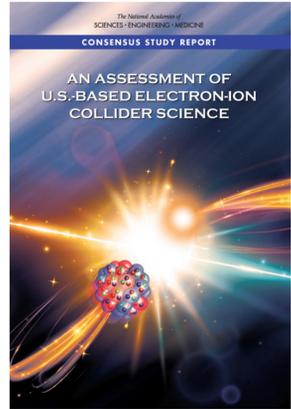


Magnetized Electron Cooling Simulations for JLEIC*

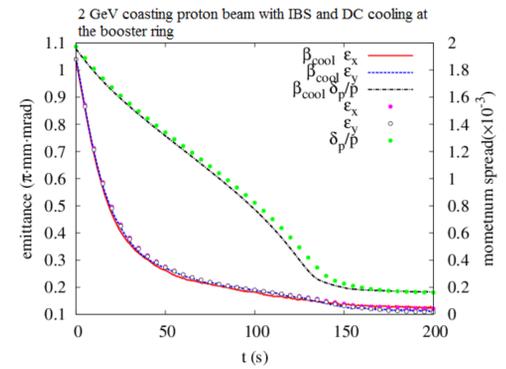
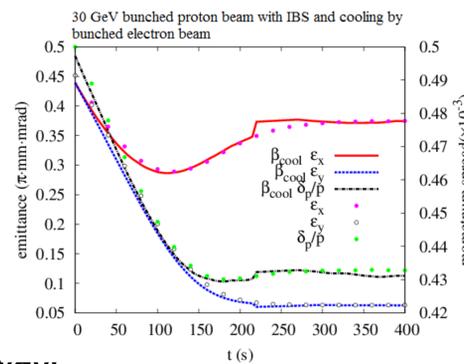
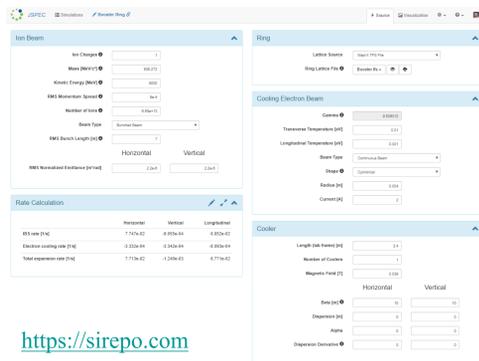


JSPEC – JLab Simulation Package for Electron Cooling

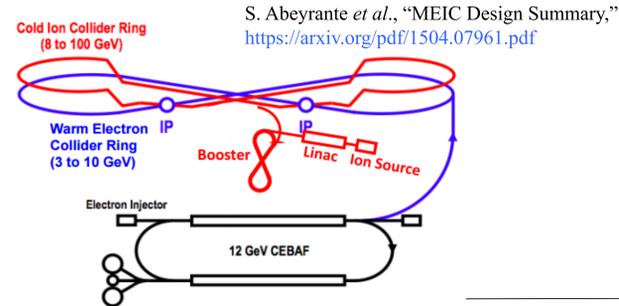
- Considers both IBS and electron cooling
 - calculate rates, simulate processes for specified ring
- C++, open source: <https://github.com/zhanghe9704/electroncooling>
- Successfully benchmarked with BETACOOOL →

H. Zhang, J. Chen, R. Li, Y. Zhang, H. Huang and L. Luo, "Development of the electron cooling simulation program for JLEIC," Proc. IPACPMW014 (2016).

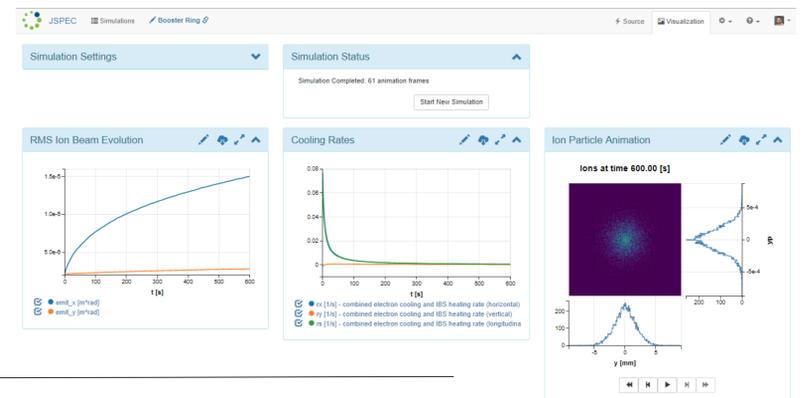
JSPEC is part of the **Sirepo** Scientific Gateway:



JLEIC Concept from Jefferson Lab

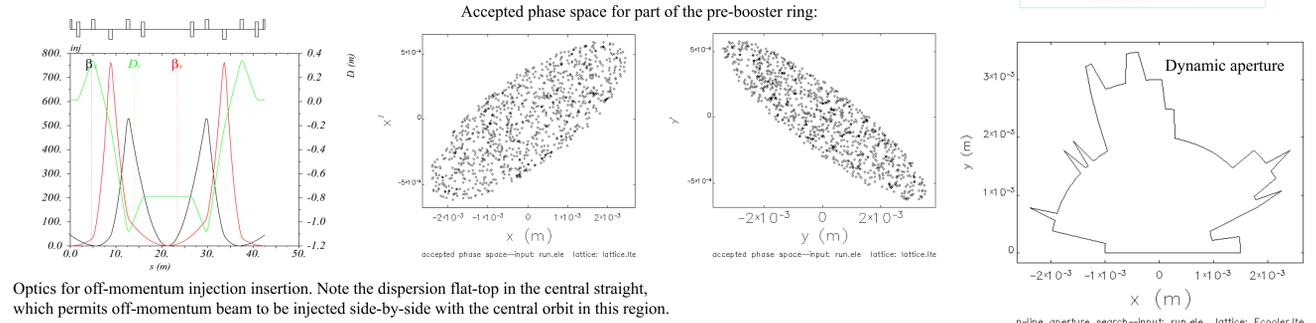


S. Abeyante *et al.*, "MEIC Design Summary," <https://arxiv.org/pdf/1504.07961.pdf>



Early concept of pre-booster for accumulation & cooling

- Suggested a few years ago by P.M. McIntyre *et al.*
 - further studied as part of this SBIR-funded project
- Ring design with MAD-X and elegant
 - elegant is also part of the Sirepo Scientific Gateway, <https://sirepo.com>



Dynamic Friction Modeling – general approach

THEORY OF ELECTRON COOLING
Yaroslav Derbenev*
Thomas Jefferson National Accelerator Facility, Newport News, VA 23606, USA

Ya. Derbenev, "Theory of Electron Cooling," arXiv (2017)

- The E-fields associated with friction must be carefully identified
 - fields from distribution of e⁻s perturbed by the presence of the ion

bulk fields friction statistical fluctuations

$$\vec{E}(\vec{r}, \vec{v}, t) = \langle \vec{E}^0 \rangle(\vec{r}, t) + \langle \Delta \vec{E} \rangle(\vec{r}, \vec{v}, t) + \vec{E}^{fl}(\vec{r}, \vec{v}, t) \quad (1.1)$$

- Friction force must be calculated along the ion trajectory:

$$\vec{F} = -ze(\Delta \vec{E})(\vec{r}, \vec{v}, t) \Big|_{\vec{r}=\vec{r}(t), \vec{v}=\vec{v}(t)} \quad (1.2)$$

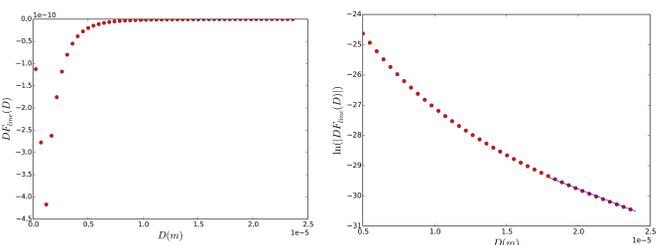
- we do this numerically for each individual ion-electron interaction
 - total force obtained by summing over e⁻ distribution (i.e. no shielding)
- bulk forces are removed by subtracting force from unperturbed e⁻s

Gyrokinetic Averaging Yields 1D Nonlinear e⁻ Oscillations

- Hamiltonian perturbation theory for single ion & e⁻
 - unperturbed motion: drifting ion and magnetized e⁻
 - primary assumption: D (impact parameter) $\gg r_L$ (Larmor radius)
 - longitudinal dynamics: $V_{ion, \parallel} = 0$ (to be relaxed in future work)
- e⁻ gyrocenters stay on cylinder of constant radius
- choose ion to be stationary at the origin (convenient coordinates)
 - gyrocenters move in a nonlinear 1D potential:
 - weak nonlinearity
 - larger amplitudes => longer periods
- shortest oscillation period:
 - both trapped and passing orbits
 - numerical simulations are required to capture these effects
- no logarithmic singularity for $D \rightarrow 0$ or $D \rightarrow \infty$

$$\tilde{z}(t) = -\frac{Ze^2}{4\pi\epsilon_0 m_e} \frac{z}{(D^2 + z^2)^{3/2}}$$

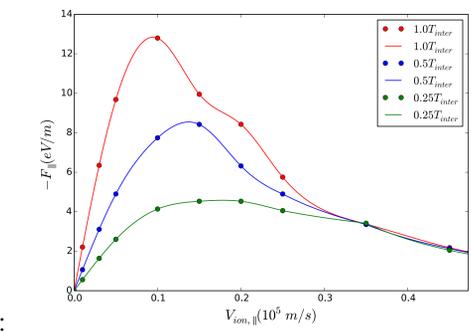
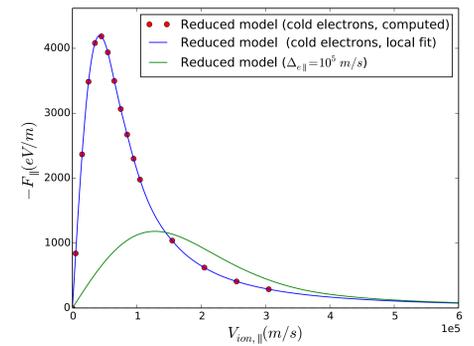
$$T_{lin} = \frac{2\pi}{c} \sqrt{\frac{D^3}{Zr_e}}$$



$$F_{\parallel}(V_{\perp} = 0) = 2\pi n_e \int_0^{\infty} dD DF_{line}(D) \equiv 2\pi n_e \int_0^{\infty} dD \int_{-\infty}^{\infty} dz_{im} F_{i-e}(z_{im}, D)$$

$F_{\parallel}(V_{ion, \parallel})$ for cold e⁻, short times: scaling in Z, T_{int}

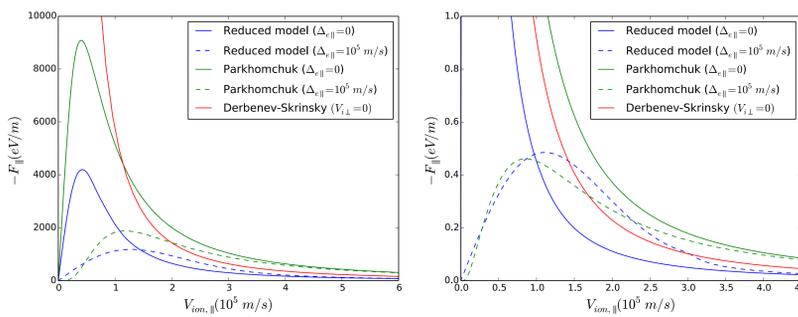
- Considered protons and Au⁺⁷⁹ ions
 - with different interaction times in the cooler
 - for small V , $dF_{\parallel}(V)/dV \approx 2Z n_e m_e r_e c^2 T_{int}$
 - at large V , $F_{\parallel} \approx 2\pi Z^2 n_e m_e (r_e c^2)^2 / V^2$
 - with no dependence on T_{int}
 - for a given T_{int} , peak friction force scales as $Z^{4/3}$
- For $T_{int} < T_{plasma}$ and small-to-moderate V_{ion}
 - $F_{\parallel}(V_{ion, \parallel})$ increases with interaction time, while the large- V tail is independent of T_{int}
- $F_{\parallel}(V_{ion, \parallel})$ is linear in n_e by construction
- Thermal effects are computed via convolution of $F_{\parallel}(V_{ion, \parallel})$ with warm longitudinal electron distribution



Upper panel: Gold ion, cold and warm e⁻s;
Lower panel: Protons, cold e⁻s, varying T_{int}

Comparison with Derbenev-Skrinsky and with Parkhomchuk models

- Comparison of new model for Au⁺⁷⁹ ions and protons with:
 - Derbenev and Skrinsky (D-S) for $V_{ion, \perp} = 0$ and large $V_{ion, \parallel}$
 - Ya. S. Derbenev and A.N. Skrinsky, Part. Accel. 8 (1978), 235
 - Parkhomchuk (P) with 0 and finite effective long. e⁻ temp
 - V.V. Parkhomchuk, Nucl. Instr. Meth. in Phys. Res. A 441 (2000), p. 9
- Consistently lower force than D-S and P for larger $V_{ion, \parallel}$
 - for lower ion velocity, details depend on Z
 - for cold electrons, new model shows consistently lower force values than Parkhomchuk at ALL velocities



Left panel: Gold ion, cold and warm e⁻s; Right panel: Protons, cold and warm e⁻s

Parametric fit yields physical insight

