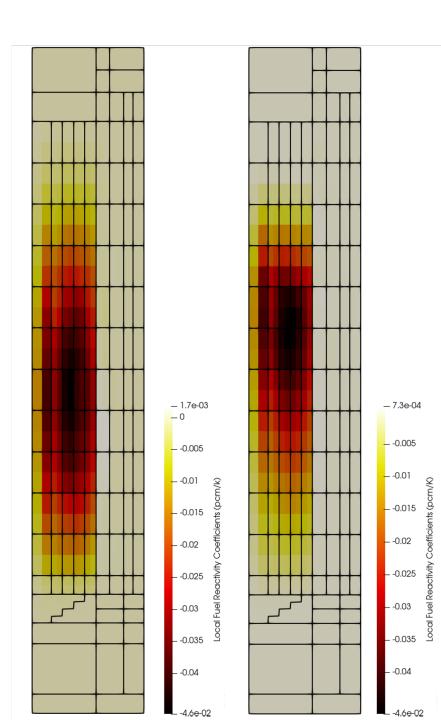


Justification of Research

- Previous research showed that temperature can influence core neutronics
 - Isothermal vs linear temperature distribution
- Temperature reactivity coefficients

Upper Region





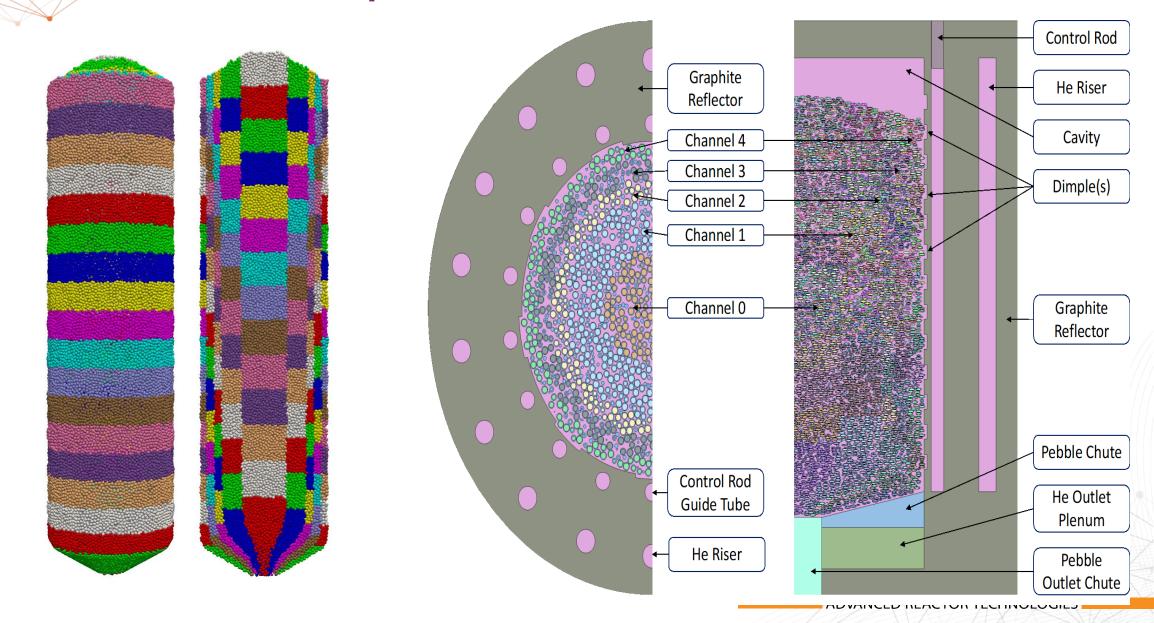
Project Goal

- Goal: Develop a methodology for incorporating multi-physics into our high-fidelity run-in simulation.
- Approach: Utilize a Python module wrapped around Serpent and Griffin/Pronghorn to simulate pebble movement through the core

Algorithm Outline

- 1. Perform burn-up step (Serpent)
- 2. Determine temperature distribution (Griffin/Pronghorn)
- 3. Shift pebbles down (Python)
- 4. Recycle/discharge pebbles (Python)
- 5. Refuel core (Python)
- 6. Update power, temperature, pebble type, etc. (Python)

GPBR200 Model: Serpent

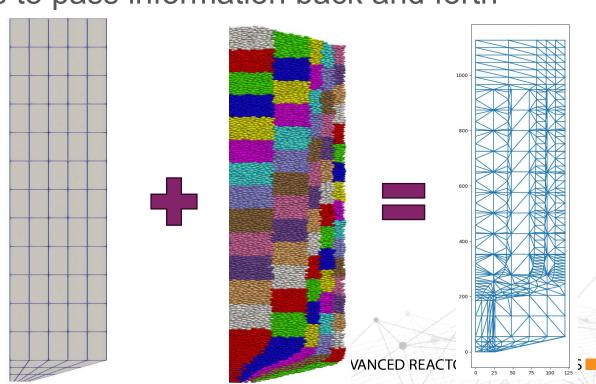


GPBR200 Model: Griffin/Pronghorn

- Captured as much of the geometry into a 2D RZ mesh
 - Varying graphite density to capture lower plenum
 - Control rod and risers implemented
- Core region
 - No streamlines (60 blocks within the core)
 - Conus region is modeled separately
- Note: Griffin blocks do not align with the Serpent axial volumes
 - Requires a mapping between the two

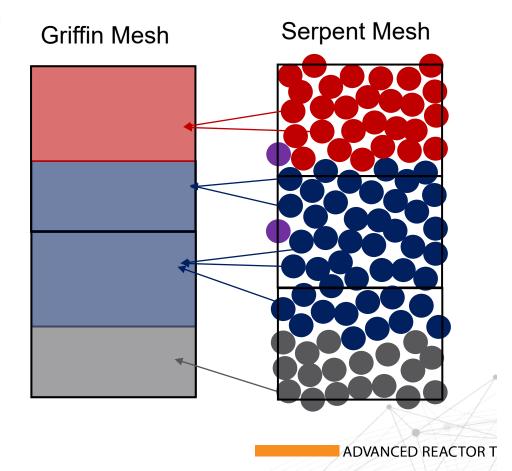
Data Transfer: Initial Attempt

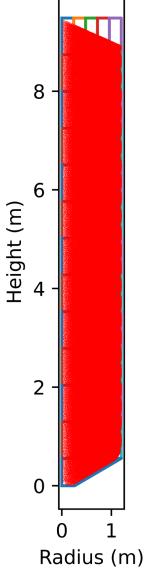
- Create a unifying mesh utilizing triangularization
 - Triangles can't cross mesh in either individual mesh (Griffin or Serpent)
- Map data in each triangle to correspond to a specific Griffin/Serpent mesh
- Volume average number densities to pass information back and forth
- Triangulation mapping became difficult
 - Mainly when vertices and segments approached each other
- Method could be possible for simpler designs



Data Transfer: Finalized Approach

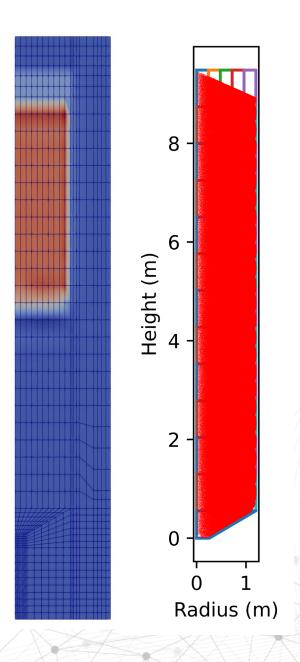
- Serpent provides explicit pebble positions, these can be easily mapped to a Griffin block and vice-versa
 - Material creation in Griffin is based on volume averaging based on the number of pebbles of each pass, fuel type, and each axial volume
- The reverse is performed to pass information in a block to a pebble





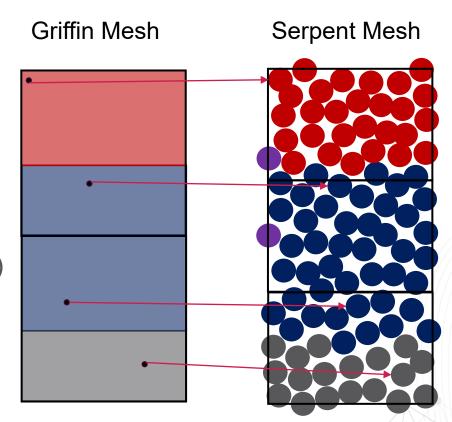
Results: Static Data Transfer

- Microscopic cross-sections have been generated from DRAGON
- Pebble isotopics can be passed from Serpent to Griffin
 - Mapping function performs the previous data transfer
- Griffin-only calculations have been performed
 - Provides confidence in the methodology
- Currently developing SPH factors
 - Help correct our cross-sections an maintain reaction rates



Results: Where are we going?

- Coupled Griffin/Pronghorn
 - Isotopic compositions
 - Generate SPH factors to conserve reaction rates
- Stand-alone Pronghorn
 - Power density
- Regardless of approach, we generate a temperature field
 - Each element will have an associated temperature
 - Pass average temperature to Griffin (50 K increments)



Conclusions & Wrap Up

- We have provided a framework for coupling a high-fidelity Monte Carlo code with a deterministic solver
 - Provides an opportunity to incorporate thermal fluid feedback into our Monte Carlo simulation
- Initial Serpent to Griffin coupling has been performed
 - Working on determining if SPH factors can be used
- Final stage will be to compare isothermal/linear temperature profile simulation with Multiphysics run-in simulation

Questions?