

Fast ion orbit modeling using a Chebyshev representation for the magnetic potential.

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- Information about the Lithium Tokamak Experiment.
- Chebyshev polynomials.
- The Chebyshev approximation in 1D and 2D.
- Computational methods.
- Simulation results.



Lithium Tokamak Experiment (LTX)

- Studies the plasma-material interaction of liquid Lithium (Li).
- B = 0.3 T.
- Neutral beam injection heating.
- Fast ion means energies > 10 keV.
- Why LTX is important?
 - ITER has a tungsten wall, which is high-Z (74) material.
 - LTX walls are coated by low-Z (3) Li to reduce recycling.
- **Recycling** when cool particles are re-introduced back to main plasma.





- Fast ions heat the plasma and supply particles to its core.
- They may escape either at first orbits or due to collisions with the plasma.
- Simulating the ion losses at collision time scales (> 10 ms) is a challenge.
- Past simulations used a bi-cubic spline approximation for calculations of magnetic field
 - Issues at large times because of discontinuous first derivative at mesh boundary.
- A Chebyshev approximation is developed here to resolve this issue.
- Goal: Simulate particle losses in a fast way



1D Chebyshev Polynomials

Definition of polynomials

 $T_0(x) = 1$ $T_1(x) = x$ $T_2(x) = 2x^2 - 1$ $T_3(x) = 4x^3 - 3x$ $T_4(x) = 8x^4 - 8x^2 + 1$ $T_5(x) = 16x^5 - 20x^3 + 5x$ $T_6(x) = 32 x^6 - 48 x^4 + 18 x^2 - 1$ $T_n(x) = \cos(n \arccos x).$ $T_{n+1} = 2xT_n - T_{n-1}, \quad n \ge 1.$ - $T_0(x)$ - $T_1(x)$ - $T_2(x)$ - $T_3(x)$ - $T_4(x)$ 0.5 0.0 -0.5-1.0-0.50.0 0.5 1.0



Complete orthonormal system.

Orthogonality identity $\sum_{k=1}^{n} T_i(\bar{x}_k) T_j(\bar{x}_k) = \begin{cases} 0 & i \neq j \\ n/2 & i = j \neq 0 \\ n & i = j = 0 \end{cases}$

Chebyshev nodes

$$\bar{x}_k = \cos \frac{\pi (2k-1)}{2n}, \quad k = 1, 2, 3, \dots, n.$$



$$f(x) \approx \sum_{i=0}^{N-1} c_i T_i(x),$$

$$c_0 = \frac{1}{N} \sum_{k=1}^{N} f(\bar{x}_k),$$

$$c_i = \frac{2}{N} \sum_{k=1}^{N} f(\bar{x}_k) T_i(\bar{x}_k),$$

$$\bar{x}_k = \cos \frac{\pi (2k-1)}{2n}, \quad k = 1, 2, 3, \dots, n.$$

- Approximate a function as a sum of the first N Chebyshev polynomials.
- The coefficients c_i are determined through summing the overlap at the Chebyshev nodes.



- M 1 N 1 $f(x,y) \approx \sum \sum c_{ij} T_i(x) T_j(y).$ $c_{00} = \frac{1}{MN} \sum_{k=1}^{M} \sum_{k=1}^{N} f(\bar{x}_k, \bar{y}_\ell),$ $c_{i0} = \frac{2}{MN} \sum_{i=1}^{M} \sum_{j=1}^{N} f(\bar{x}_k, \bar{y}_\ell) T_i(\bar{x}_k),$ $c_{0j} = \frac{2}{MN} \sum_{k=1}^{M} \sum_{j=1}^{N} f(\bar{x}_{k}, \bar{y}_{\ell}) T_{j}(\bar{y}_{\ell}),$ $c_{ij} = \frac{4}{MN} \sum_{k=1}^{M} \sum_{j=1}^{N} f(\bar{x}_k, \bar{y}_\ell) T_i(\bar{x}_k) T_j(\bar{y}_\ell).$
- This extends to 2D, where the functions are multiples of the Chebyshev polynomials.
 - The same orthogonality identities hold, so the coefficients c_{ij} are determined the same way.
 - Our code evaluates a 2D approximation of the magnetic potential.
 - First partial derivatives are used to calculate the magnetic field.



- GPU is much faster through massive parallelism.
- Useful for calculating the Chebyshev representation for > 100,000 particles.



- Varying particle number and measuring time to evaluate the magnetic field.
- Initial offset is due to loading variables onto GPU.





- Codes written in C and CUDA C.
- Use the Soloviev solution to the Grad-Shafranov equation to find the magnetic potential.
- Approximate the magnetic potential using the Chebyshev representation and find the first partial derivatives with respect to r and z.
- For each particle,
 - Calculate the magnetic fields using the Chebyshev representation.
 - Generate the time derivatives of particle position and velocities.
 - Use Runge-Kutta-4 algorithm for full orbit calculations.

Example particle orbits



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Example particle orbit in the r-z plane.







- I learned how to use C and CUDA to successfully build a partial kinetic code!
- GPU speeds up calculations.
- Observed orbits are consistent with theory.
- Issues at large time steps. Limited accuracy of RK4
- Fully kinetic codes are computationally intensive (1 million steps for each particle).
 - It is not feasible to simulate all particles, just the few important ones.
- The Chebyshev approximation, being compact and smooth, is suitable for GPU.
- We need more testing with realistic magnetic configurations.



Thank you!