Computational Modelling of the Rotamak

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Fields Used in the Model

For $z^2 < z \downarrow s \uparrow 2$:

For $z^2 \ge z is 12$ (escaped):

 $\boldsymbol{B} = (B\omega\cos(\omega t) - 1/2 \ x\pi Bm/zs) \\ \sin(\pi z/zs) \boldsymbol{i}$

B=(2*Bm*+*B*0)**k**

E=0

+ $(B\omega \sin(\omega t) - 1/2 \gamma \pi Bm/zs \sin(\pi z/zs))$

+ $(Bm(1-\cos(\pi z/zs))+B0)k$

 $\boldsymbol{E} = \omega B \omega (\boldsymbol{x} \cos(\omega t) + \boldsymbol{y} \sin(\omega t) \boldsymbol{k})$

Simulation Parameters

 $B_{\omega} = 10 \text{ Gauss}$ $B_0 = 200 \text{ Gauss}$ $R_{m} = 10 (B_{m} = 900 \text{ Gauss})$ $\omega = 5 \times 10^7 \text{ rad/s}$ $n = 1 \times 10^{19} \, \text{m}^{-3}$ $\ln\Lambda = 10$ $r_s = 5 \text{ cm}$ $z_{s} = 45 \text{ cm}$ $v = n\pi e t 4 \ln \Lambda / m t^2 v t^3$ (electron-ion collisions only)

Electrons initialize with random positions on the z=0 disk (with radius r_s), and are given spherically distributed random velocities, all with identical energies.



Escaping Electrons

- Electrons starting with higher energies are far more likely to be trapped.
- Most low energy electrons escape only after two cycles of the Rotamak (~250 ns).

% of Particles Escaping after 2 Cycles vs. Initial Kinetic Energy (sampled from 1000 particles per run)



Initial Kinetic Energy (eV)

1500

2000

2500

1000

10

0

0

500



Electron Heating

- Electrons initially under ~500eV heat rapidly within two cycles, and then level off.
- Further very gradual heating is due to scattering, taking around a quarter of a millisecond until most of the particles have escaped.



Heating Within the First Two Cycles



Future Work

- Find out how the heating rates change with the simulation parameters, especially with the value of B_{ω} .
- Initialize the particles outside of the bottle like in the machine.
- Find rate of energy dissipation via escaping particles.