

FULL CORE POWER AND ISOTOPIC OSCILLATIONS WITH VARIOUS DEPLETION SCHEMES

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OUTLINE



- Governing depletion equations
- Summary of depletion coupling schemes
- Full core analysis
- Single assembly convergence study
- Conclusions
- Discussion

- Bateman equations (Bell and Glasstone 1970)

$$\frac{dN(r, t)}{dt} = \underline{\underline{M}}(R(r, t))N(r, t)$$

- $R(r, t)$ contains the reaction rates
- $\underline{\underline{M}}$ is a matrix describing transmutation of isotopes
- Can solve for $N(r, t)$ by taking exponent of matrix
- Transmutation trajectory analysis (TTA) (Cetnar 2006)
- Chebyshev rational approximation method (CRAM) (Pusa and Leppänen 2010; Lago and Rahnema 2017; Pusa 2011)
- Reaction rates and nuclide distributions are strongly coupled
- Temperature dependence through Doppler broadening

Predictor - CE

1. Beginning of step (BOS) flux depletes to obtain N_{i+1}

Predictor-corrector – CE/LI

1. BOS flux depletes to obtain N_{i+1}^P
2. Predictor end-of-step (EOS) number density is used to obtain a corrector BOS flux ϕ_i^C
3. Average of fluxes is used to deplete across interval and obtain EOS isotopics

$$\overline{\phi_i} = \frac{1}{2} (\phi_i^P + \phi_i^C)$$
$$\overline{\phi_i} \rightarrow N_{i+1}$$

Higher order methods utilize previous EOS flux (Isotalo and Aarnio 2011a)

- Extrapolate using ϕ_{i-1} and ϕ_i to obtain flux throughout depletion interval ϕ_i^P
- Deplete with ϕ_i^P and obtain N_{i+1}^P
- Set the next BOS number density using interpolation

$$N_{i+1} = w_{i-1}N_{i-1} + w_iN_i + w_{i+1}N_{i+1}^P$$

- Linear extrapolation/Quadratic interpolation – LE/QI

Substep methodology (Isotalo and Aarnio 2011b)

- Assume that reaction rates are piecewise constant
- Many depletion sub steps inside depletion interval
- No additional transport solutions

Stochastic Implicit Euler (SIE) (Dufek, Kotlyar, and Shwageraus 2013)

- Multiple transport solutions per depletion interval
 - Total histories equally distributed
- Iteration-wise EOS flux obtained using iteration EOS number density

$$N_i^{(n)} \rightarrow \phi_i^{(n)}$$

- Updated end of step flux is average of iteration-wise fluxes

$$\overline{\phi_i^{(n)}} = \frac{1}{n} \sum_{j=1}^n \phi_i^{(j)}$$

- Averaged flux used to deplete and obtain isotopics for next iteration

$$\overline{\phi_i^{(n)}} \rightarrow N_i^{(n+1)}$$

- After N iterations

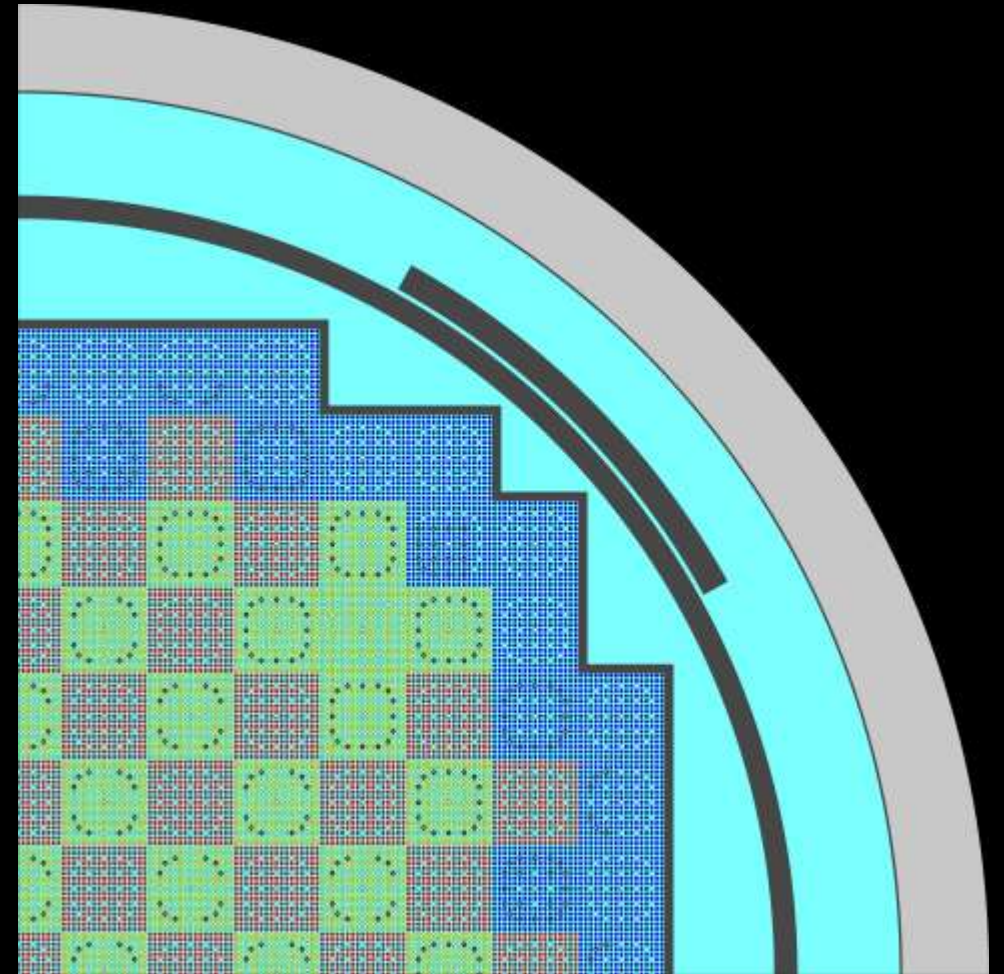
$$\phi_{i+1}^{(0)} = \overline{\phi_i^{(N)}}$$

Explicit Euler variations

- Use flux ϕ_i to obtain the number density N_{i+1} separated by some Δt
- Assume that all other parameters are constant
 - Reaction rates
 - Reactor state
- Increasing Δt means fewer transport calculations
 - Accurate in the limit of small Δt
- Isotopic oscillations have been observed (Dufek et al. 2013; Kotlyar and Shwageraus 2013)
 - Multi-region problems

FULL CORE ANALYSIS

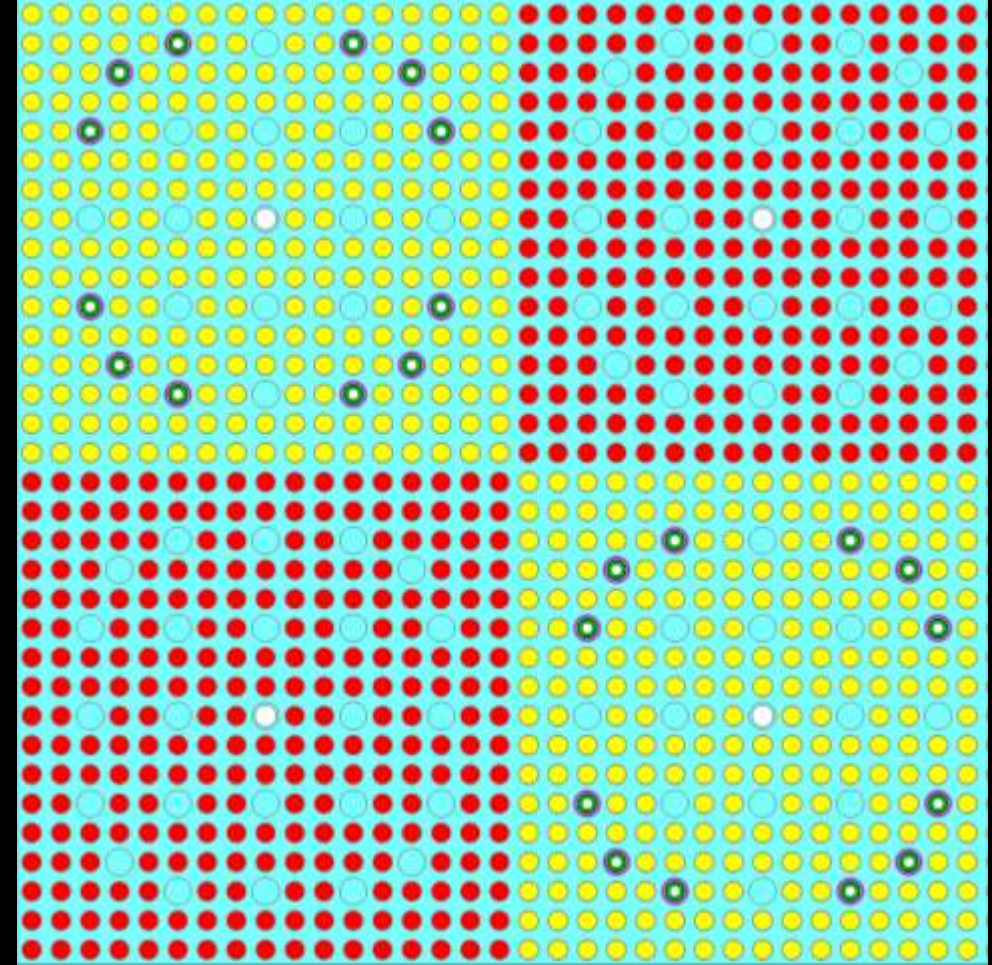
- BEAVRS reactor (Horelik et al. 2013)
- 150000 particles/cycle
- 100/500 inactive/active cycles
- ENDF/B-VII.0



Quarter of the BEAVRS v2 SERPENT Model

FULL CORE - MODELING

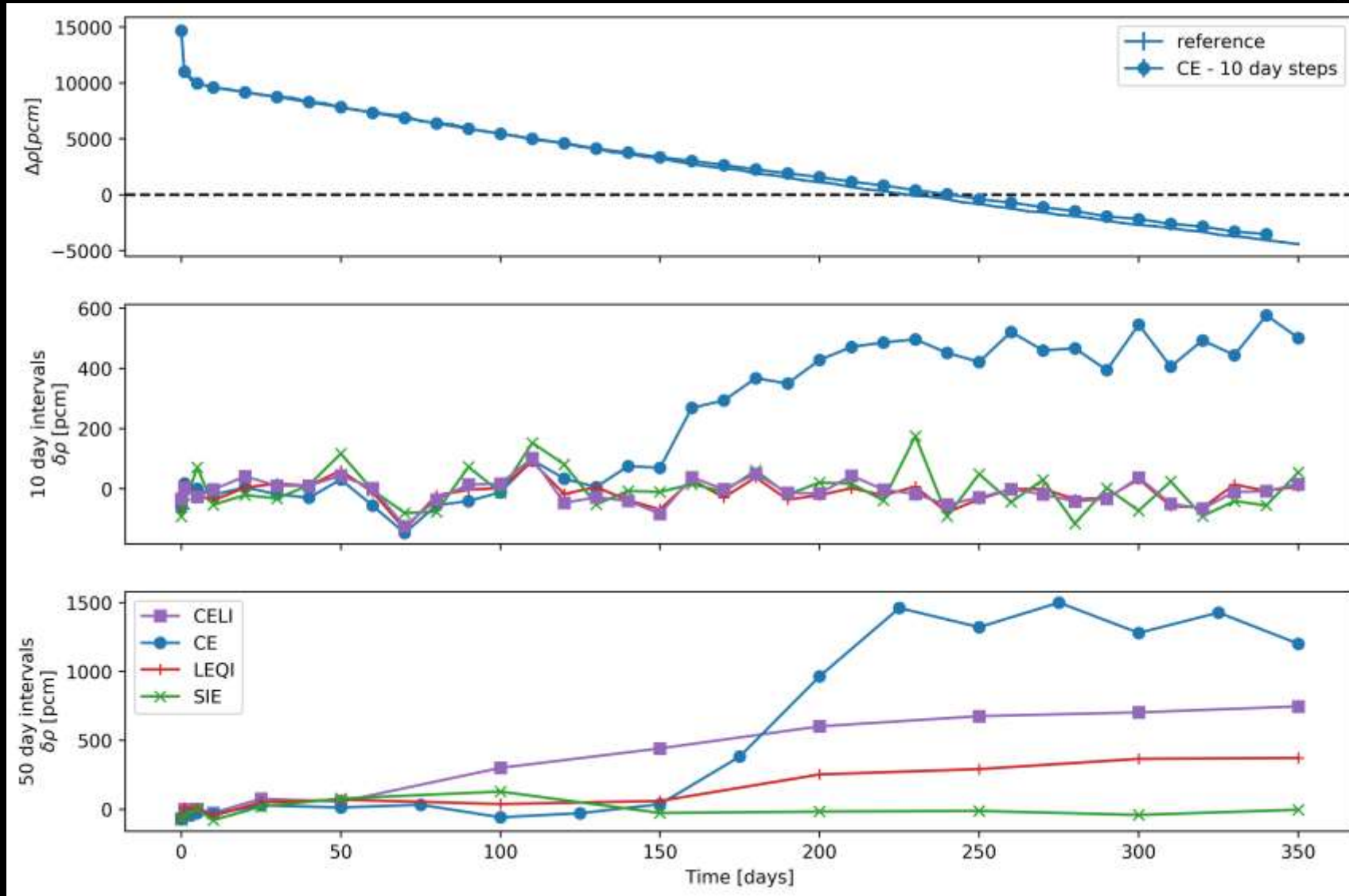
1. $1/8^{\text{th}}$ core explicitly modeled
2. 12 axial layers of unique materials
3. Unique burnable materials per assembly
4. No spacer grids nor soluble boron



3.1 wt% (red) and 2.4 wt% (yellow) enriched assemblies

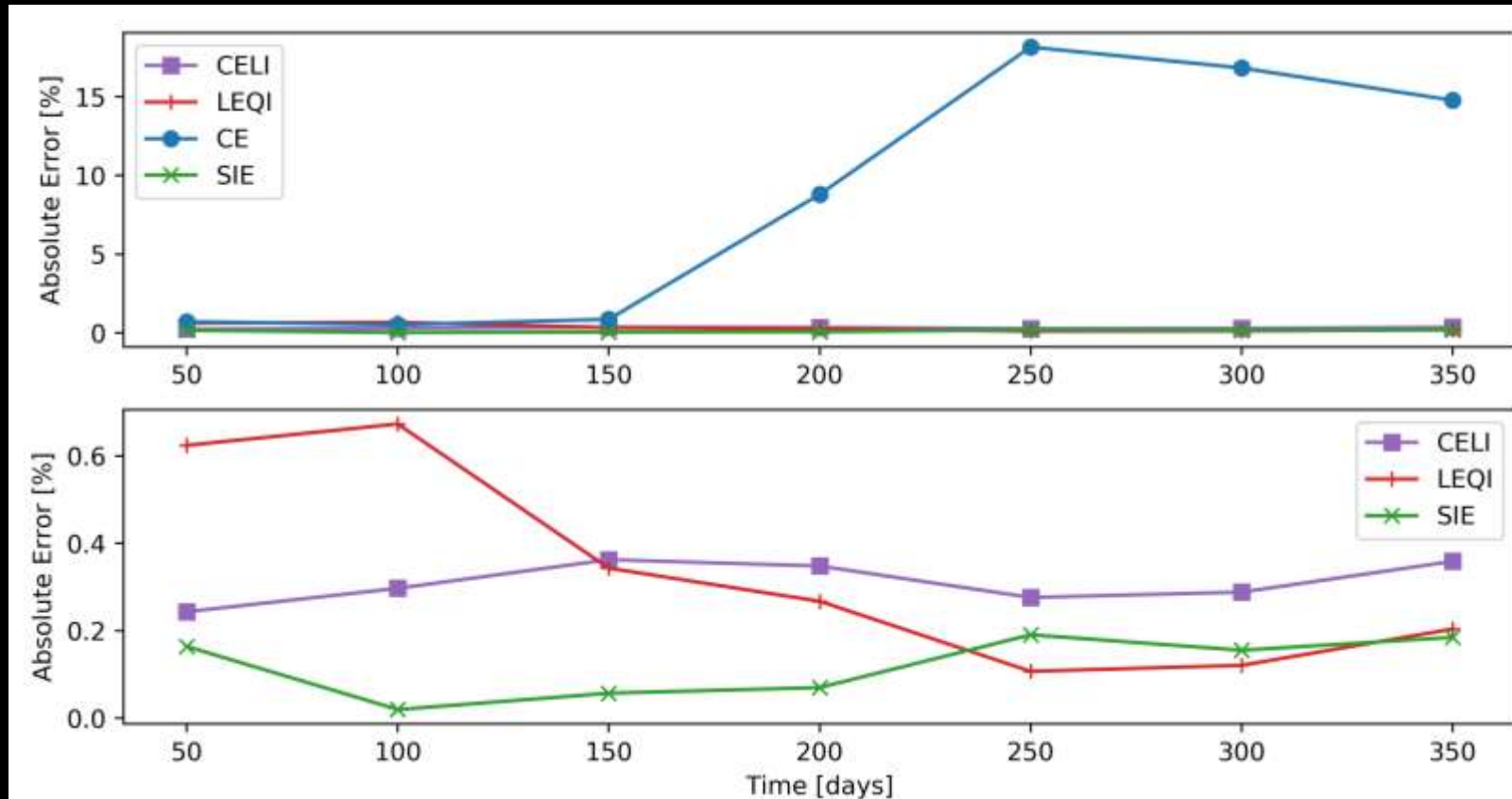
- 350 day depletion schedule
 - 1, 4, and 5 day initial depletion intervals
- Three separate depletion schemes
 - 5 day time steps
 - 10 day time steps
 - 50 day time steps
- Reference: SIE with 5 day time steps
- Tested nearly all schemes offered by SERPENT (Leppänen et al. 2015)
 - CE, CE/LI, LE/QI, SIE
- SIE utilized 10 iteration per interval
 - $1/10^{\text{th}}$ as many particles per cycle
- CE used twice as many depletion events
 - Equal number of transport calculations

FULL CORE – KEFF



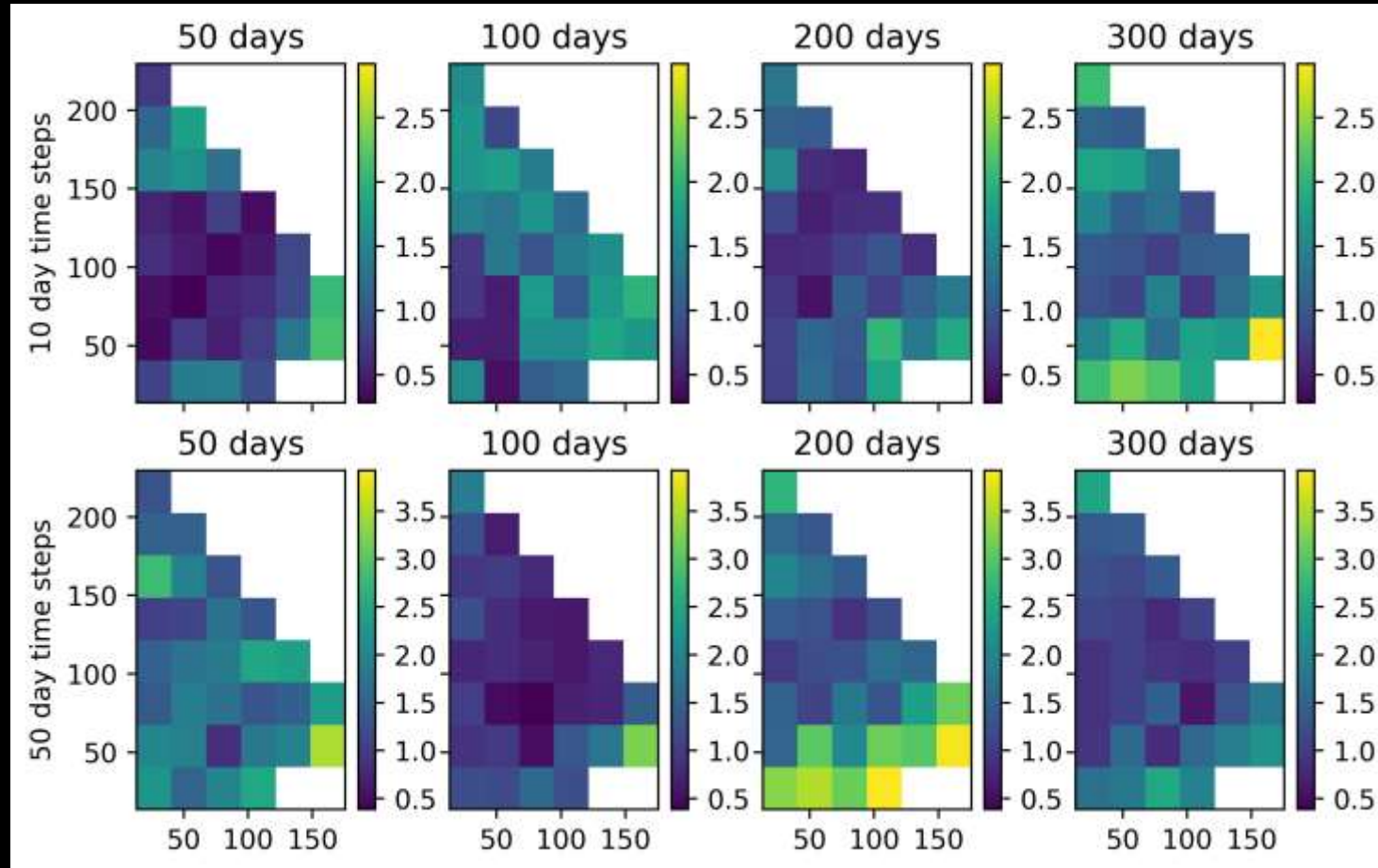
k_{eff} and $\delta\rho$ for full core analysis

FULL CORE – CUMULATIVE XENON



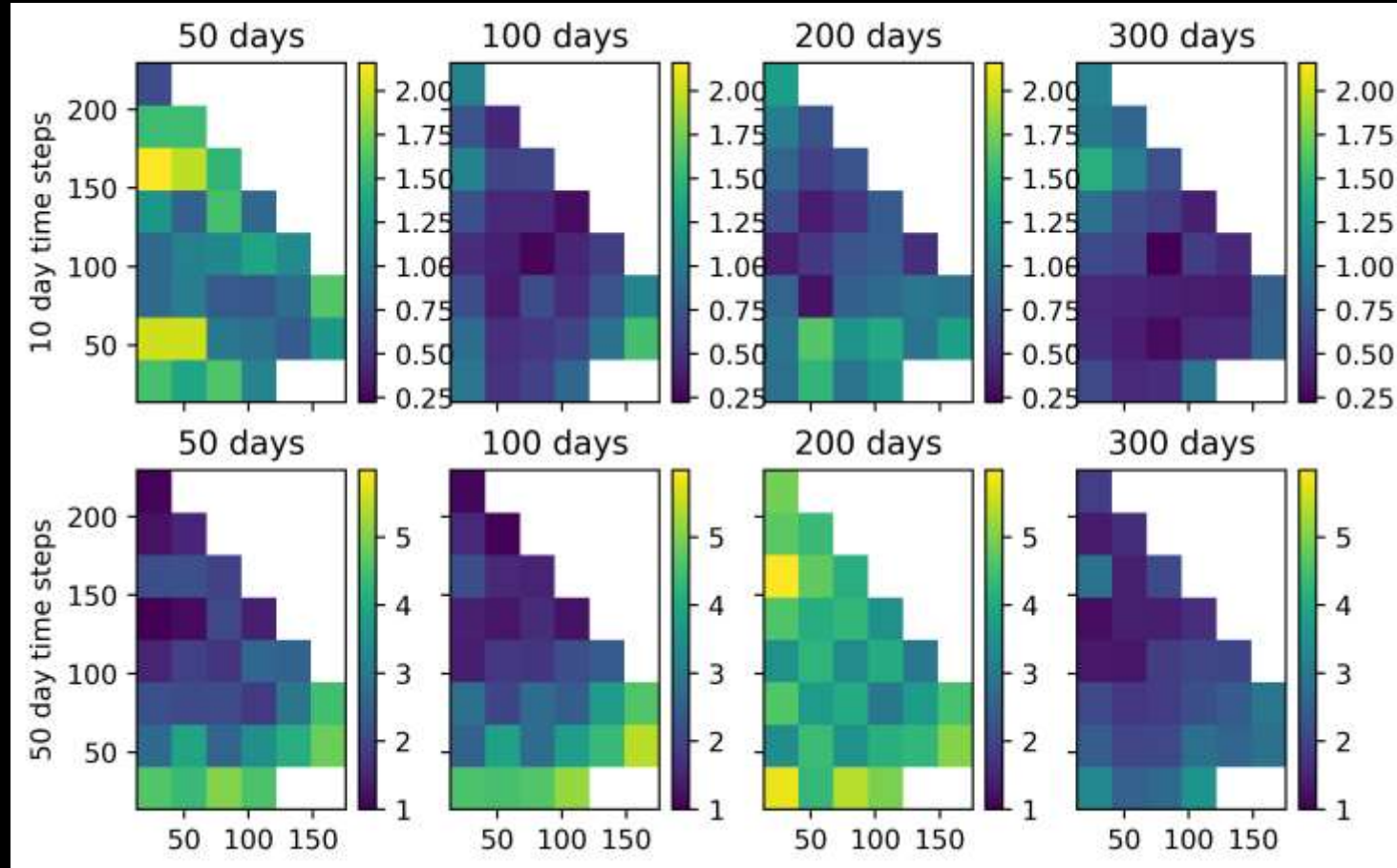
Absolute error in full core cumulative
Xe-135 density 50 day steps

FULL CORE – ASSEMBLY XENON ERRORS



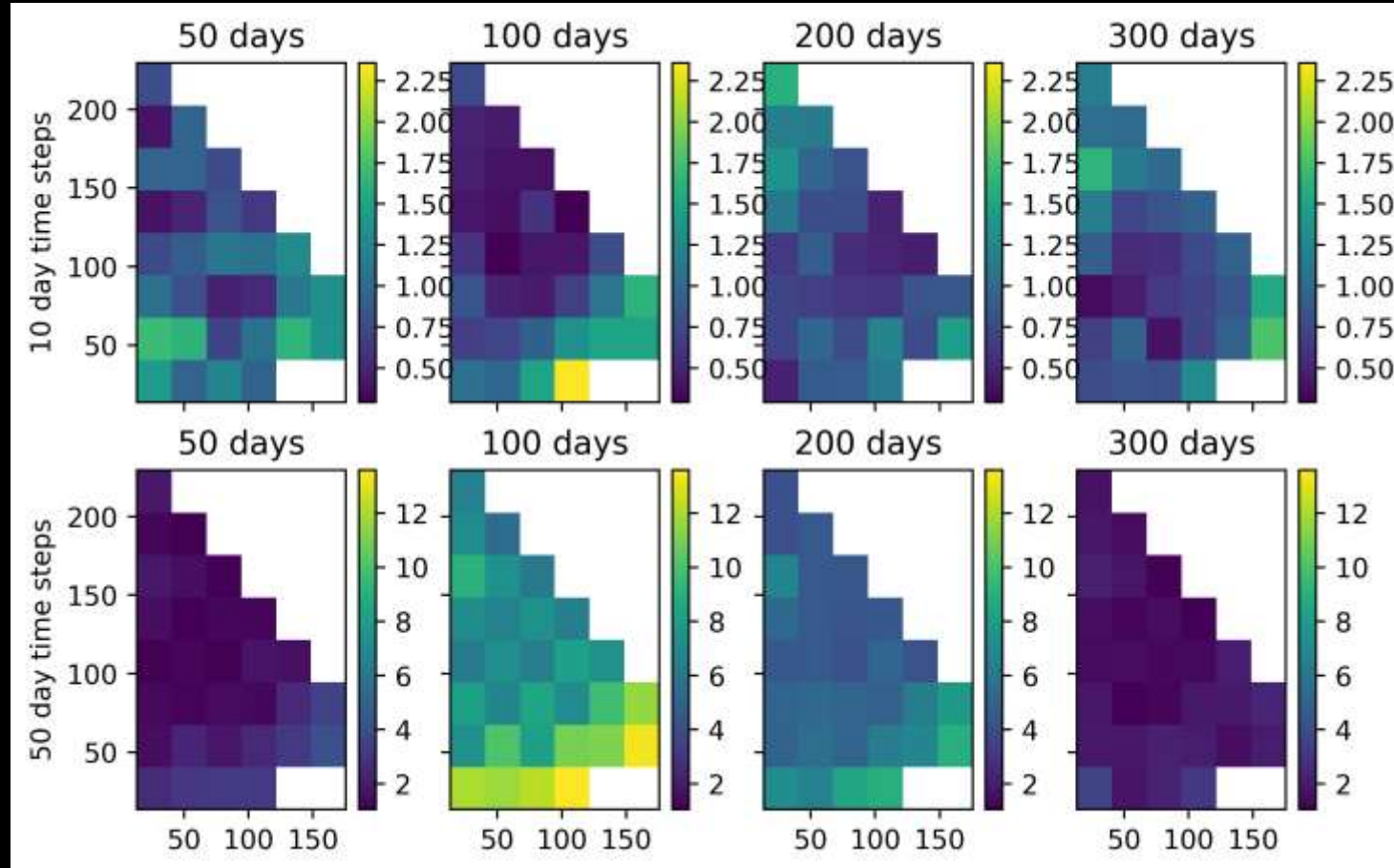
Maximum nodal xenon density error
using SIE method

FULL CORE – ASSEMBLY XENON ERRORS



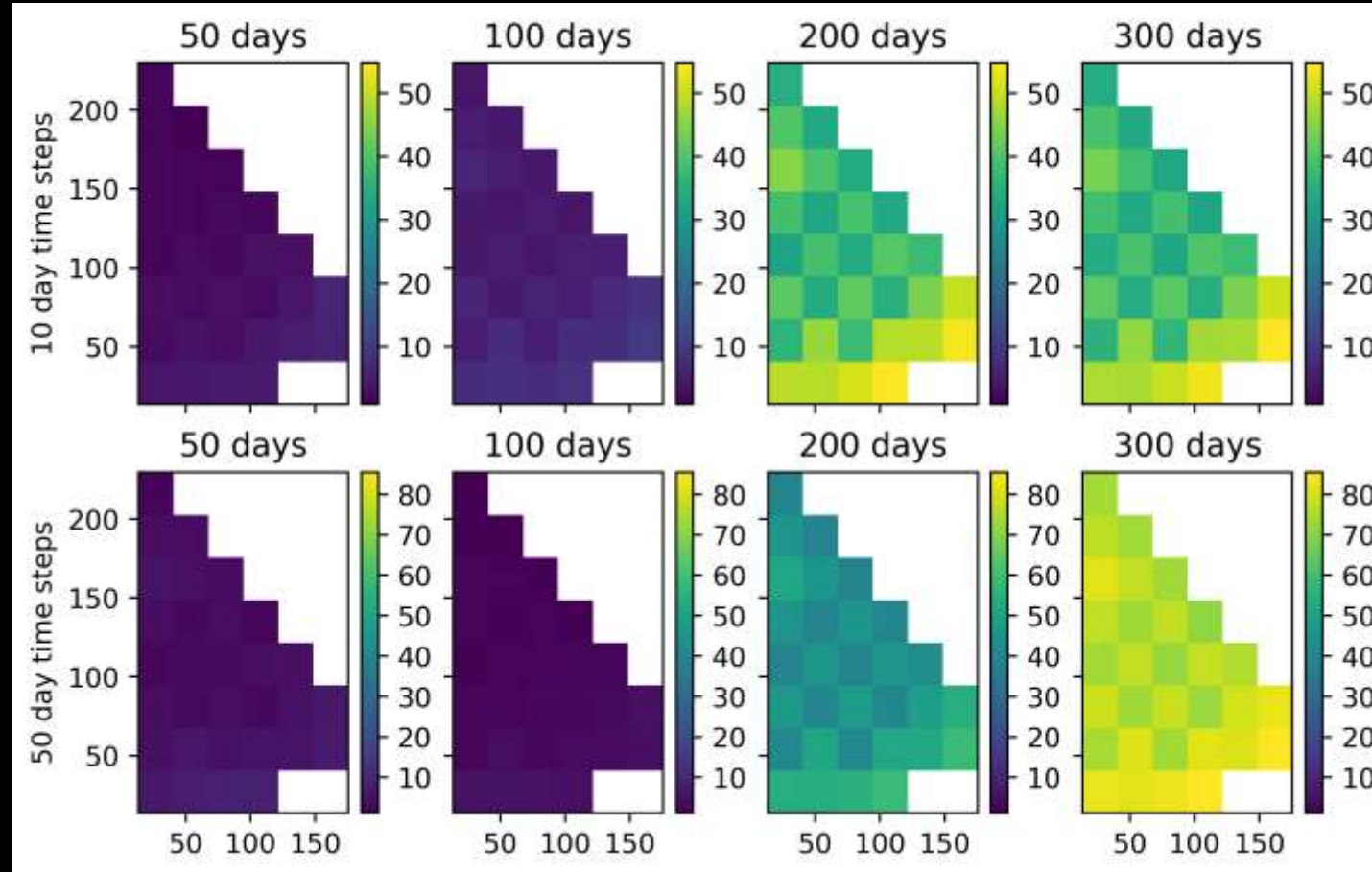
Maximum nodal xenon density error
using LE/QI method

FULL CORE – ASSEMBLY XENON ERRORS



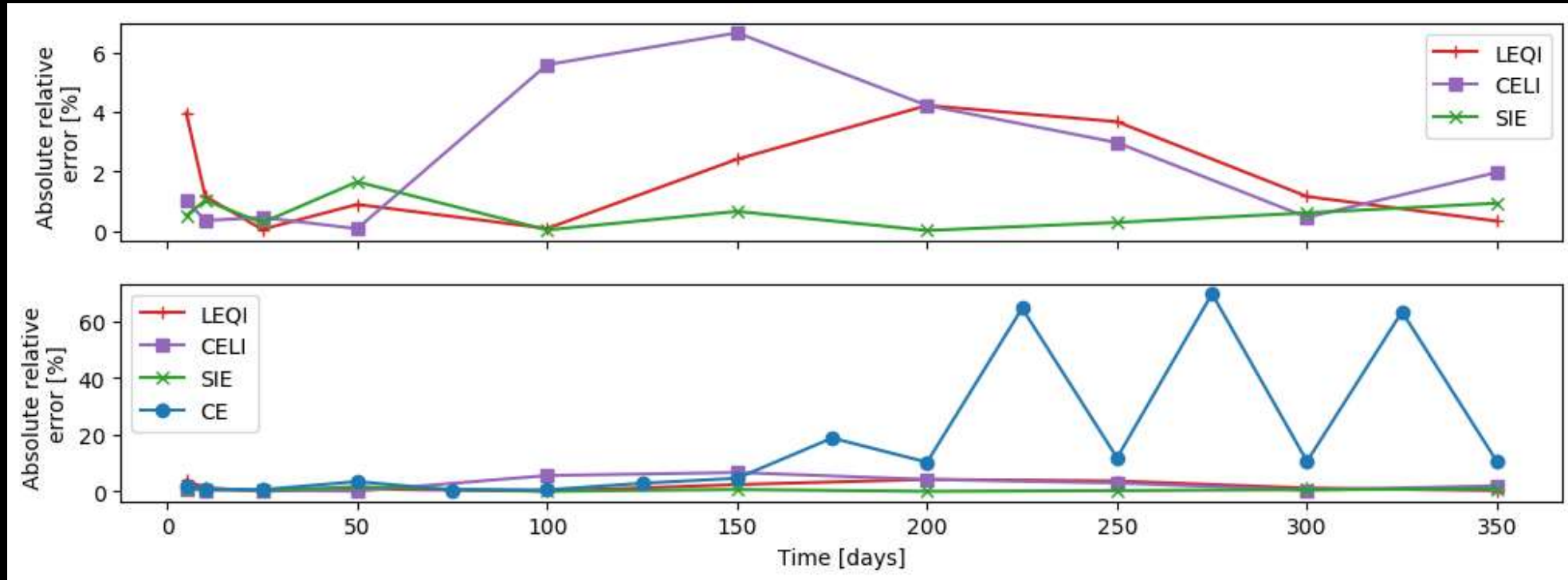
Maximum nodal xenon density error
using CE/LI method

FULL CORE – ASSEMBLY XENON ERRORS



Maximum nodal xenon density error
using CE method

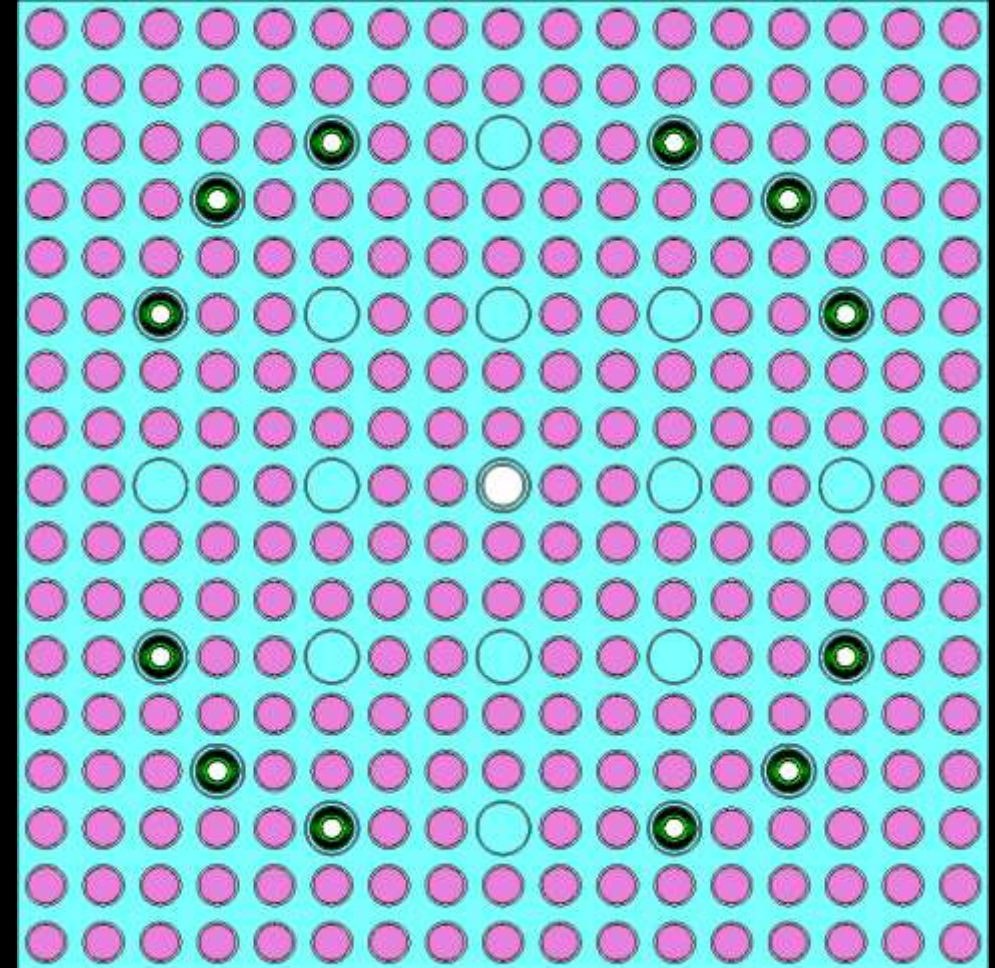
FULL CORE – NODAL XENON ERRORS



Xenon density and error for a single
node – 50 day schedule

CONVERGENCE STUDY

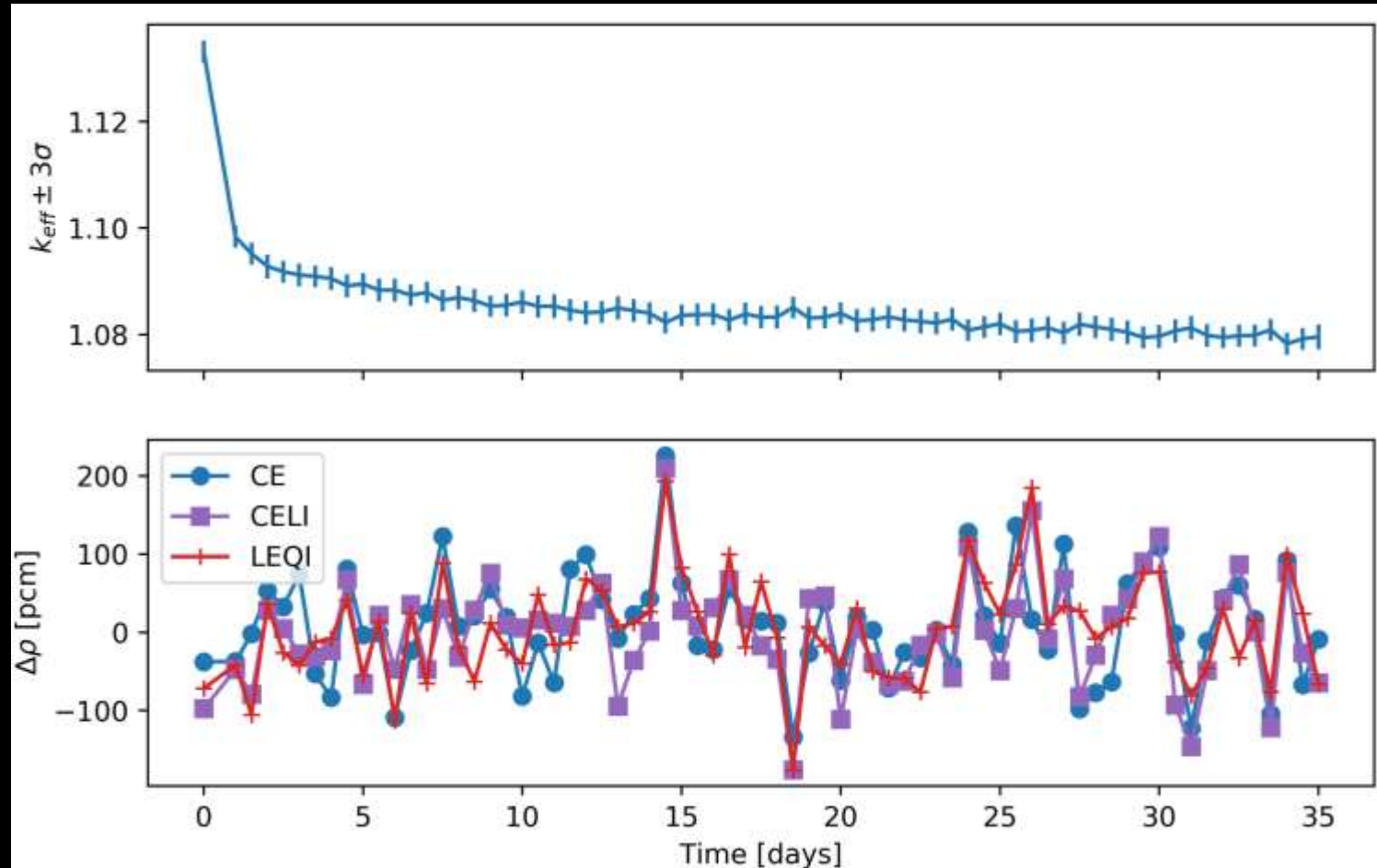
- Further investigate the effect of decreasing time steps
- Single assembly, 12 burnable absorbers from reactor
- 16 axial layers
- 7500 particles/cycle
- 100/500 inactive/active cycles



12BA assembly for convergence study

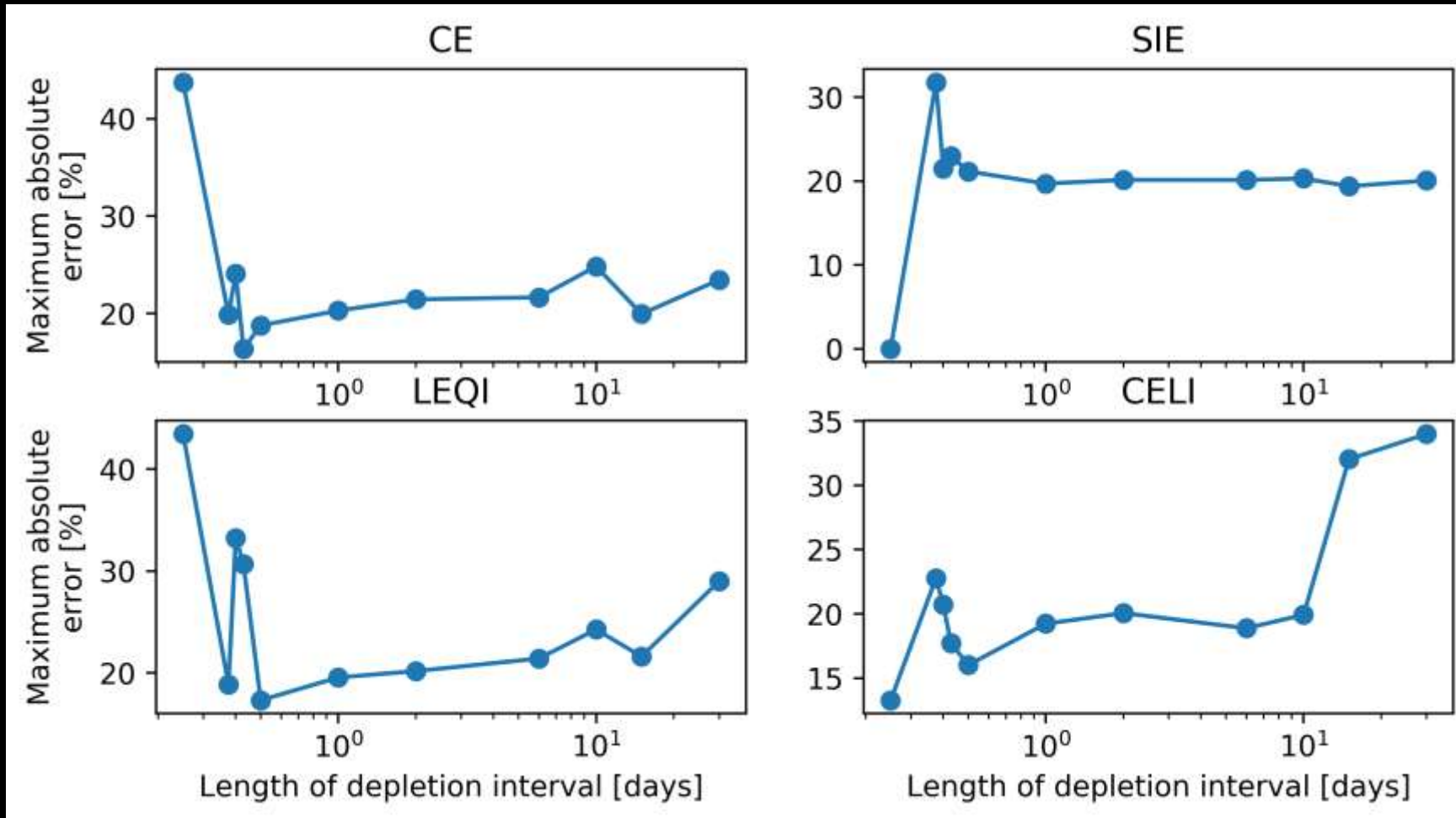
- 35 day schedule
 - 0.5 days * 10 for initial FP build up
 - Remaining 30 days divided into integer number of equal-width intervals
 - 30 days per step down to 0.25 days per step
- 10 substeps per depletion interval – CE/LI and LE/QI
 - 10 iterations per depletion interval – SIE
 - Reference – 0.25 days per depletion interval

CONVERGENCE STUDY

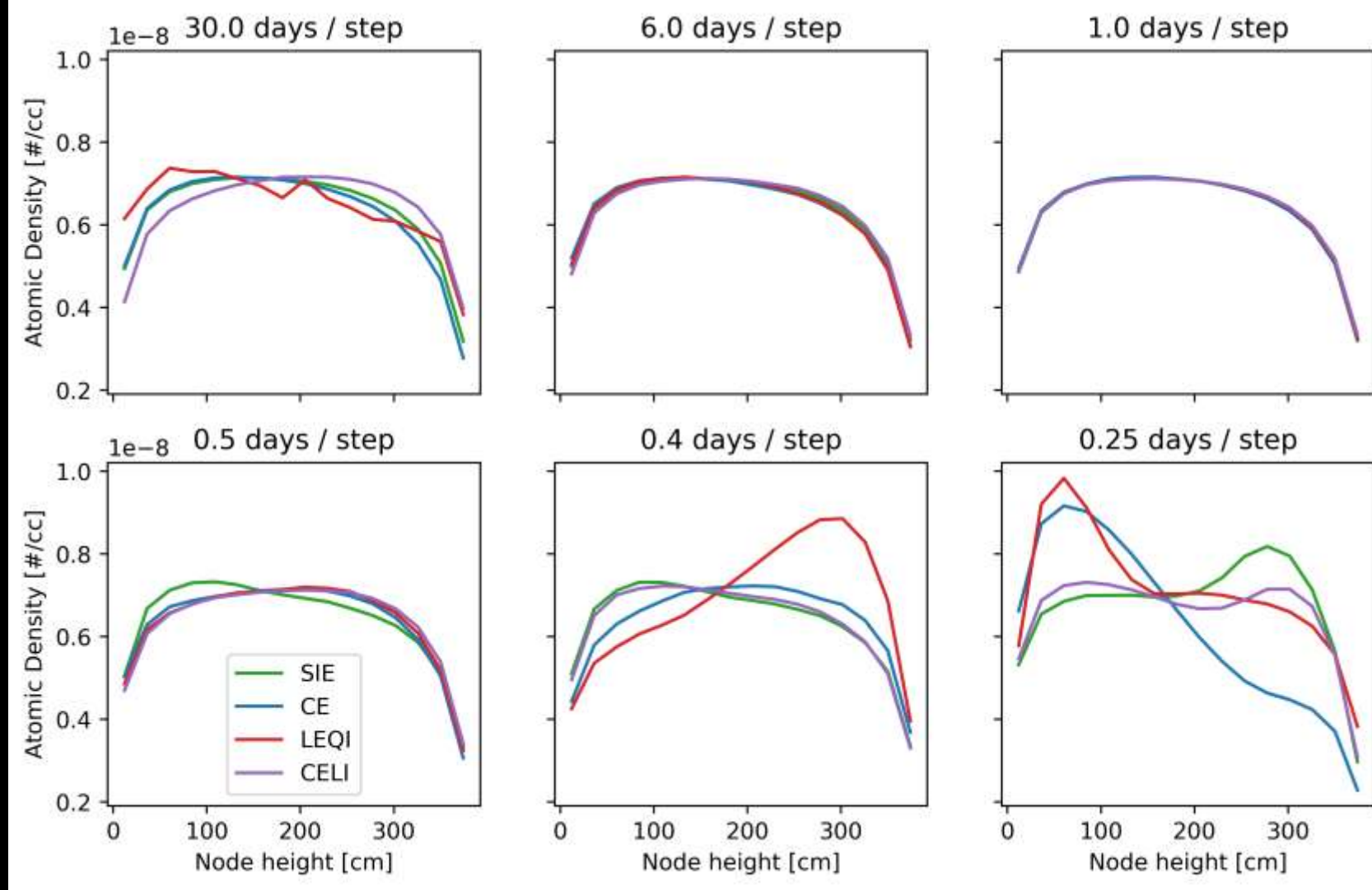


k_{eff} and $\delta\rho$ for methods with 0.5 day steps

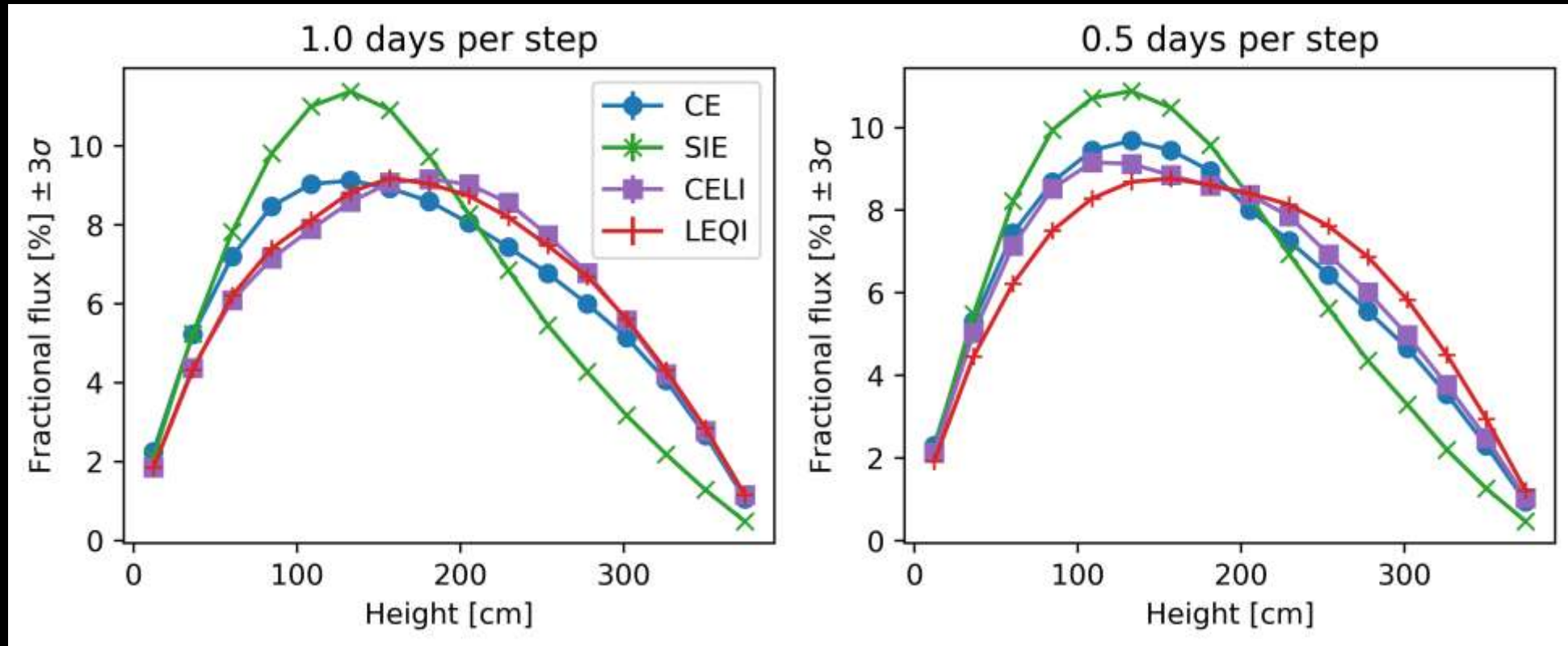
ASSEMBLY– XENON ERROR CONVERGENCE



Maximum nodal xenon density error at EOL vs.
length of depletion interval



Xenon distribution at end of cycle for various depletion interval lengths



Nodal flux share at EOL for 1 and 0.5 day steps for all methods

Full Core Analysis

- ≥ 500 pcm error in k_{eff}
- 20% error in local xenon density
 - Similar results were obtained for other fission products
- Methods appeared to converge towards the reference solution at EOL

Assembly Analysis

- Some indication of error convergence for reducing time steps
- Divergence for time steps below one day
- 10 % deviation in nodal flux

- As with this work, reference cases are often those with the smallest time step
- Trend towards transient and life-cycle analysis with Monte Carlo methods
- Accuracy of depletion method will drive accuracy of transport solution
- Further work on improved coupling schemes
- Perturbation theory-based methods
- More points for extrapolation/interpolation
- Benchmarking

REFERENCES



Bell, George I., and Samuel Glasstone. 1970. *Nuclear Reactor Theory* (Van Nostrand Reinhold Company: New York).

Dufek, Jan, Dan Kotlyar, Eugene Shwageraus, and Jaakko Leppänen. 2013. 'Numerical stability of the predictor-corrector method in Monte Carlo burnup calculations of critical reactors', *Annals of Nuclear Energy*, 56: 34-38.

Horelik, Nicholas, Bryan Herman, Benoit Forget, and Kord Smith. 2013. "Benchmark for Evaluation and Evaluation of Reactor Simulations (BEAVRS), v1.0.1." In *International Conference on Mathematics and Computational Methods Applied to Nuclear Science and Engineering*. Sun Valley, Idaho.

Isotalo, A. E., and P. A. Aarnio. 2011. 'Higher order methods for burnup calculations with Bateman solutions', *Annals of Nuclear Energy*, 38: 1987-95.

Isotalo, A. E., and P. A. Aarnio. 2011. 'Substep methods for burnup calculations with Bateman equations', *Annals of Nuclear Energy*, 38: 2509-14.

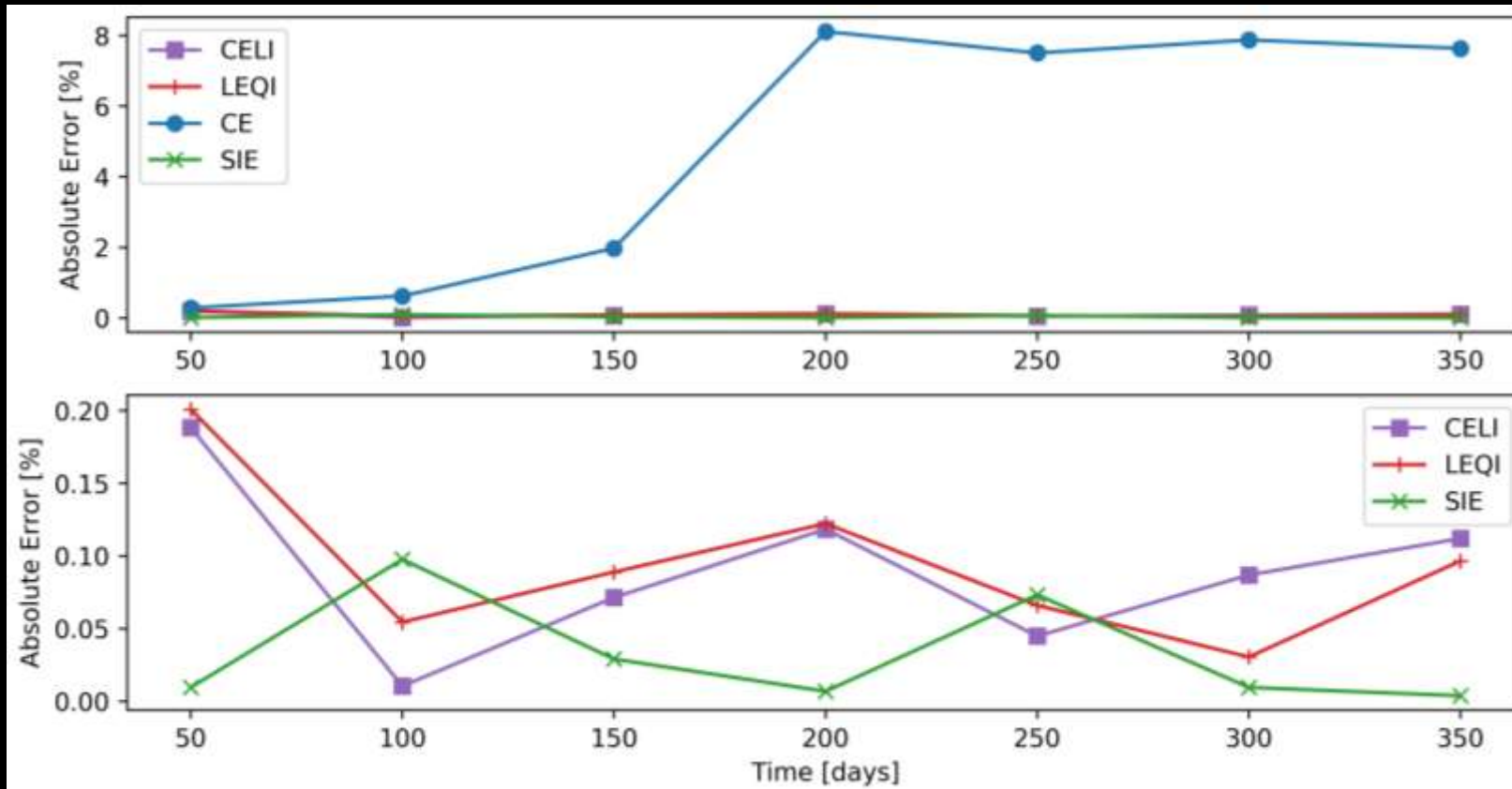
Lago, Daniel, and Farzad Rahnema. 2017. 'Development of an Application Programming Interface for Depletion Analysis (APIDA)', *Annals of Nuclear Energy*, 103: 163-72.

Leppänen, Jaakko, Maria Pusa, Tuomas Viitanen, Ville Valtavirta, and Toni Kaltiaisenaho. 2015. "The Serpent Monte Carlo code: Status, development, and applications in 2013." In *Annals of Nuclear Energy*.

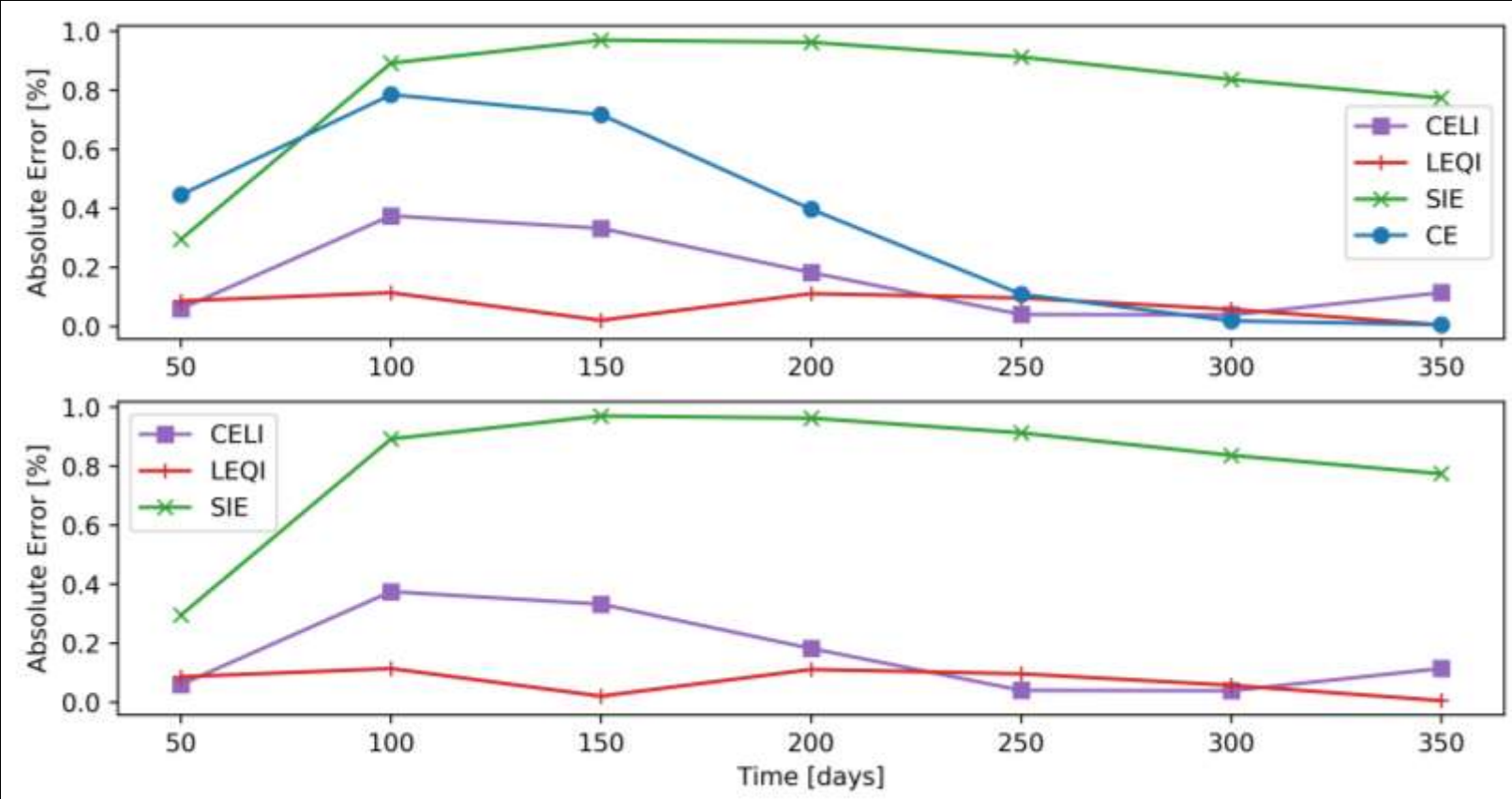
Pusa, Maria. 2011. 'Rational Approximations to the Matrix Exponential in Burnup Calculations', *Nuclear Science and Engineering*, 196: 155-67.

Pusa, Maria, and Jaakko Leppänen. 2010. 'Computing the Matrix Exponential in Burnup Calculations', *Nuclear Science and Engineering*, 164: 140-50.

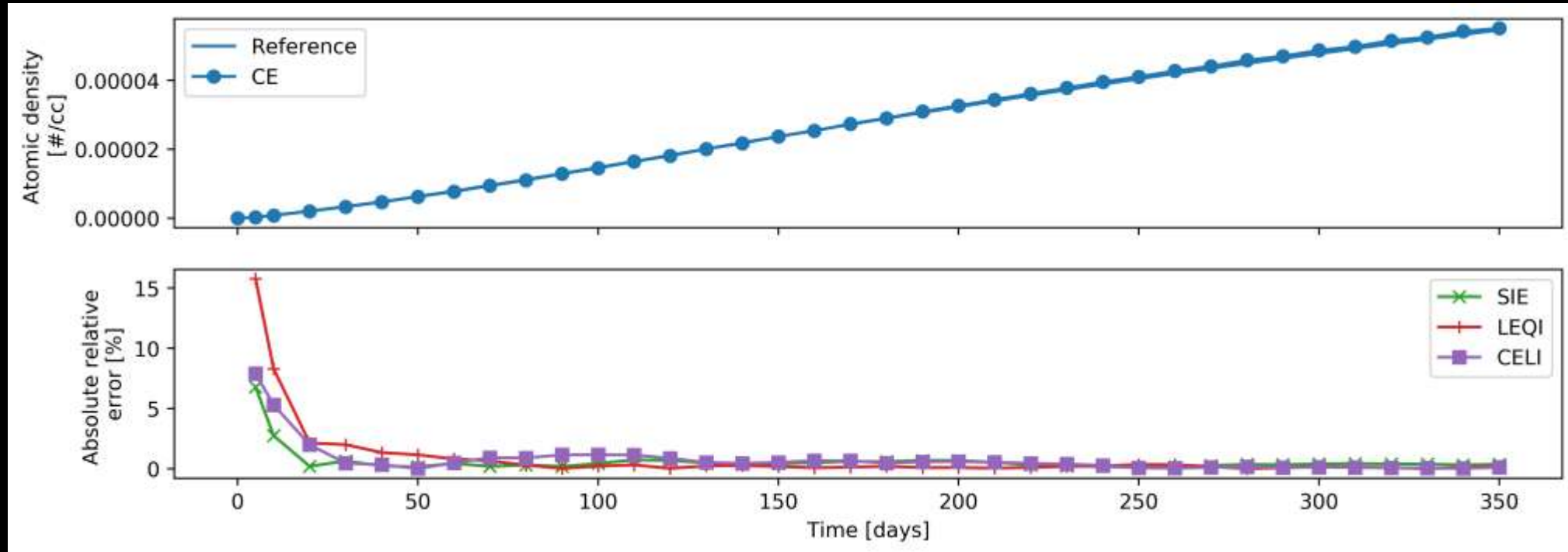
THANK YOU



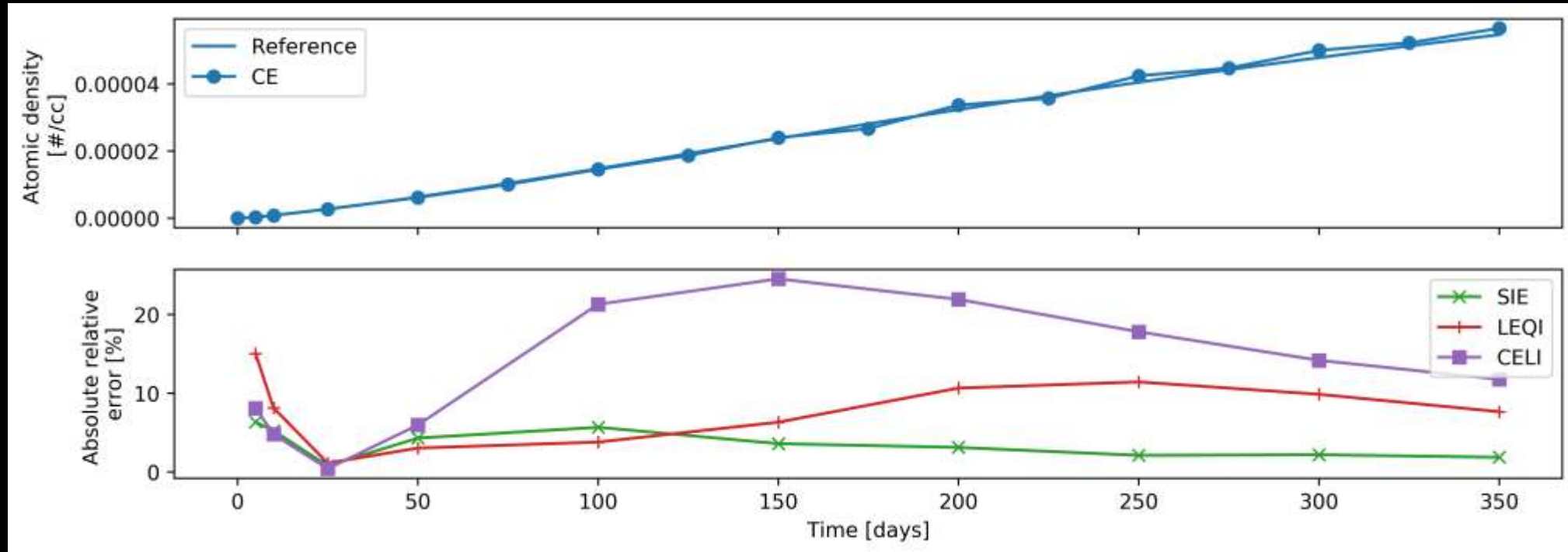
Error in cumulative Xenon density for full core.
10 day steps, plotted at 50 day increments



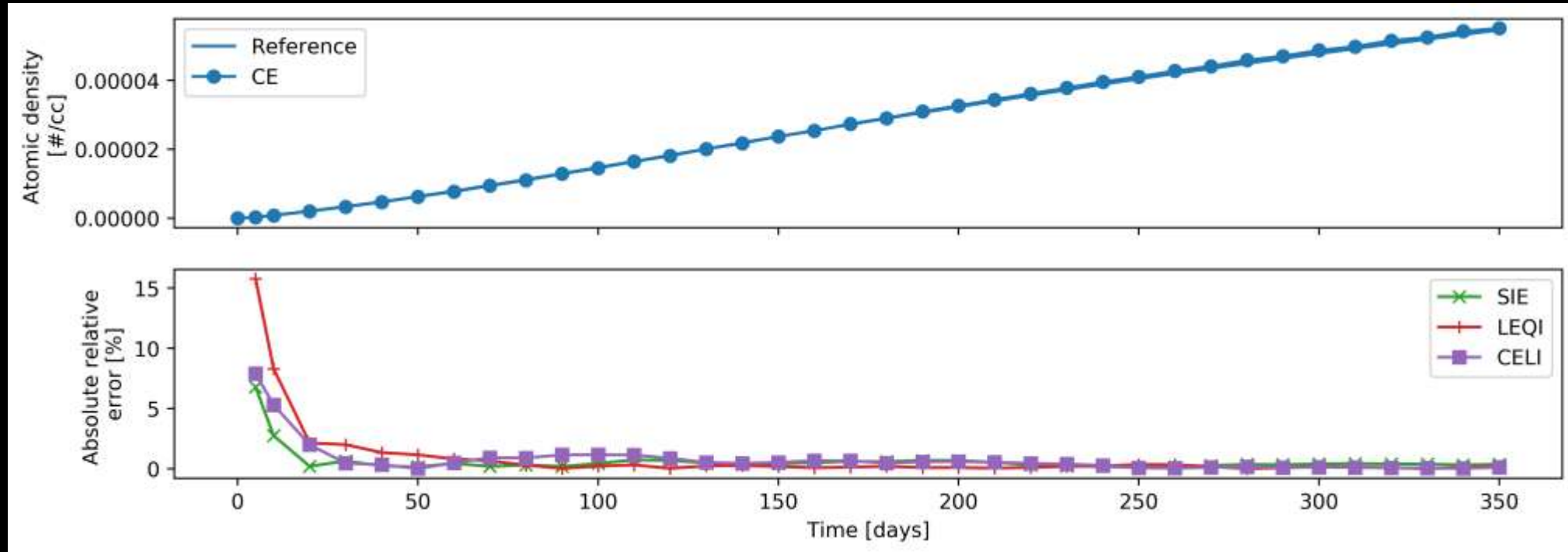
Cumulative error in Pu-239 density - 50 day steps



Absolute error in nodal Pu-239 using 10 day time steps



Absolute error in nodal Pu-239 using 50 day time steps



Absolute error in nodal Pu-239 density with 10 day steps