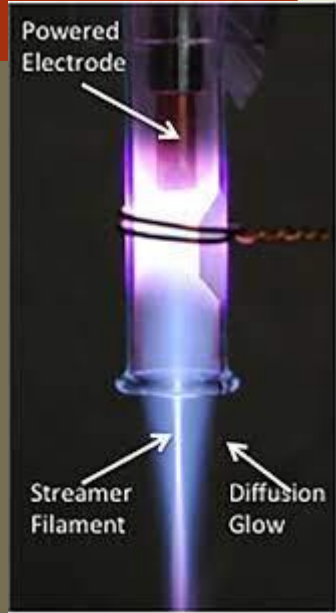
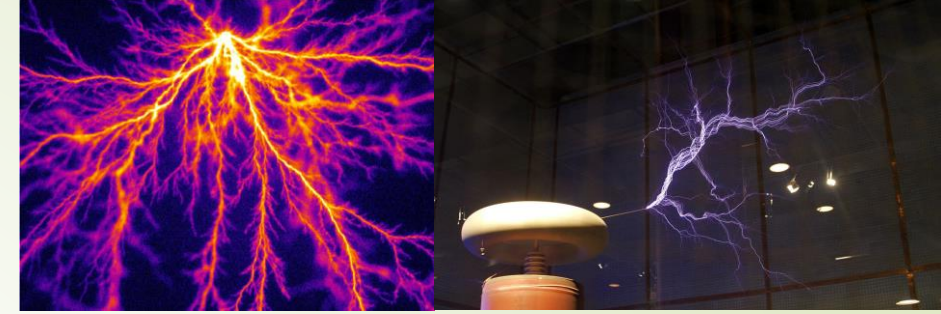




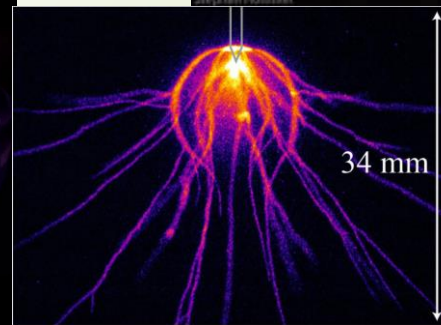
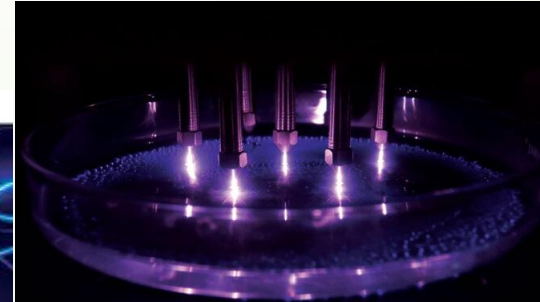
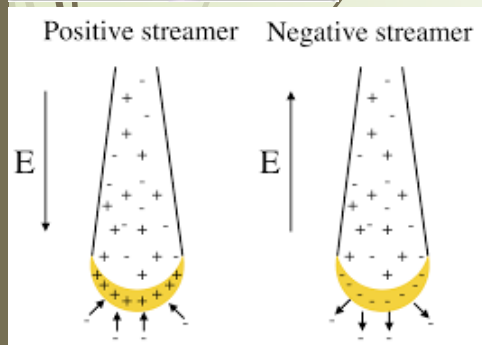
Simulations of positive and negative streamers in the AMReX environment


Ilija Simonović, Danko Bošnjaković and Saša Dujko

What are streamers?



- Thin channels of weakly-ionized nonstationary plasma produced by an ionization front that moves through non-ionized matter
- Streamers in nature: lightning and sprite discharges in the upper planetary atmospheres
- Applications of streamers:
 - ignition of high-intensity discharge lamps
 - treatment of polluted gases and water
 - plasma medicine



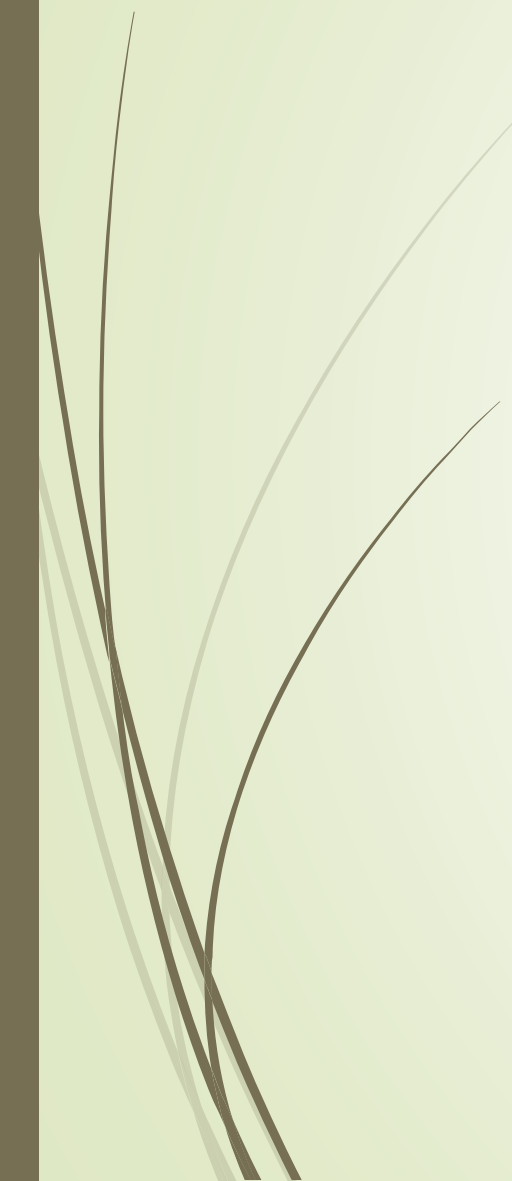


Streamer modelling

► Motivation:

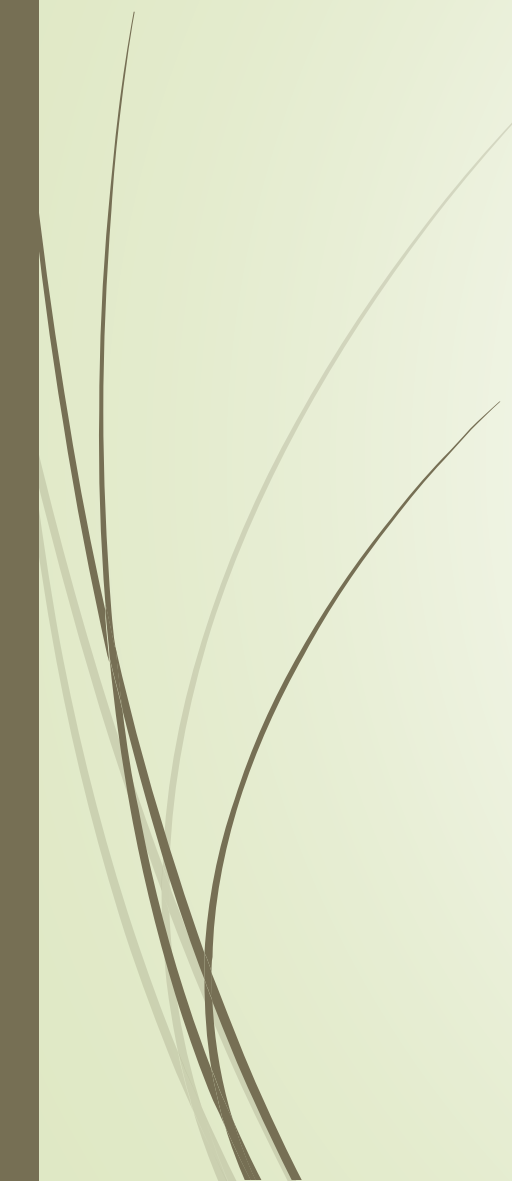
- better understanding of streamer physics
- optimization of streamer applications

► Types of streamer models

- Particle models
 - Fluid models
 - Hybrid models
- 



Fluid models of streamers

- ▶ Number densities of charged particles are represented by continuous functions
 - ▶ Time evolution of these number densities are represented by the fluid equations
 - ▶ Fluid models are generally more computationally efficient than the particle models and are much simpler than the hybrid models
- 

First-order fluid model

Advection diffusion reaction equation for the time evolution of the number density of electrons:

$$\frac{\partial n_e}{\partial t} + \nabla(n_e \mathbf{W} - \mathbf{D} \nabla n_e) = n_e(\alpha - \eta)|\mathbf{W}| + S_{ph}$$

Reaction equations for the time evolution of the number densities of ions:

$$\frac{\partial n_p}{\partial t} = n_e \alpha |\mathbf{W}| + S_{ph}$$

$$\frac{\partial n_n}{\partial t} = n_e \eta |\mathbf{W}|$$

Local field approximation

Total electric field:

$$\mathbf{E} = \mathbf{E}_{applied} - \nabla \Phi_{space_charge}$$

$$\Delta \Phi_{space_charge} = -q_e \frac{n_p - n_e - n_n}{\epsilon_0}$$

Photoionization

➤ In air photons emitted from excited nitrogen molecules can ionize oxygen molecules

➤ Zheleznyak model:

$$S_{ph}(\mathbf{r}) = \int d^3r' \frac{I(\mathbf{r}')f(|\mathbf{r}-\mathbf{r}'|)}{4\pi|\mathbf{r}-\mathbf{r}'|^2}$$

➤ The photon source term

$$I(\mathbf{r}) = \frac{p_q}{p+p_q} \xi^B \frac{\nu_u}{\nu_i} S_i(\mathbf{r})$$

➤ Ionization source term:

$$S_i(\mathbf{r}) = \alpha |\mathbf{W}|$$

➤ The absorption function:

$$\int_0^\infty f(r)dr = 1$$

Photoionization

- The absorption function can be represented as:

$$f(|\mathbf{r} - \mathbf{r}'|) = p_{O_2}^2 |\mathbf{r} - \mathbf{r}'| \sum_{j=1}^N A_j^B e^{-\lambda_j^B p_{O_2} |\mathbf{r} - \mathbf{r}'|}$$

- Leading to the decomposition of the source term:

$$S_{ph} = \sum_{j=1}^N S_{ph,j}$$

$$S_{ph}(\mathbf{r}) = \frac{p_q}{p+p_q} \xi^B \frac{\nu_u}{\nu_i} \sum_{j=1}^N p_{O_2}^2 A_j^B \int d^3 r' \frac{S_i(\mathbf{r}') e^{-\lambda_j^B p_{O_2}^2 |\mathbf{r} - \mathbf{r}'|}}{4\pi |\mathbf{r} - \mathbf{r}'|}$$

- The components of the source term can be determined by solving a set of Helmholtz equations:

$$(\Delta - (p_{O_2} \lambda_j^B)^2) S_{ph,j} = - \left(A_j^B p_{O_2}^2 \frac{p_q}{p+p_q} \xi^B \frac{\nu_u}{\nu_i} \right) S_i$$

- Parametrizations: Luque, Bourdon 2 term, Bourdon 3 term



Spatial discretization: Finite volume method

- Scalar variables are defined at cell centers
- Vector variables are defined at cell faces
- Electric field components at the cell centers are interpolated from the cell faces
- Interpolation of the number density of electrons from the cell centers to the cell faces, to calculate electron fluxes, is not trivial!
- First order upwind scheme creates too much numeric diffusion, while linear interpolation (central differencing) creates strong numeric oscillations.
- For this reason, flux limiting schemes are used.

Koren flux limiter

- The flux of electrons is calculated as:

$$f_{i+\frac{1}{2}} = \begin{cases} W_{i+\frac{1}{2}} \left(n_i + \frac{1}{2} \phi(r_{i+\frac{1}{2}}^+) (n_i - n_{i-1}) \right), & \text{for } W_{i+\frac{1}{2}} > 0 \\ W_{i+\frac{1}{2}} \left(n_{i+1} + \frac{1}{2} \phi(r_{i+\frac{1}{2}}^-) (n_{i+1} - n_{i+2}) \right), & \text{for } W_{i+\frac{1}{2}} < 0 \end{cases}$$

- The Koren flux limiter is defined as:

$$\phi(r) = \max \left(0, \min \left(2r, \min \left(\frac{1}{3} + \frac{2r}{3}, 2 \right) \right) \right)$$

- The upwind ratio of consecutive solution gradients:

$$r_{i+\frac{1}{2}}^+ = \frac{n_{i+1} - n_i + \epsilon}{n_i - n_{i-1} + \epsilon}$$

$$r_{i+\frac{1}{2}}^- = \frac{n_i - n_{i+1} + \epsilon}{n_{i+1} - n_{i+2} + \epsilon}$$

Time integration

- ▶ Time integration is performed by employing the 2nd order Runge-Kutta method.
- ▶ Electric potential of space charge, resulting electric field, and transport coefficients are calculated at each stage of the Runge-Kutta method
- ▶ Time step restrictions:

- The CFL condition:

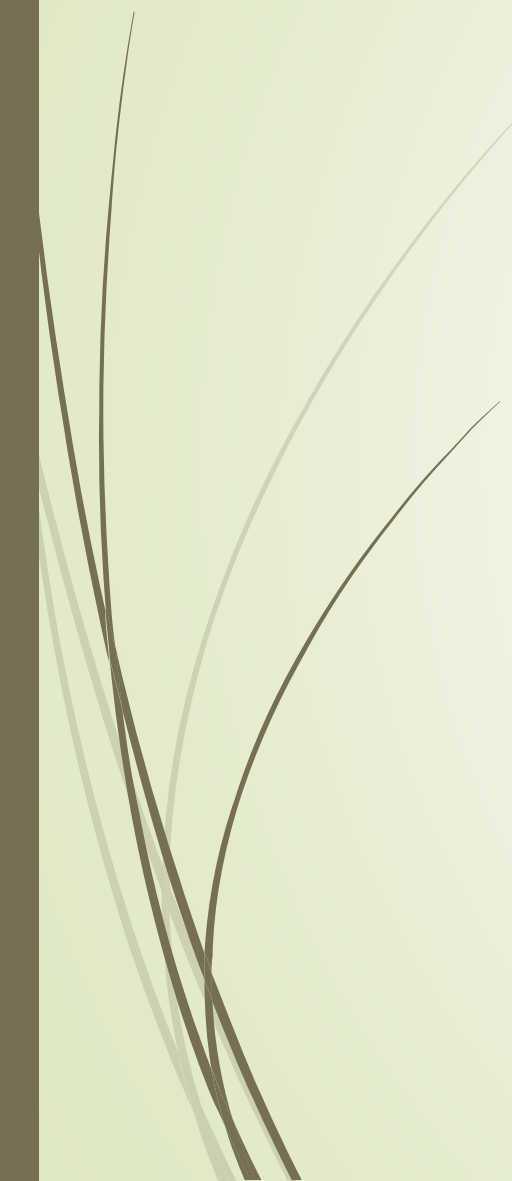
$$\Delta t \left(\sum_{i=1}^{N_{dim}} \frac{|W_i|}{\Delta x} + \frac{2N_{dim}D}{\Delta x^2} \right) < \text{CFL}, \quad \text{CFL}=0.4$$

- The dielectric relaxation time:

$$\Delta t < \frac{\epsilon_0}{q_e \mu_e n_e}, \quad \mu_e = \frac{|W(|E_{total}|)|}{|E_{total}|}$$

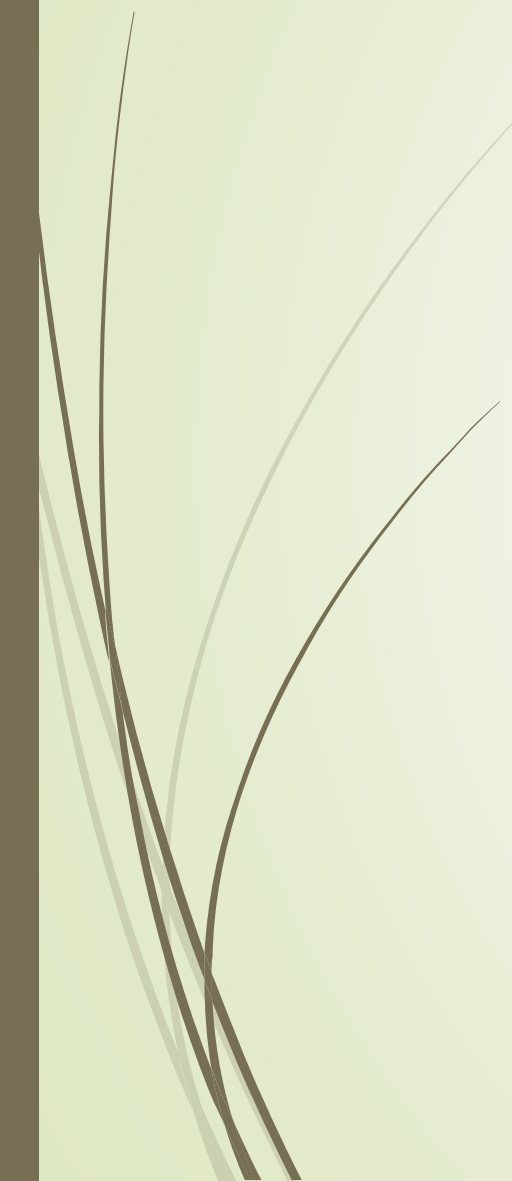


The AMReX library

- An open-source C++ library for massively parallel, block-structured adaptive mesh refinement (AMR) applications
 - Has inbuilt geometric multigrid solvers
 - Has many inbuilt classes which enable a convenient implementation of both grid and particle data
 - Allows both MPI and OpenMP parallelization, as well as parallelization on graphic processing units
- 



Adaptive mesh refinement

- Adaptive mesh refinement is used when high precision is required only in the subset of the calculation domain
 - Using a uniform mesh with high resolution is impractical and inefficient under such circumstances
 - The streamer dynamics is determined by the electron dynamics in the narrow region at the streamer front
 - Thus, a high-resolution mesh is required at the streamer front, while a coarser mesh can be used in the streamer channel, and in the other parts of the domain, which are far from the streamer front
- 

Refinement criteria

- The refinement criterion due to ionization frequency:

$$\alpha_{eff}(c_1|\mathbf{E}_{total}|)\Delta x > c_0, \quad \alpha_{eff} = \alpha - \eta \quad c_1 = 1.2, \quad c_0 = 0.8$$

- The refinement criterion due to number density of space charges:

$$\frac{\Delta x^2 |\rho|}{\epsilon_0} > c_2, \quad c_2 = 0.1V$$

- The de-refinement criterion:

$$\alpha_{eff}(c_1|\mathbf{E}_{total}|)\Delta x < 0.1 \wedge \Delta x < 30\mu m$$



Solving elliptic equations

- The Poisson equation and the Helmholtz equation are elliptic equations
- Their solution in each part of the domain depends on the values of the right-hand side in the entire domain
- Geometric multigrid method is very efficient in solving elliptic equations
- The method consists of applying a simple iterative procedure (like Gauss-Seidel or Jacobi) across a hierarchy of grids with reducing resolution
- AMReX includes inbuilt geometric multigrid solvers for the Poisson equation and the Helmholtz equations.
- These multigrid solvers can be easily applied across the entire hierarchy of adaptive mesh refinement levels



Results: Simulation conditions

- Axisymmetric model
- Domain size: [0 mm, 16 mm] along both coordinates
- Number of points along each axis at the coarsest level: 64
- Boundary conditions:
 - For the number density of electrons: Zero Neumann conditions at all boundaries
 - For the electric potential of space charge and photoionization source terms: Zero Neuman conditions at boundaries perpendicular to the radial axis and zero Dirichlet conditions at boundaries perpendicular to the axial coordinate
- Results of our AMReX-streamer code are compared to the results of the open-source Afivo-streamer code

Two-headed streamer in the Titan mixture

- The Titan mixture: 98.4% N₂, 1.6% CH₄
- The number of AMR levels 7 (from 0 to 6)
- No photoionization
- The applied reduced electric field: $E/N_0 = 147 \text{ Td}$,
 $1\text{Td} = 10^{-21} \text{ V/m}$, $N_0 = 2.50475764 \cdot 10^{25} \text{ m}^{-3}$.
- The initial condition: Neutral Gaussian given by:

$$n_e = n_p = n_{background} + n_0 \cdot e^{-\frac{(z-z_0)^2 + r^2}{\sigma^2}}, \quad n_n = 0 \text{ m}^{-3}$$

$$n_{background} = 10^{14} \text{ m}^{-3}$$

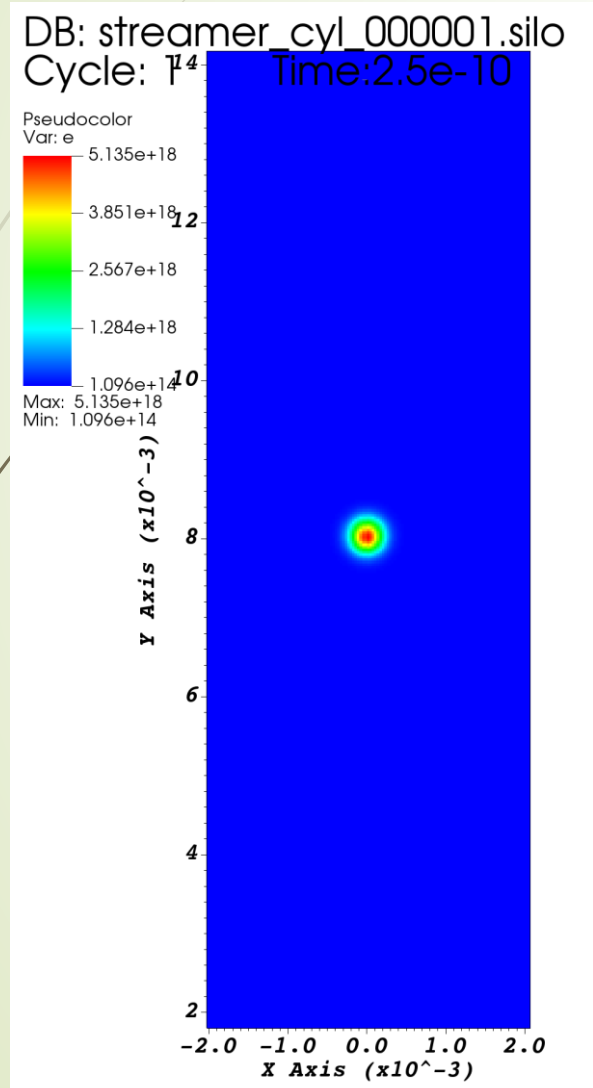
$$n_0 = 5 \cdot 10^{18} \text{ m}^{-3}$$

$$z_0 = 8 \text{ mm}$$

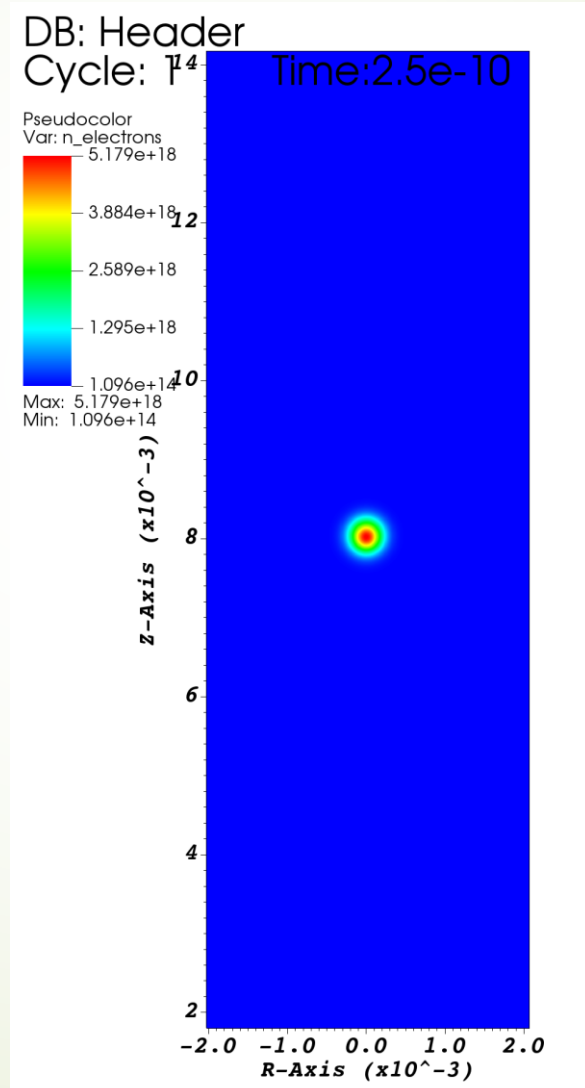
$$\sigma = 4 \cdot 10^{-4} \text{ m}$$

Number density of electrons at $t = 0.25$ ns

Afivo-streamer

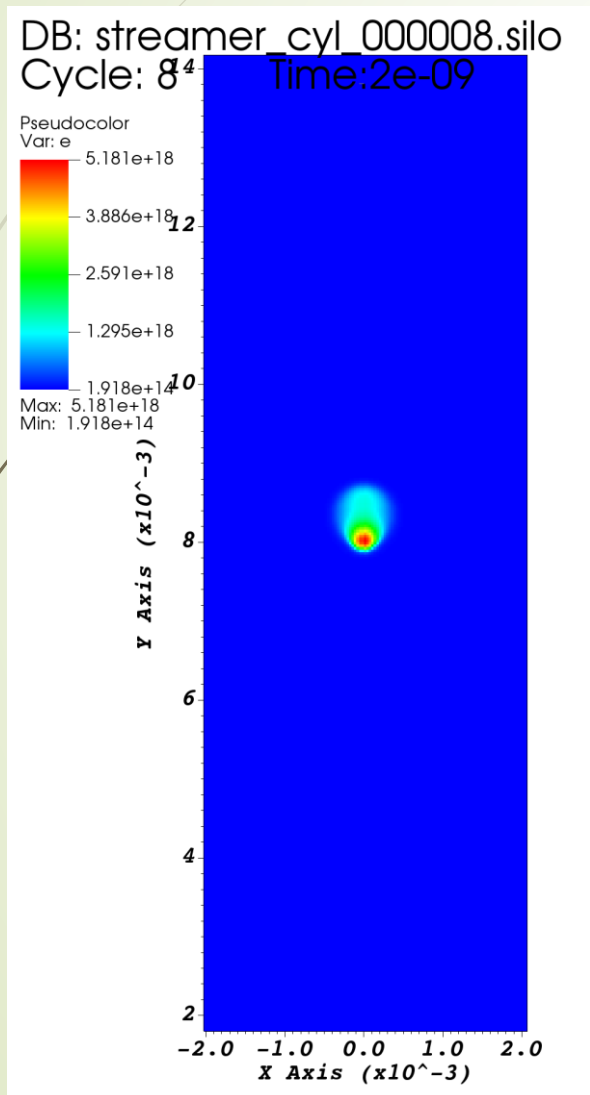


AMReX-streamer

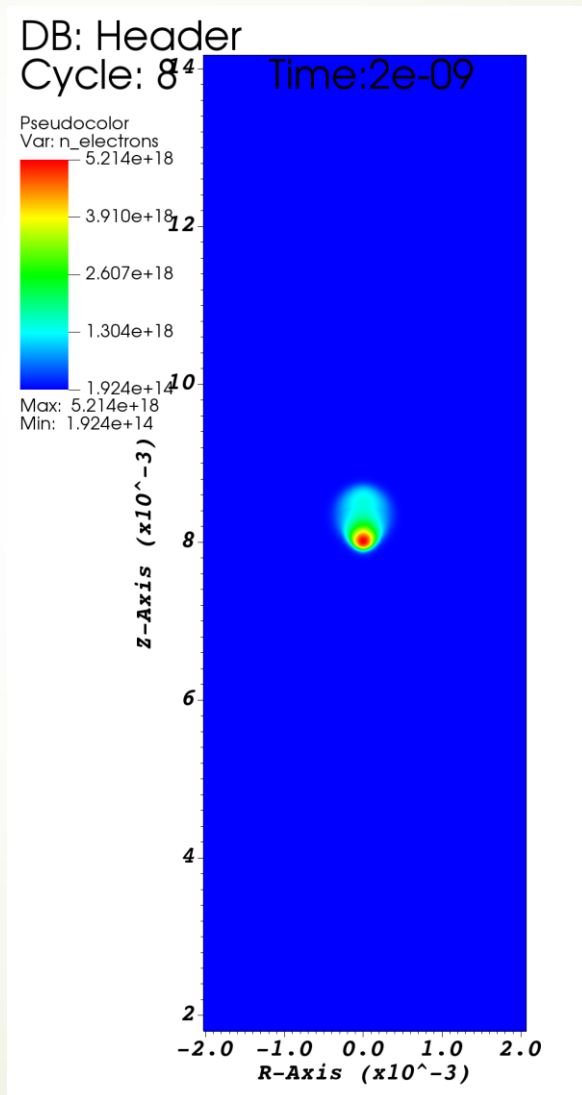


Number density of electrons at $t = 2 \text{ ns}$

Afivo-streamer

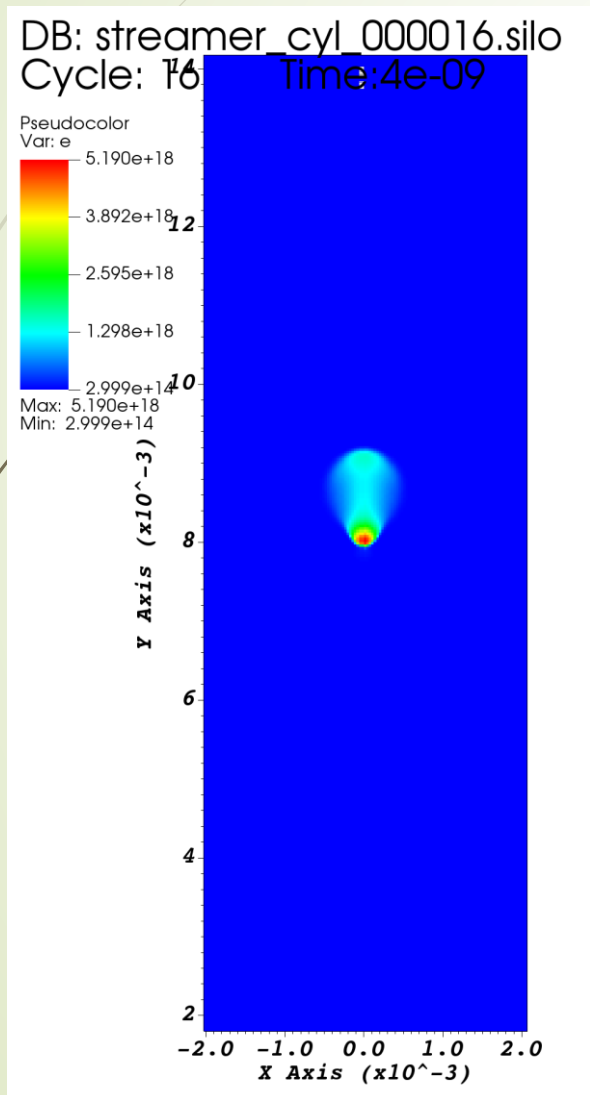


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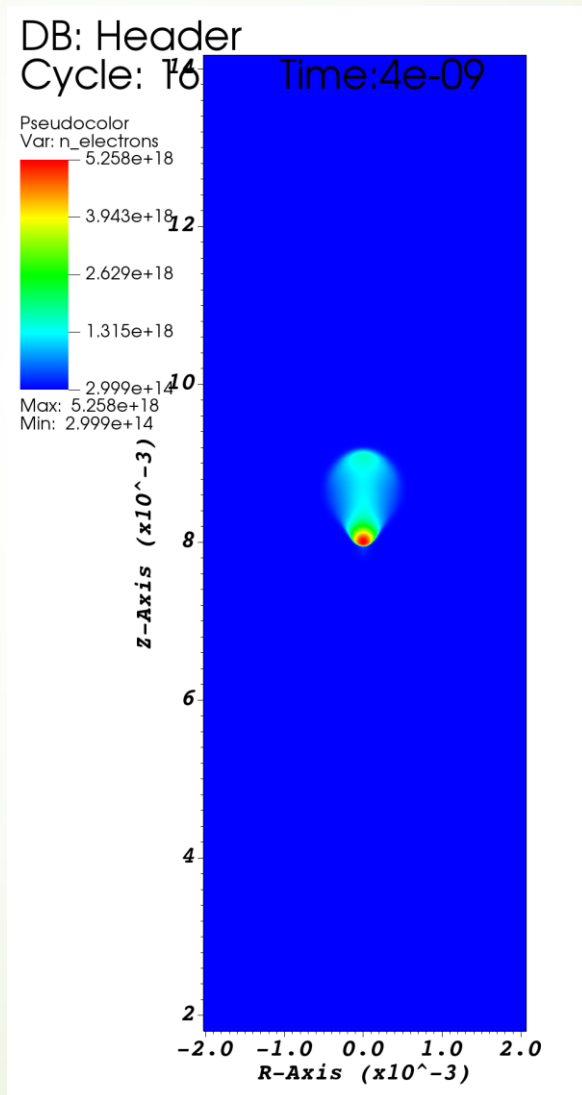


Number density of electrons at $t = 4$ ns

Afivo-streamer

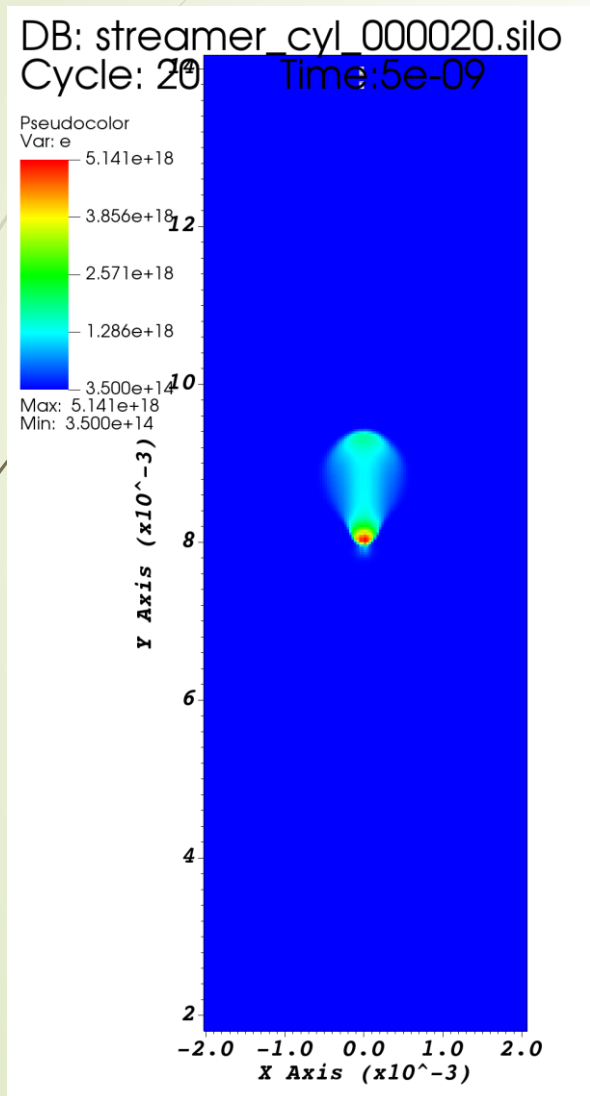


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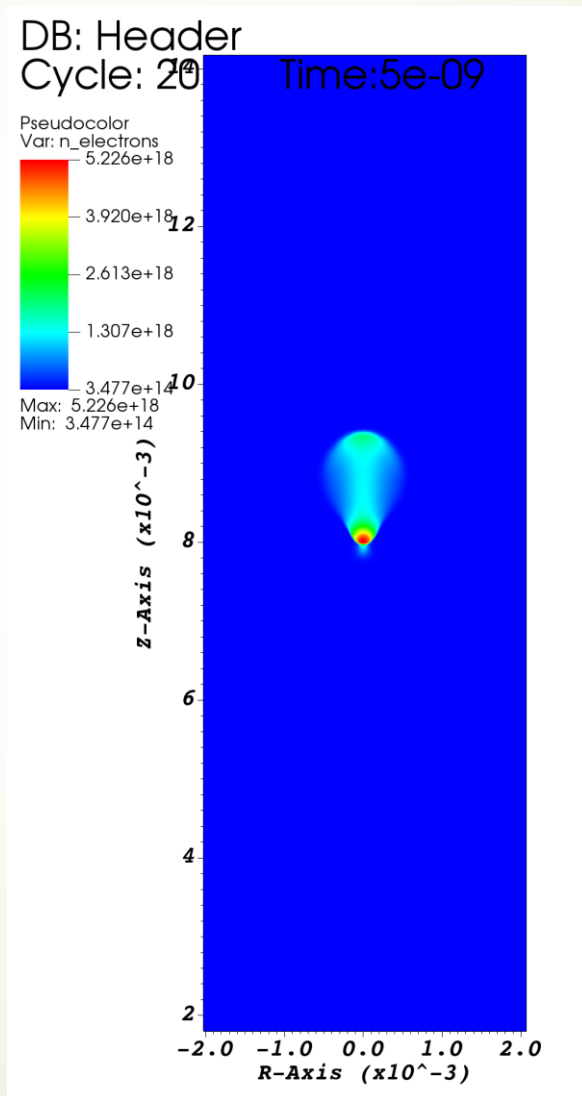


Number density of electrons at $t = 5$ ns

Afivo-streamer

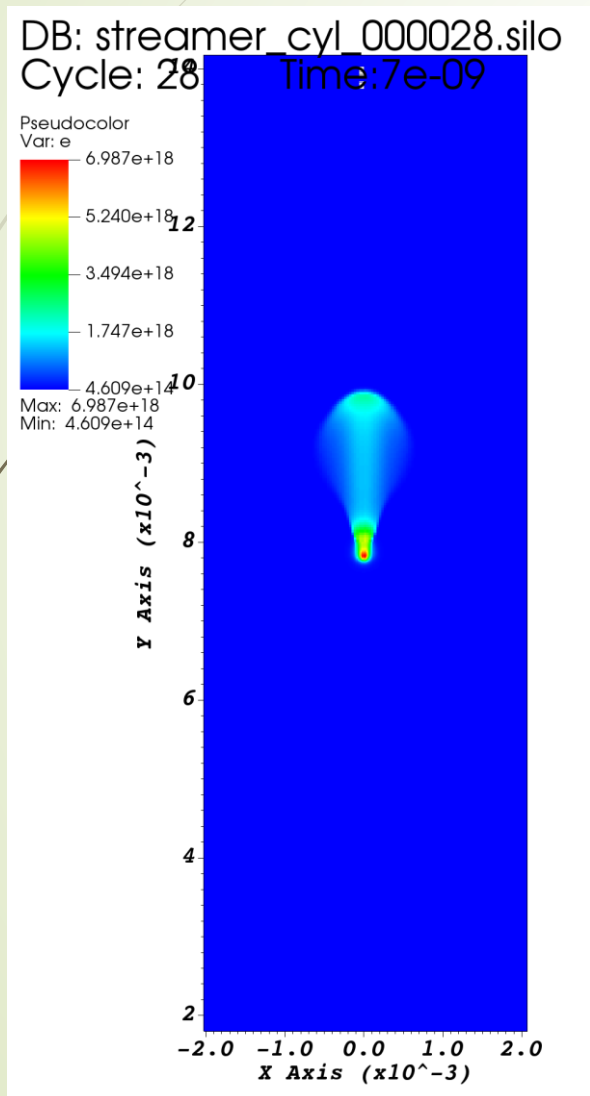


AMReX-streamer

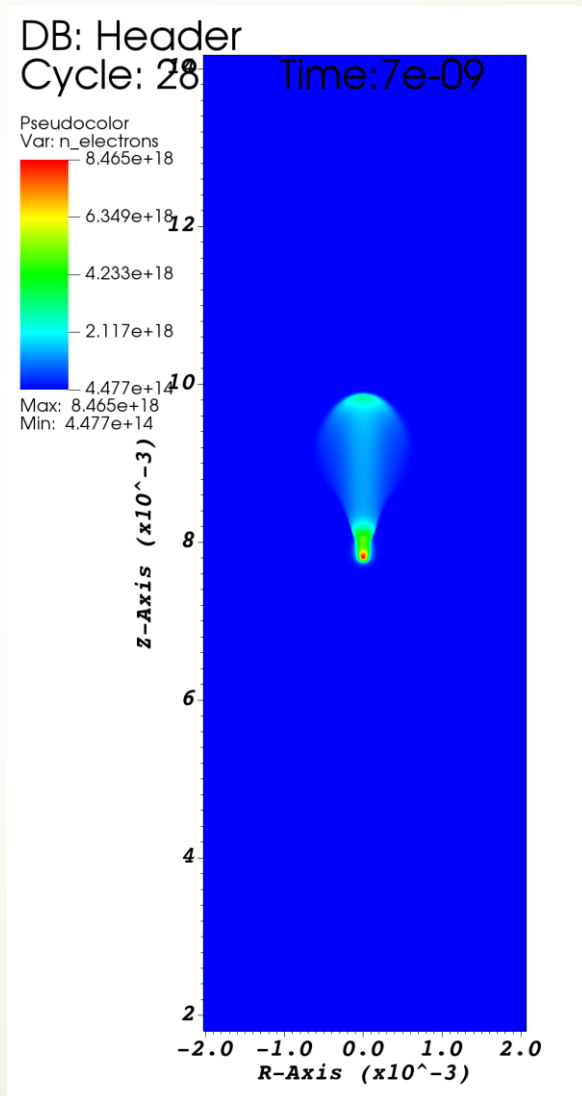


Number density of electrons at $t = 7$ ns

Afivo-streamer

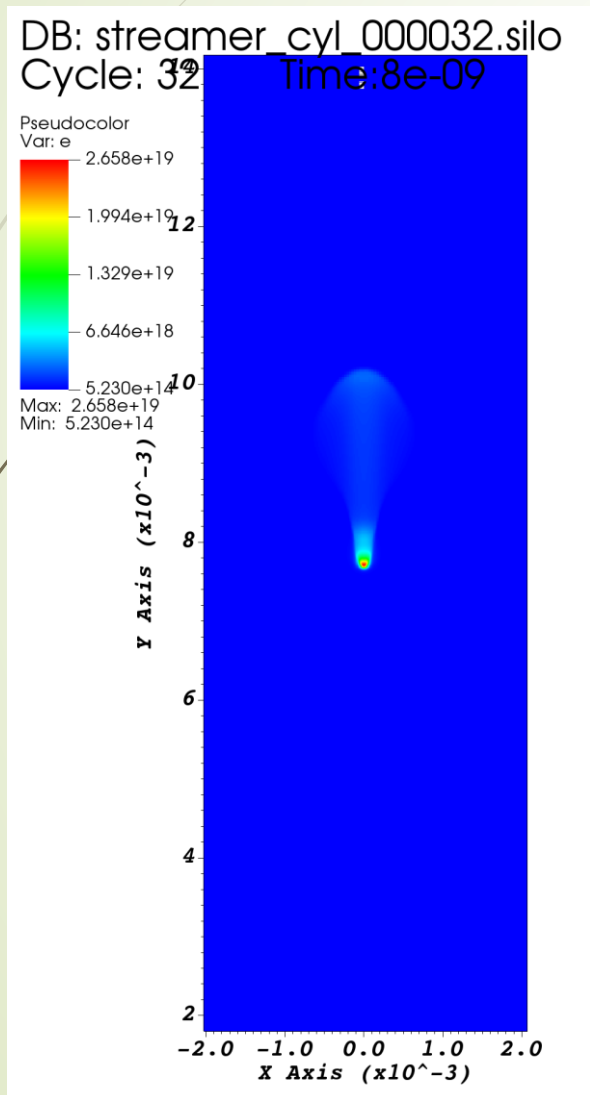


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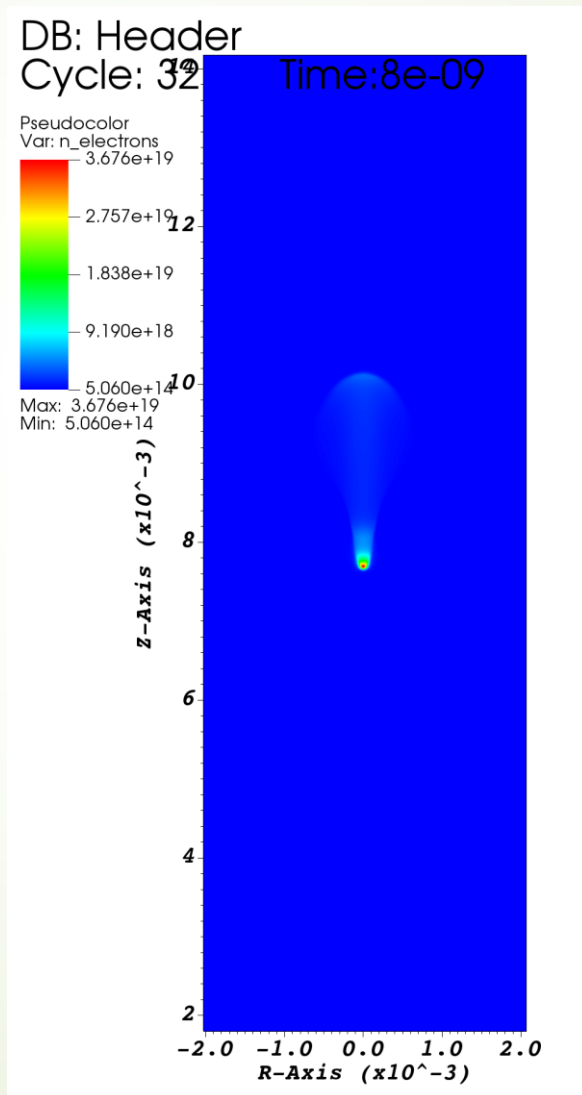


Number density of electrons at $t = 8$ ns

Afivo-streamer

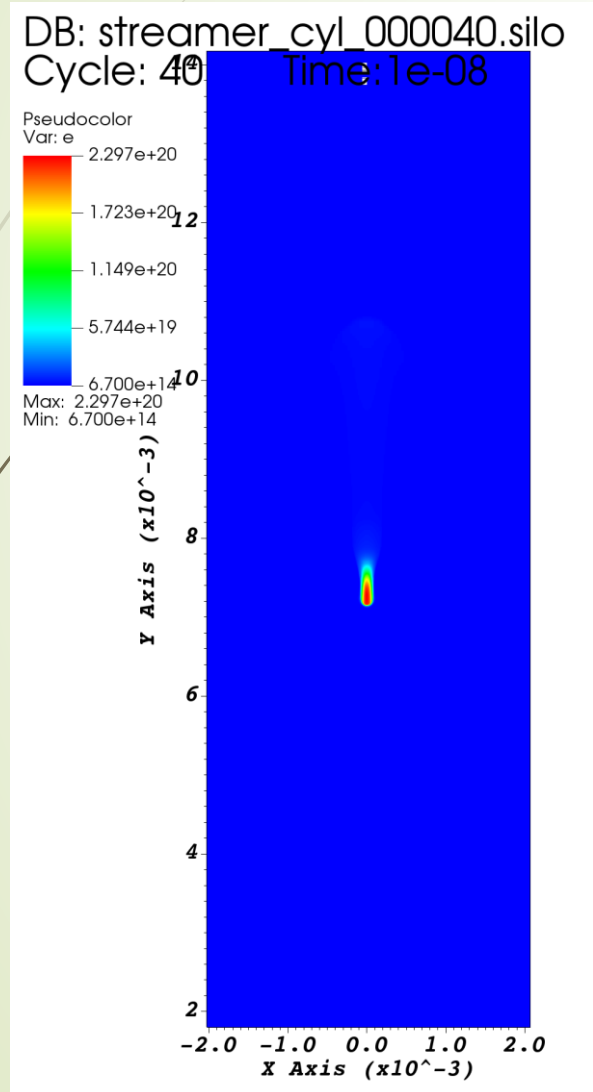


AMReX-streamer

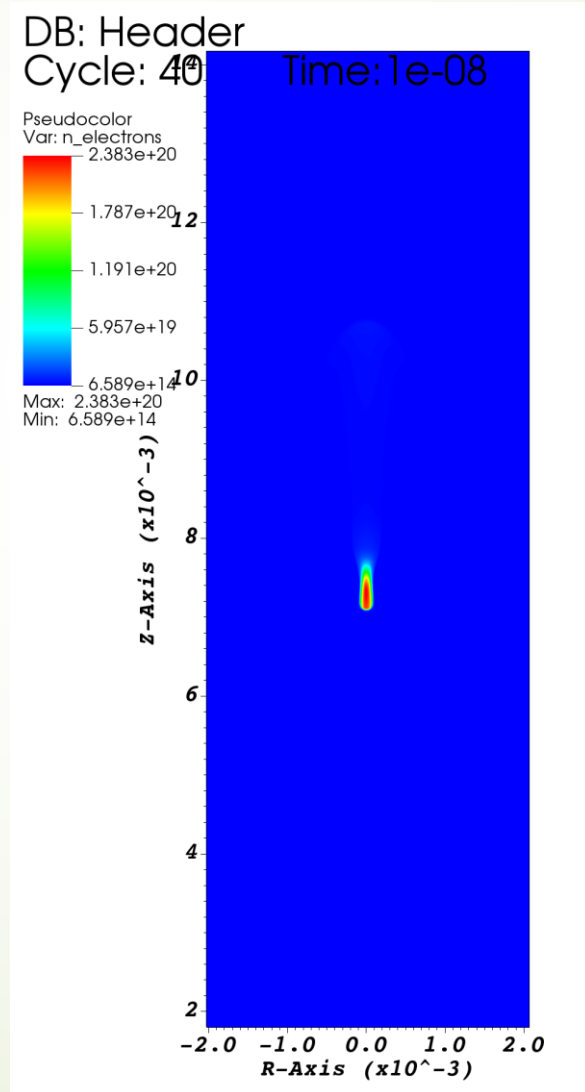


Number density of electrons at $t = 10$ ns

Afivo-streamer

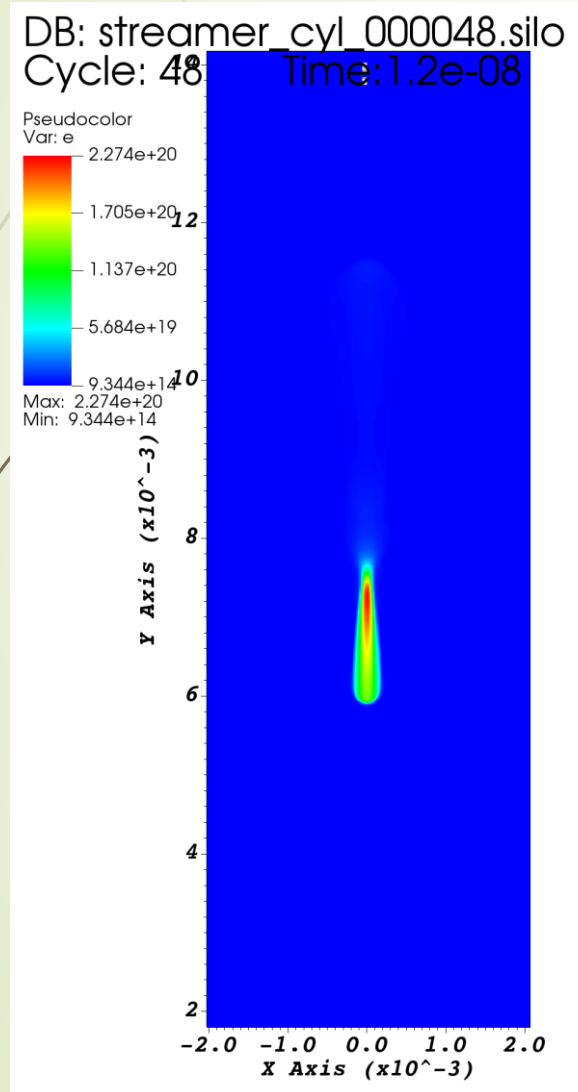


AMReX-streamer

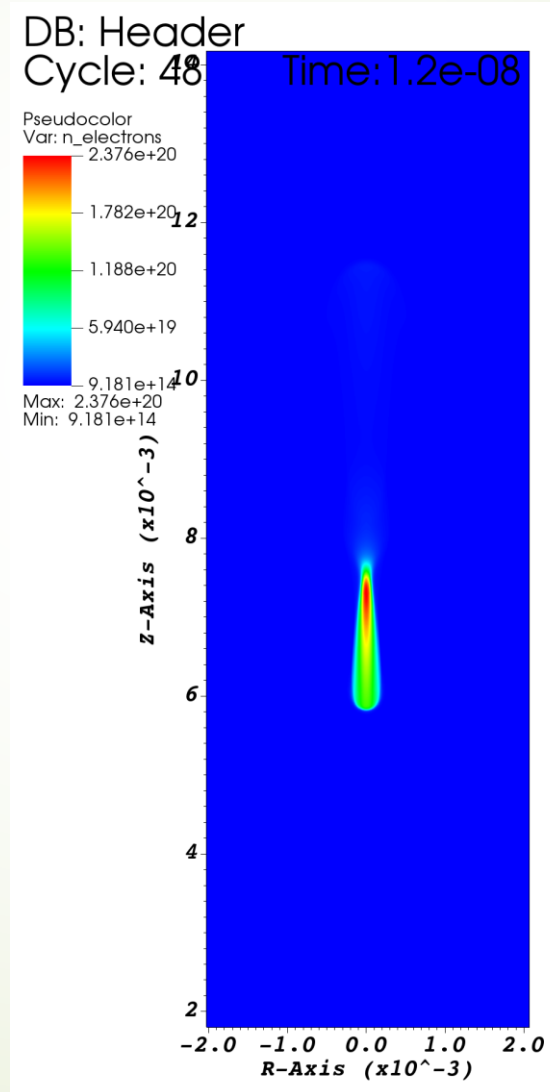


Number density of electrons at $t = 12 \text{ ns}$

Afivo-streamer

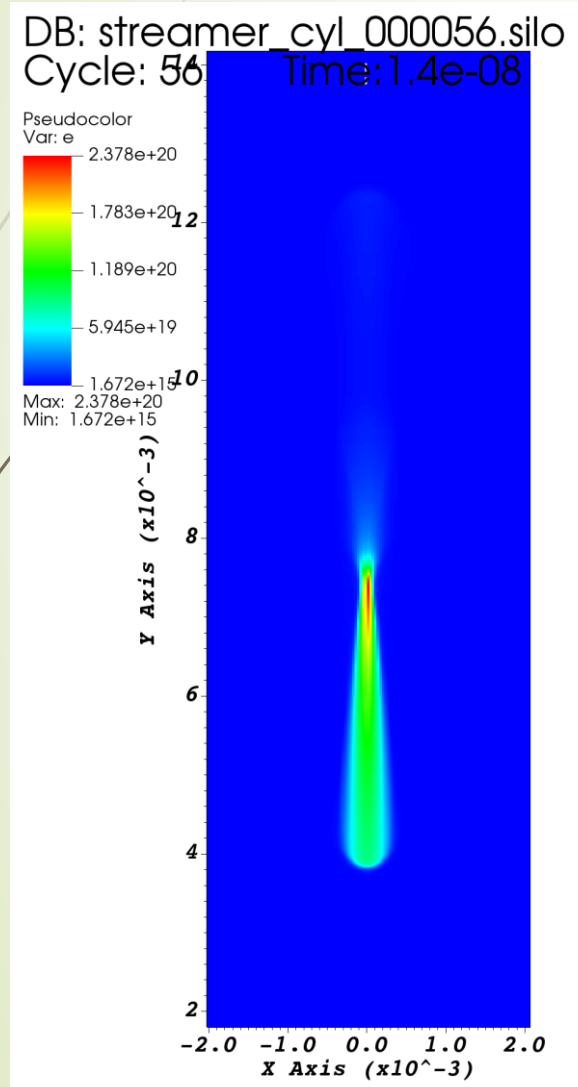


AMReX-streamer

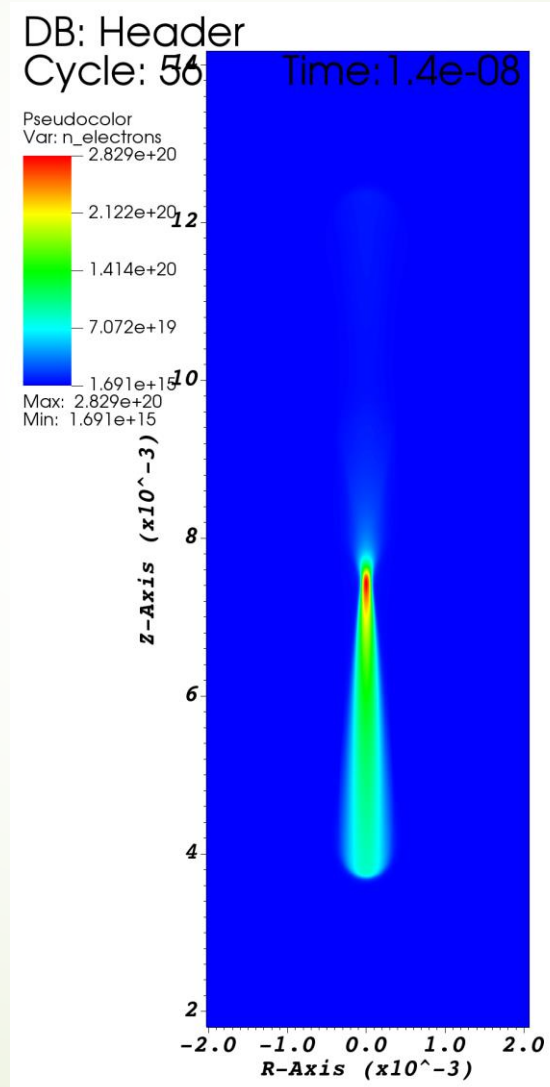


Number density of electrons at $t = 14$ ns

Afivo-streamer

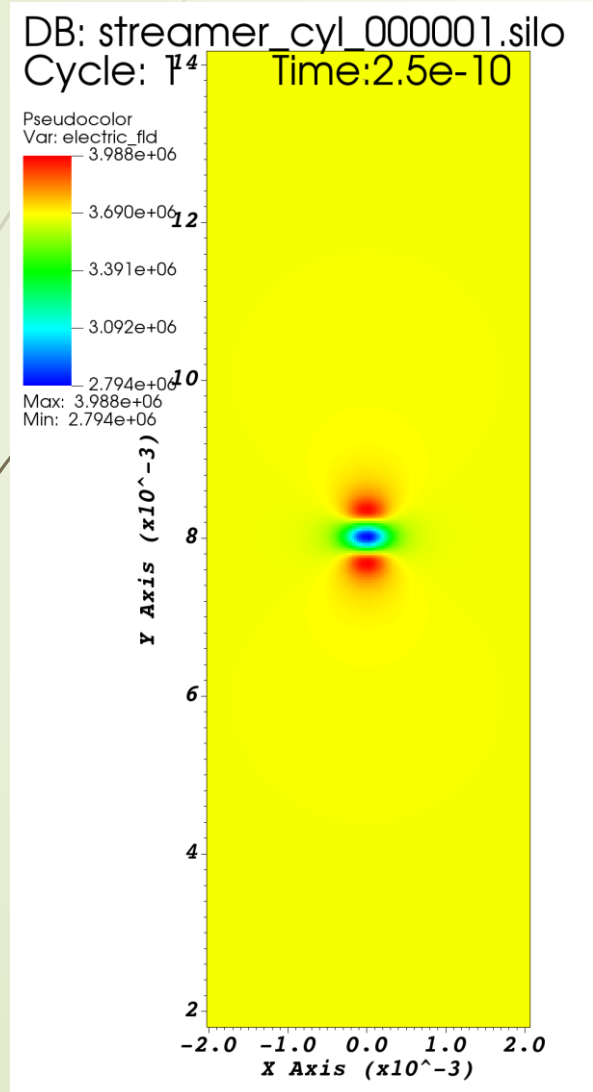


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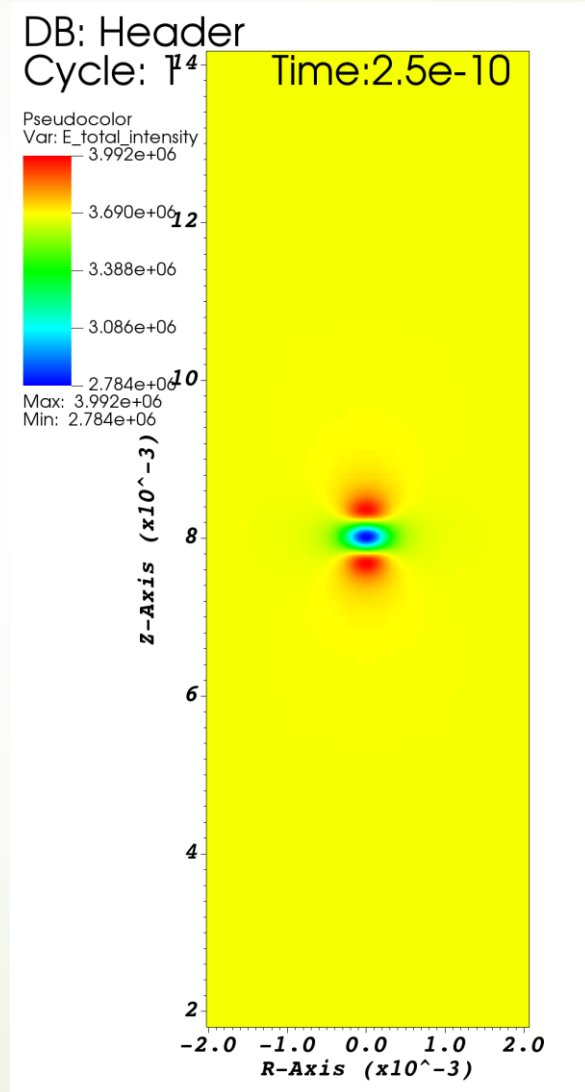


Electric field intensity at $t = 0.25$ ns

Afivo-streamer

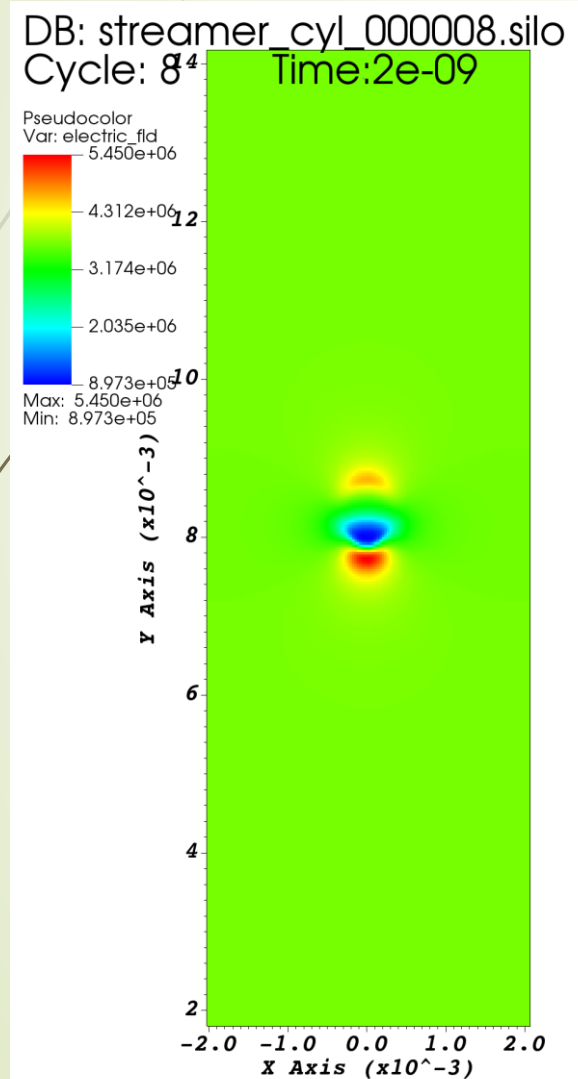


AMReX-streamer

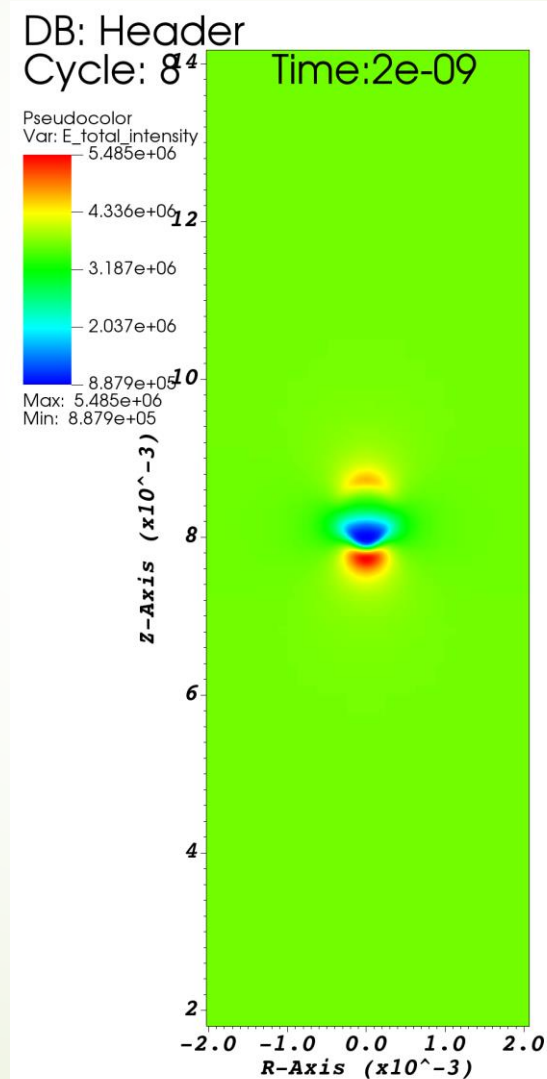


Electric field intensity at $t = 2 \text{ ns}$

Afivo-streamer

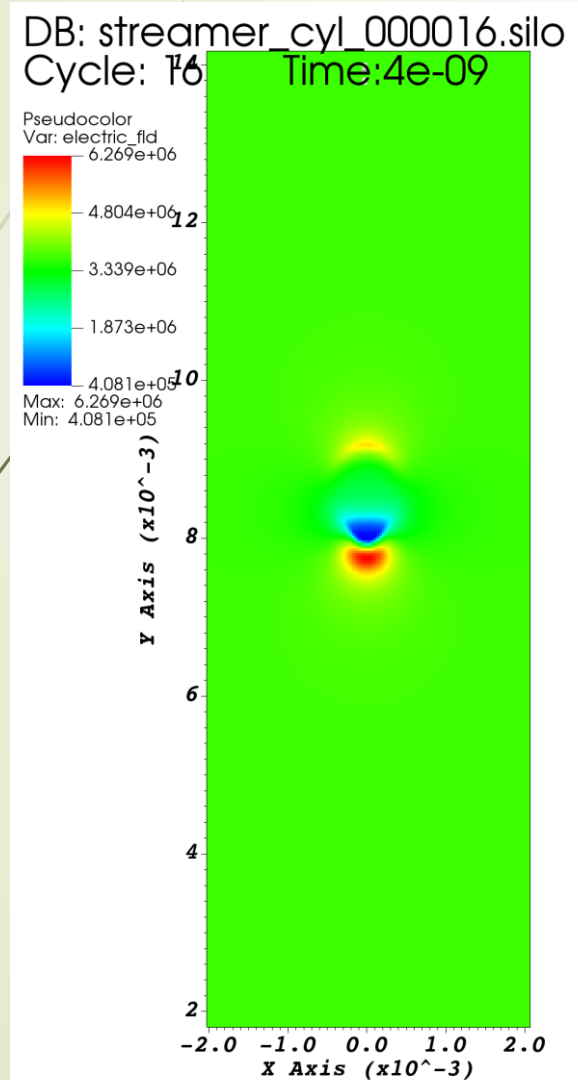


AMReX-streamer

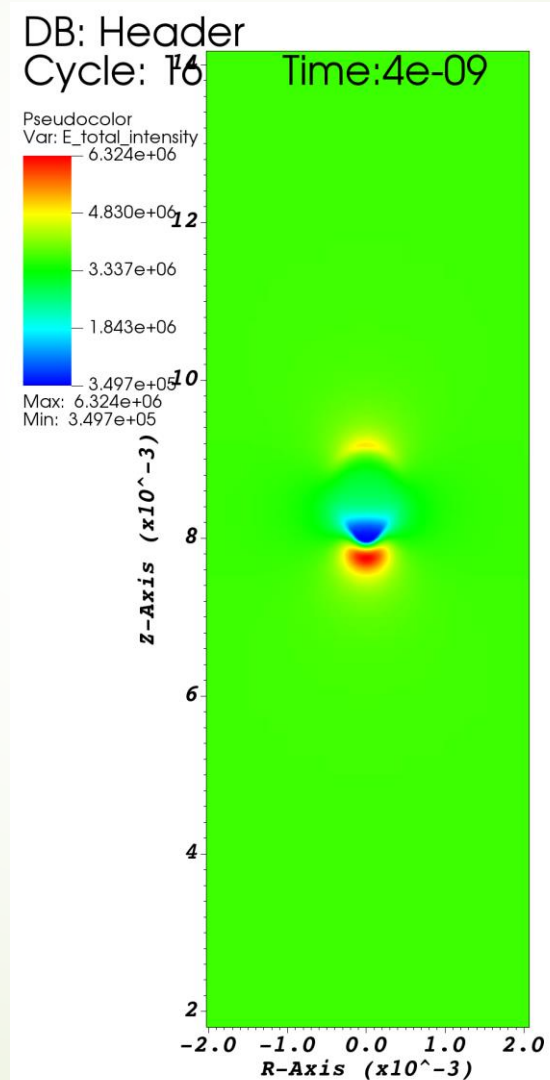


Electric field intensity at $t = 4 \text{ ns}$

Afivo-streamer

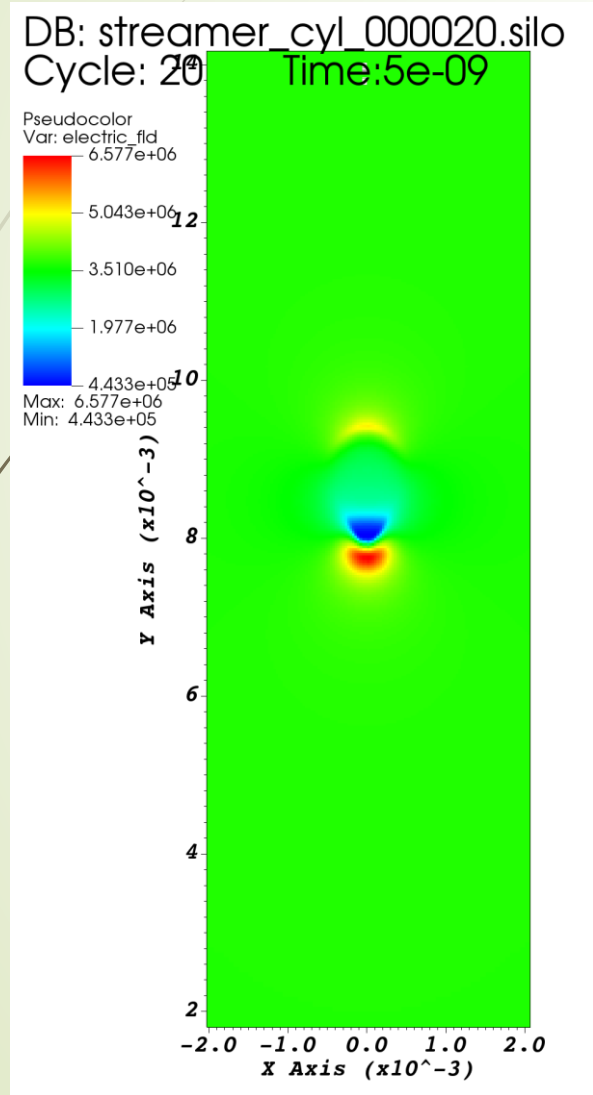


AMReX-streamer

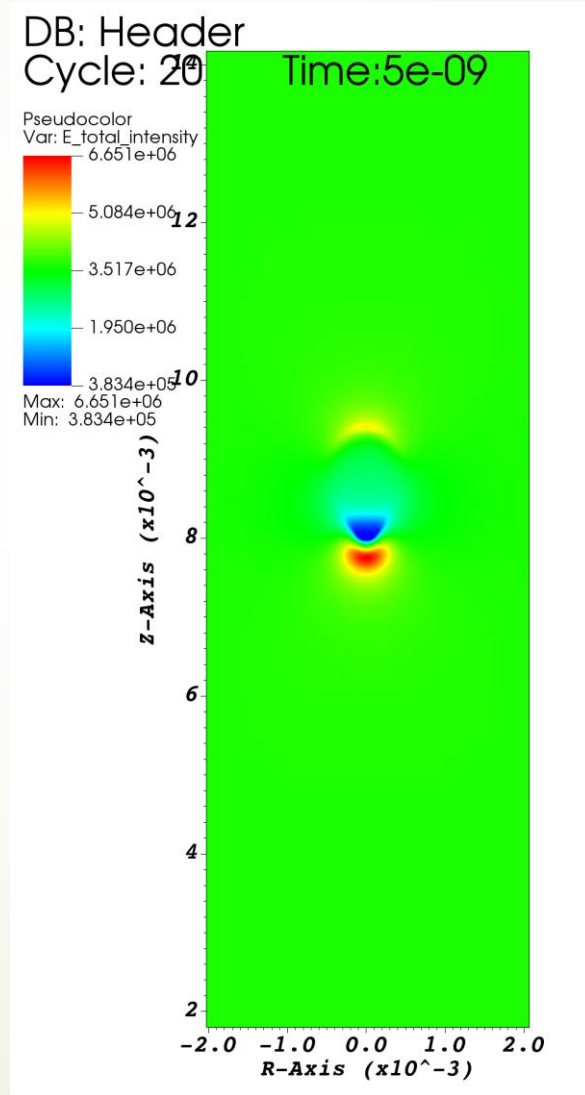


Electric field intensity at $t = 5 \text{ ns}$

Afivo-streamer

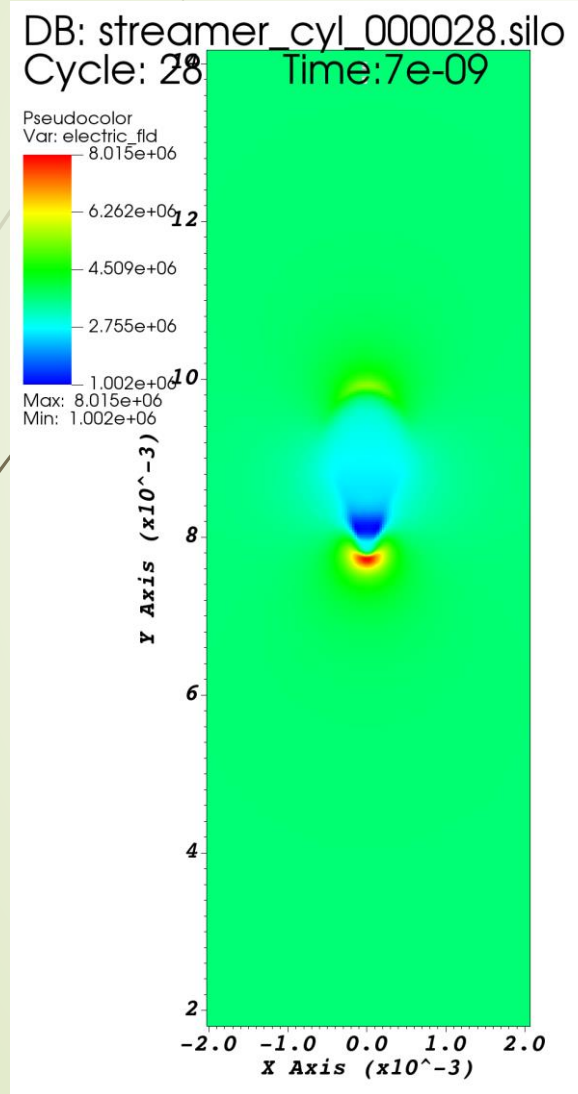


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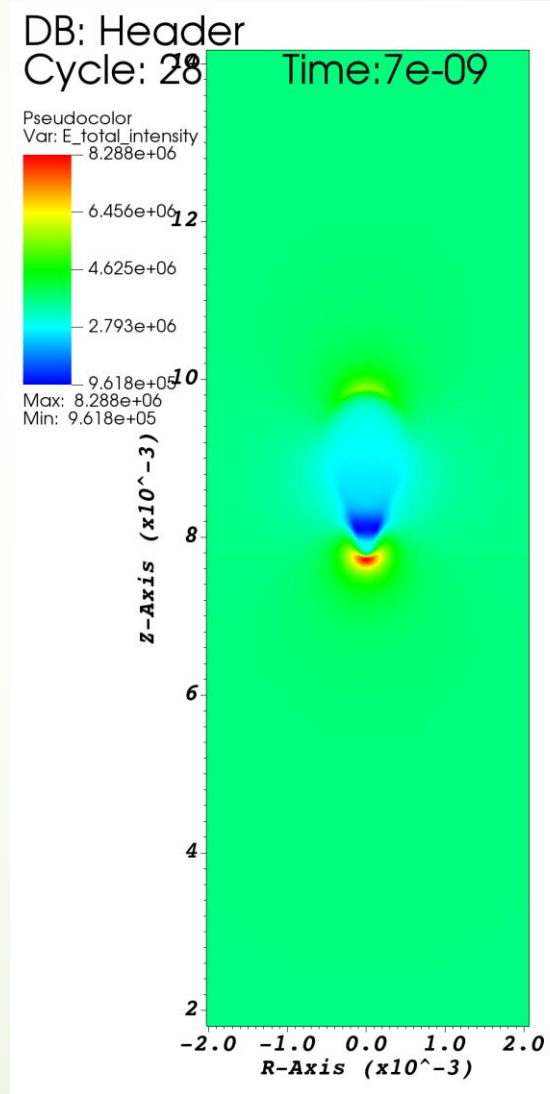


Electric field intensity at $t = 7$ ns

Afivo-streamer

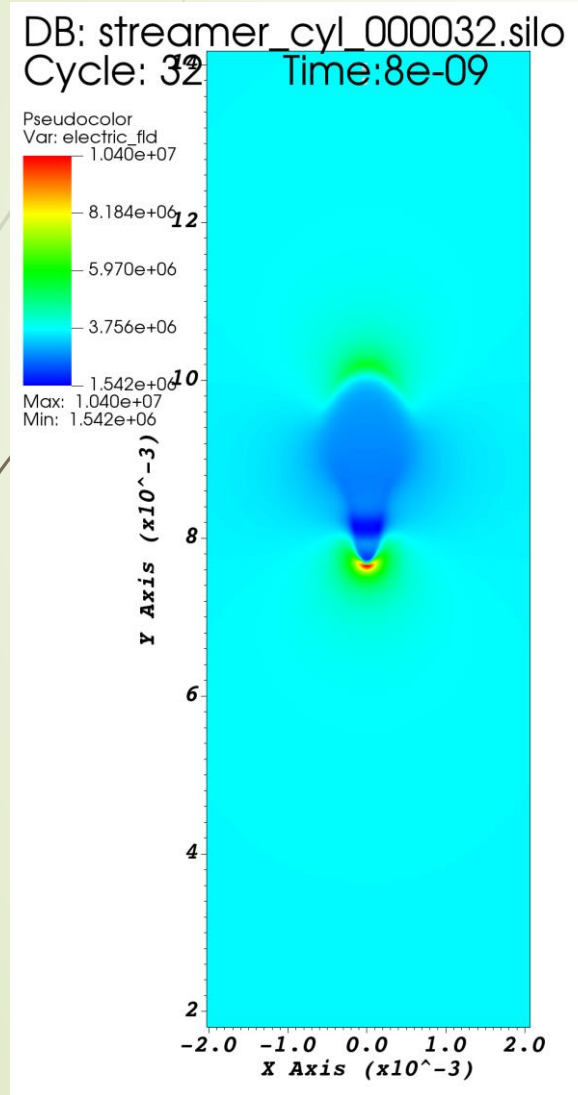


AMReX-streamer

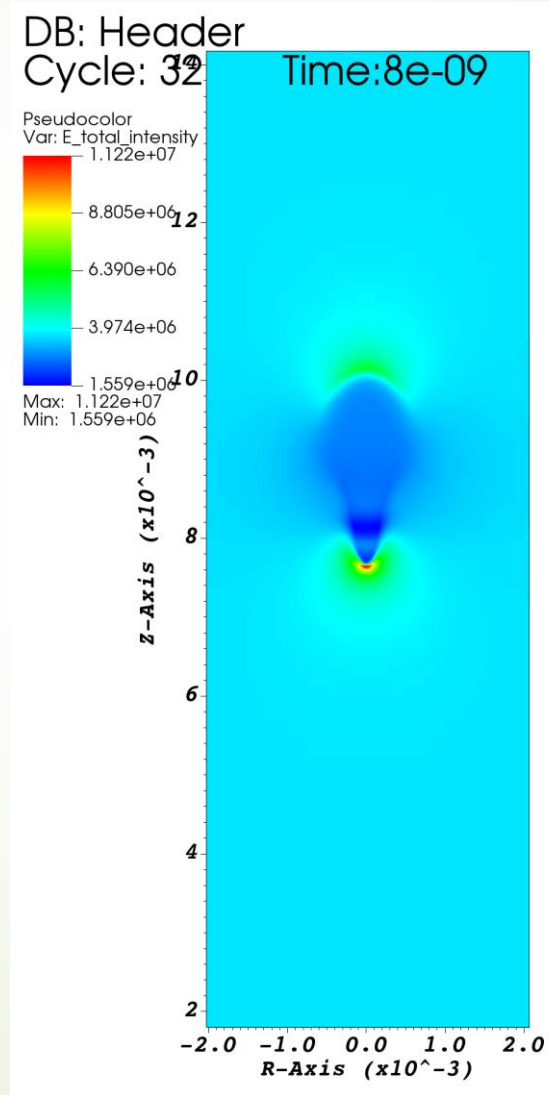


Electric field intensity at $t = 8 \text{ ns}$

Afivo-streamer

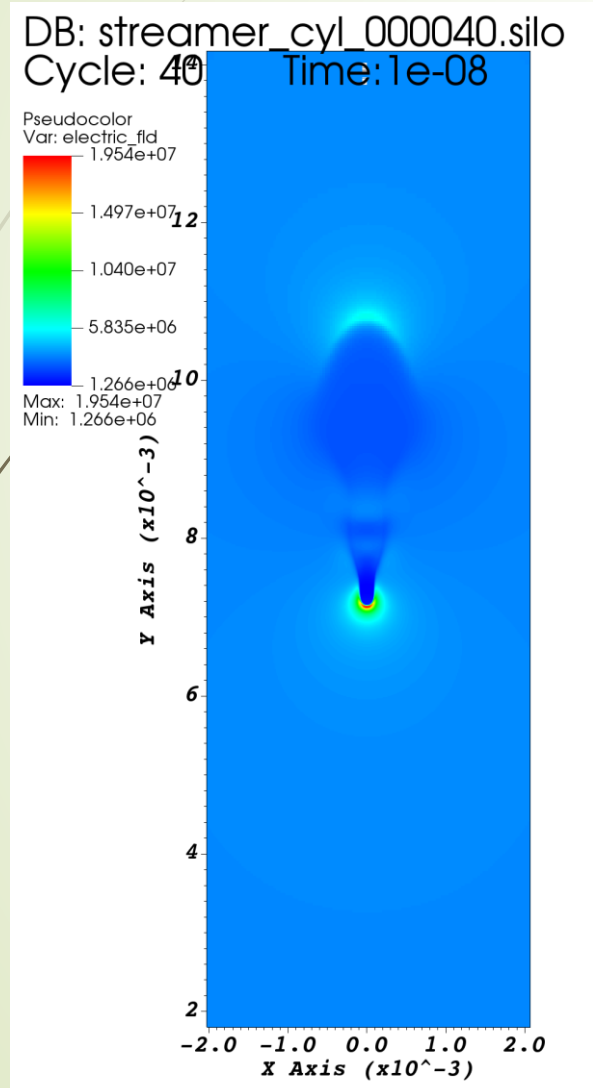


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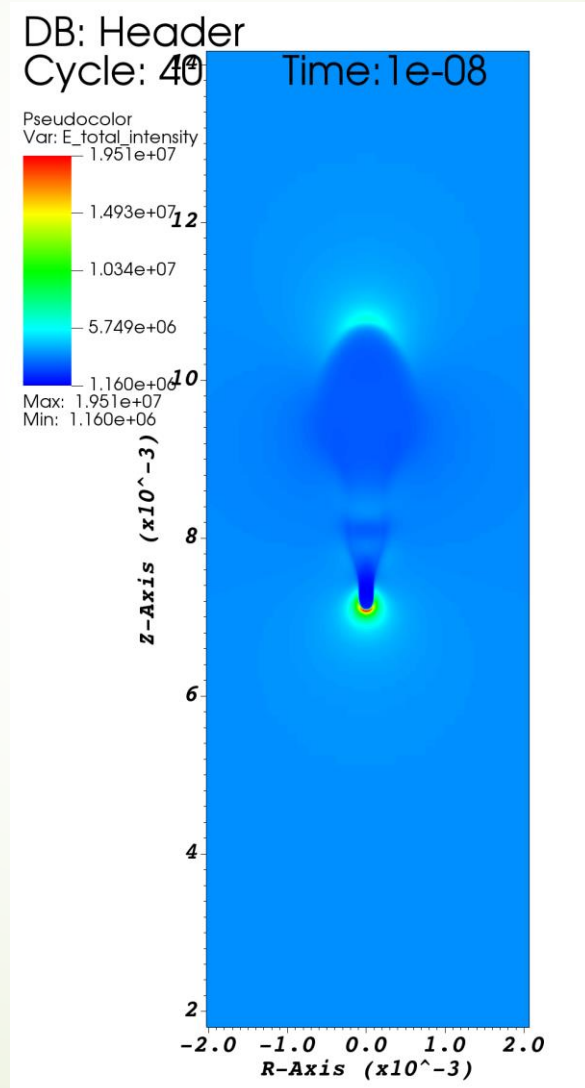


Electric field intensity at $t = 10$ ns

Afivo-streamer

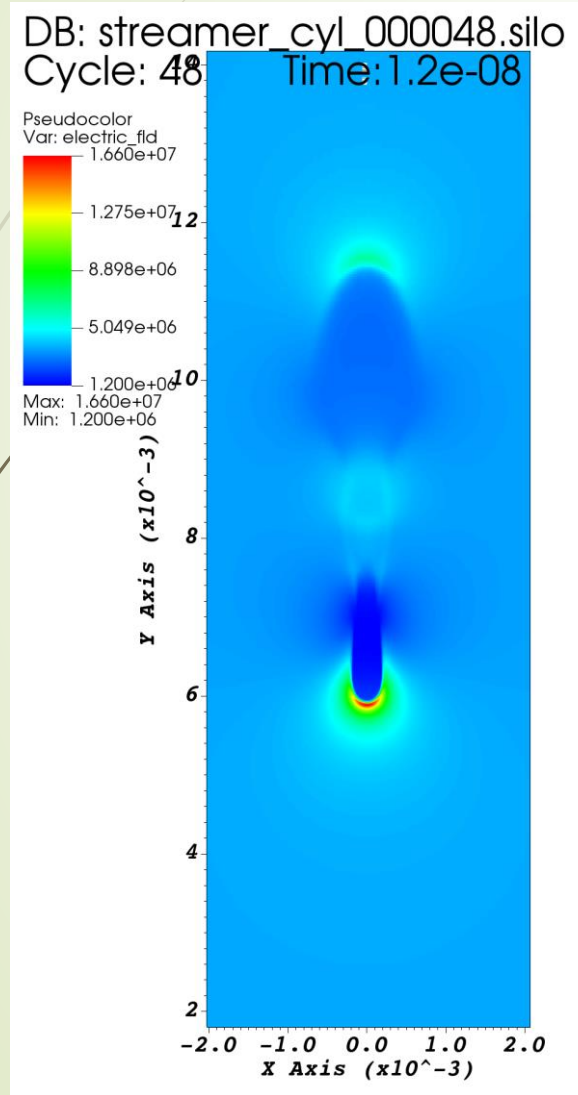


AMReX-streamer

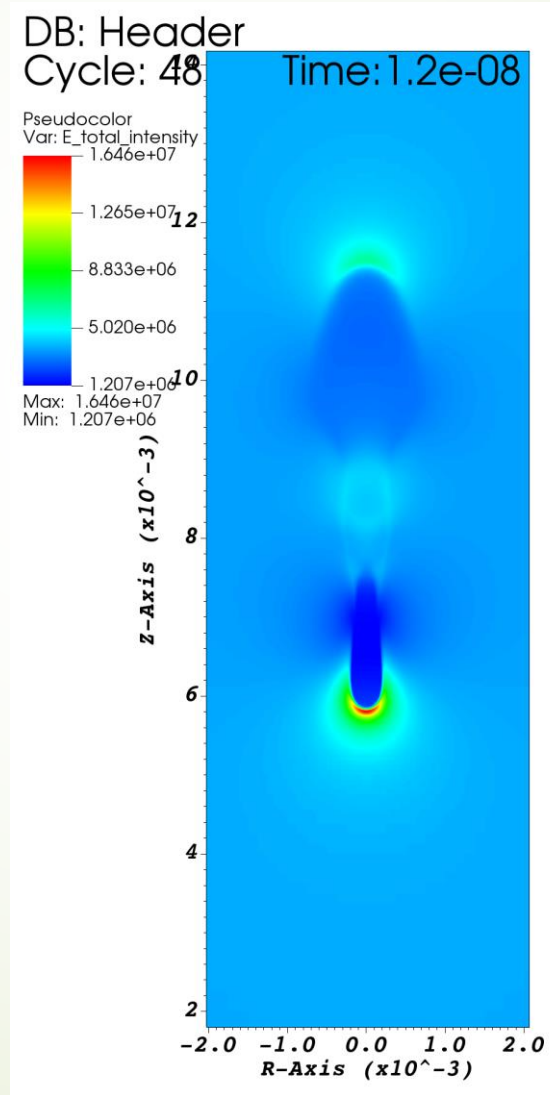


Electric field intensity at $t = 12 \text{ ns}$

Afivo-streamer

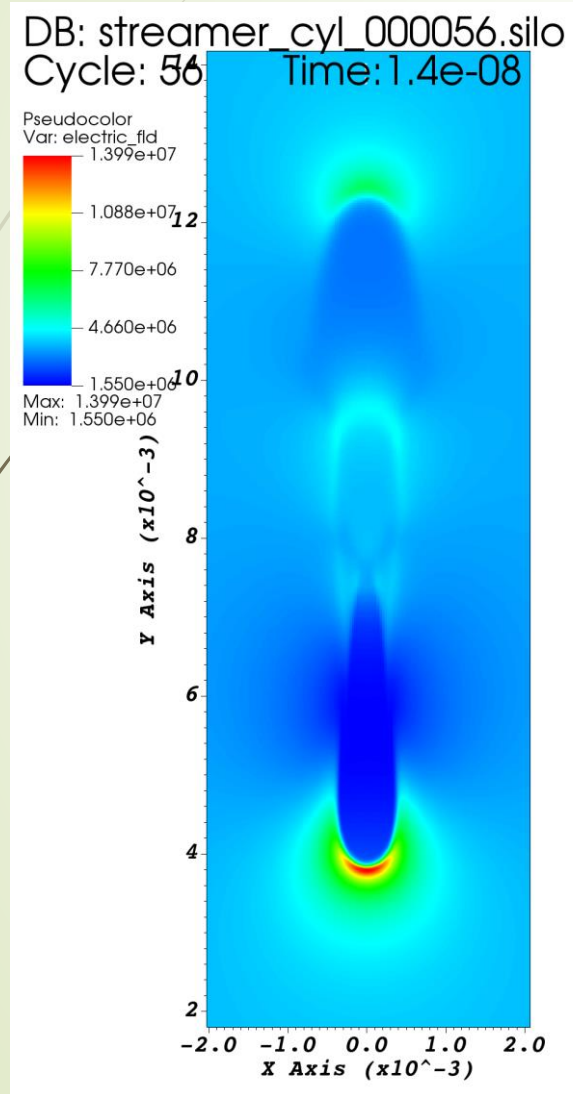


AMReX-streamer

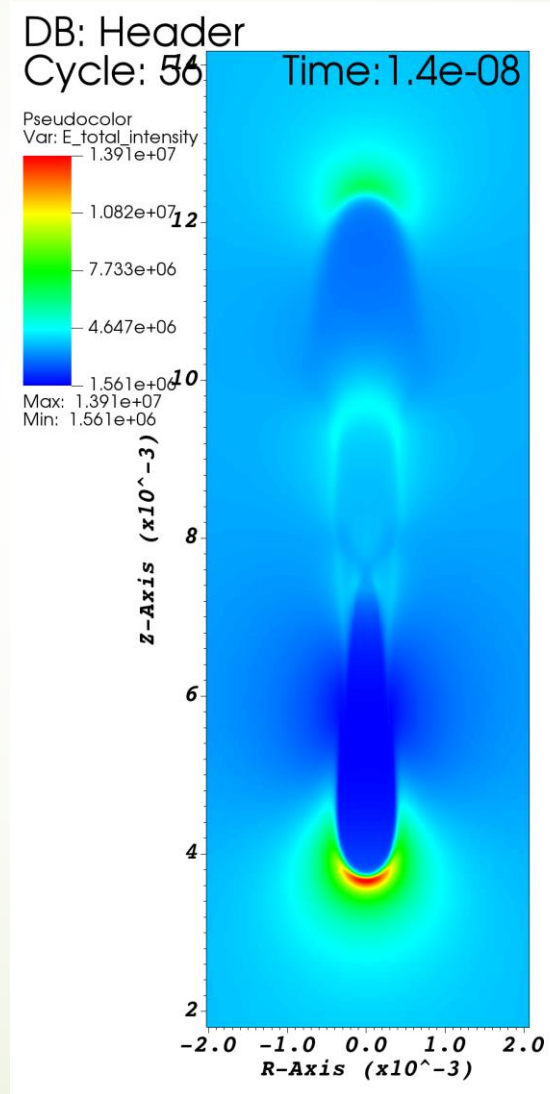


Electric field intensity at $t = 14$ ns

Afivo-streamer



AMReX-streamer





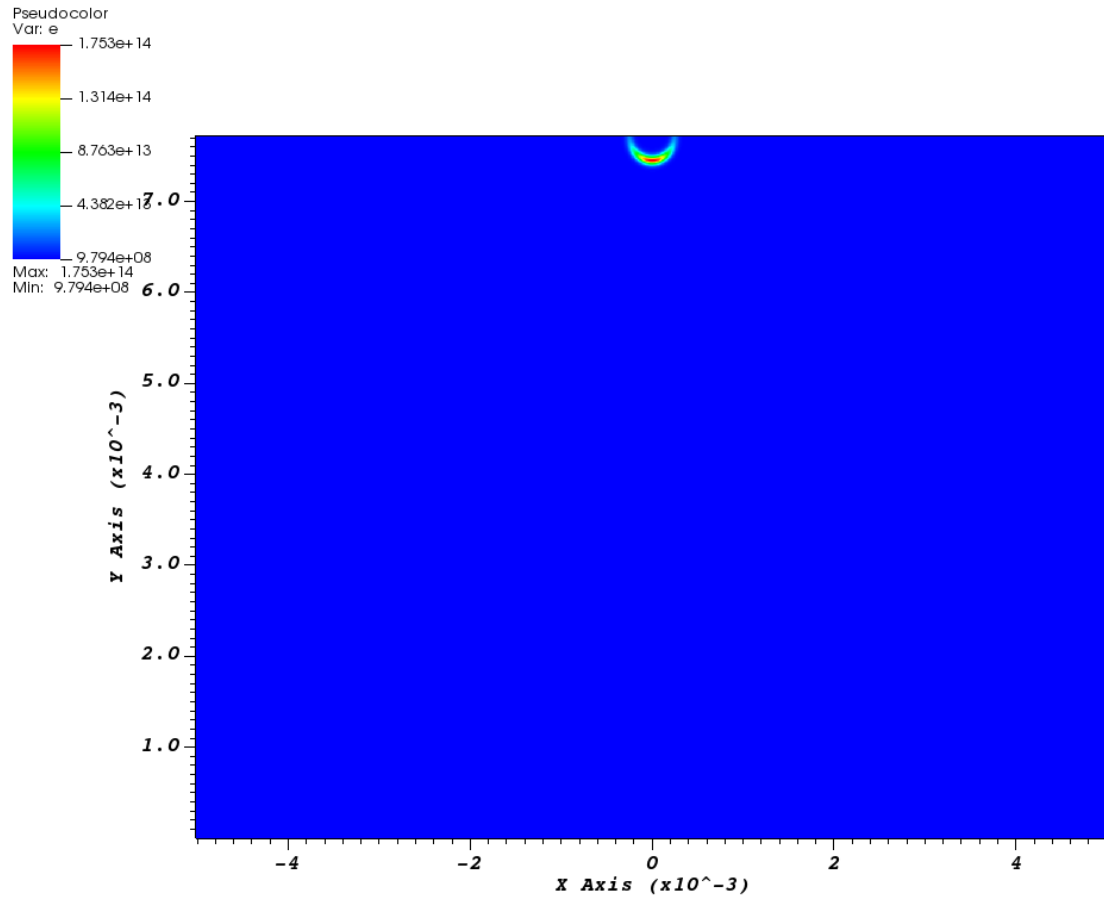
Positive streamer in the air

- The dry air: 80.0% N₂, 20.0% O₂ .
- The number of AMR levels 7 (from 0 to 6)
- Photoionization parametrization: Bourdon 3 term
- The applied reduced electric field: $E/N_0 = 80 \text{ Td}$
- The initial condition: Positive line segment with a length of 0.64 mm and a width of 0.4 mm, centered at the center of the z coordinate range
- Background ionization: 10^9 m^{-3} .
- Neutral gas density: $2.50475764 \cdot 10^{25} \text{ m}^{-3}$.

Number density of electrons at $t = 0.25$ ns

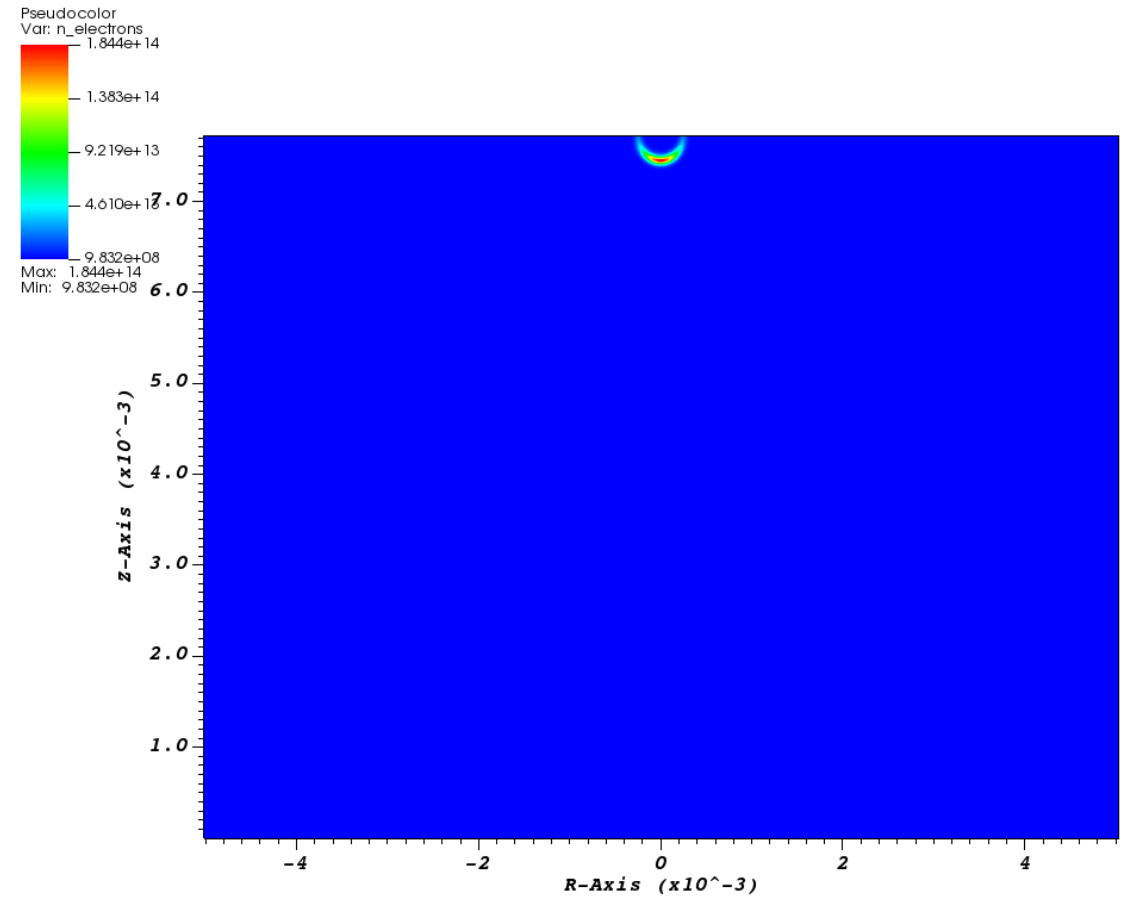
Afivo-streamer

DB: streamer_cyl_000001.silo
Cycle: 1 Time: 2.5e-10



AMReX-streamer

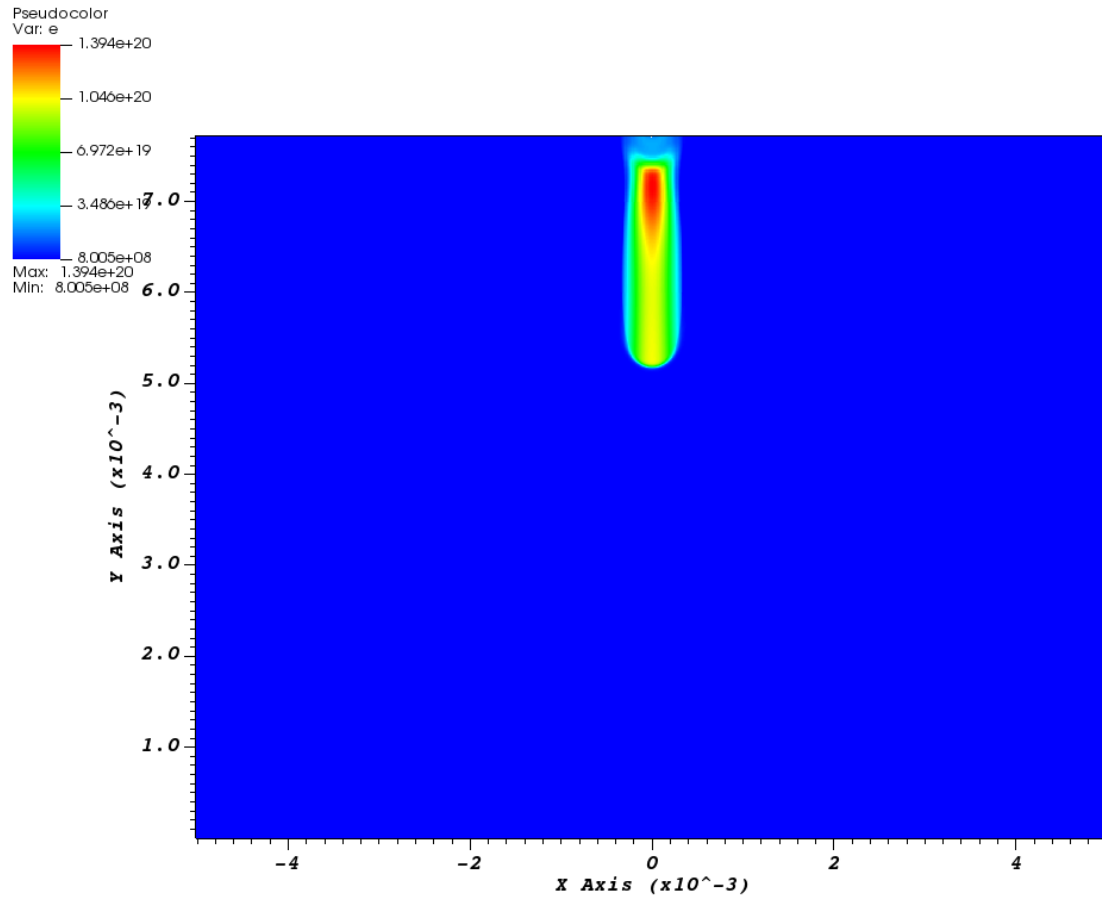
DB: Header
Cycle: 1 Time: 2.5e-10



Number density of electrons at $t = 4 \text{ ns}$

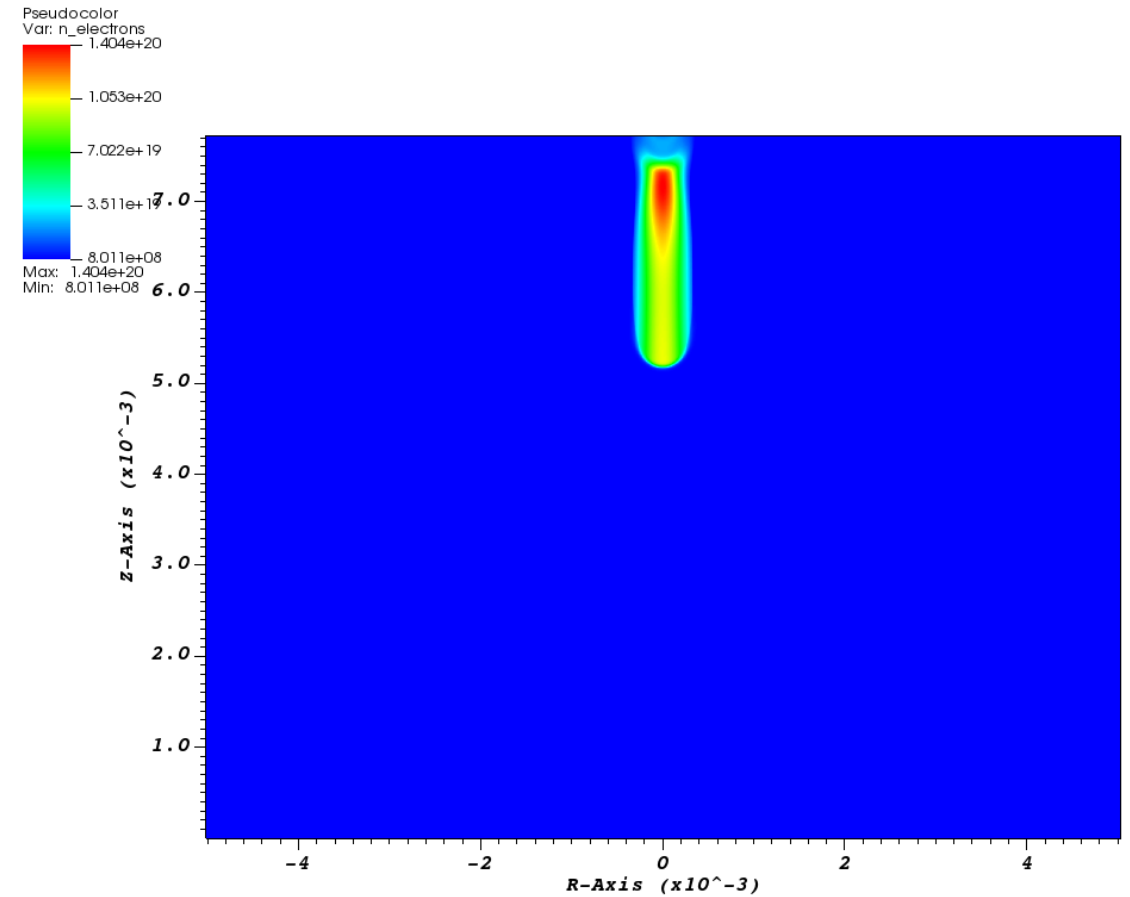
Afivo-streamer

DB: streamer_cyl_000016.silo
Cycle: 16 Time: 4e-09



AMReX-streamer

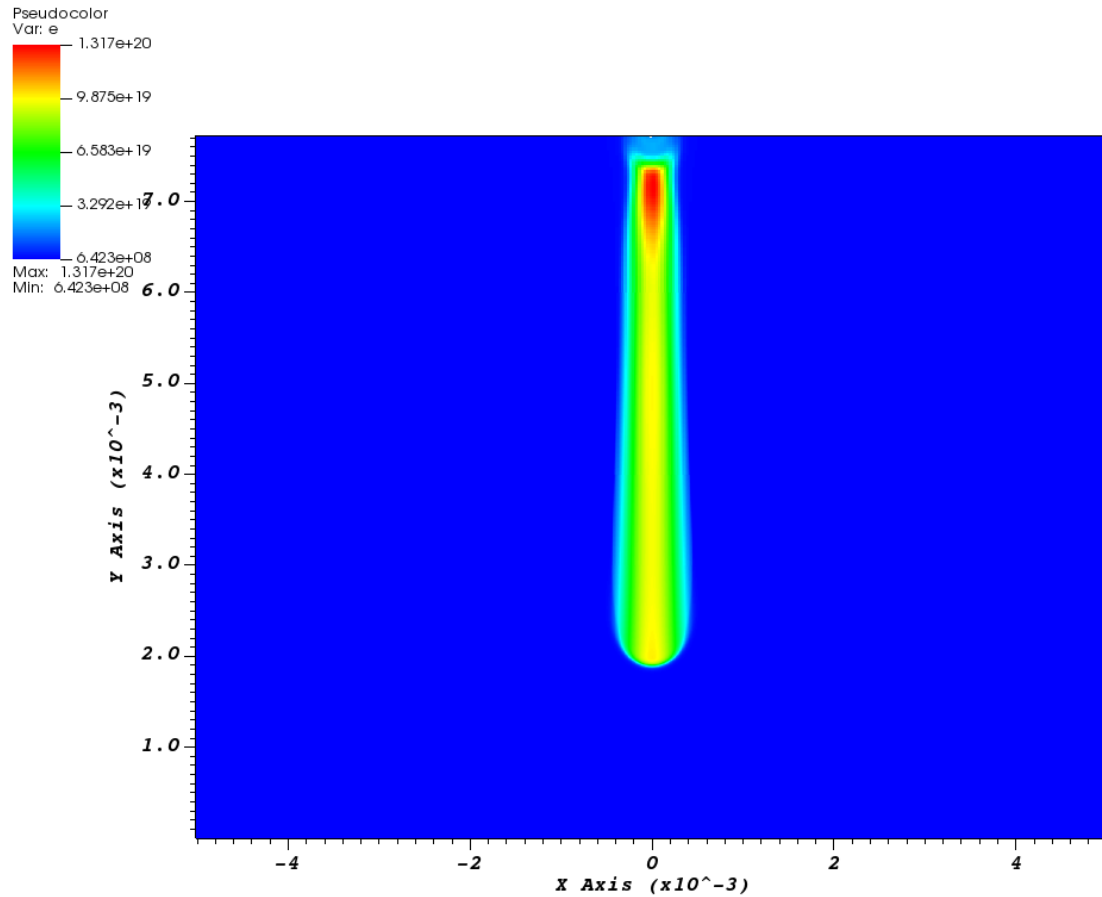
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Number density of electrons at $t = 8$ ns

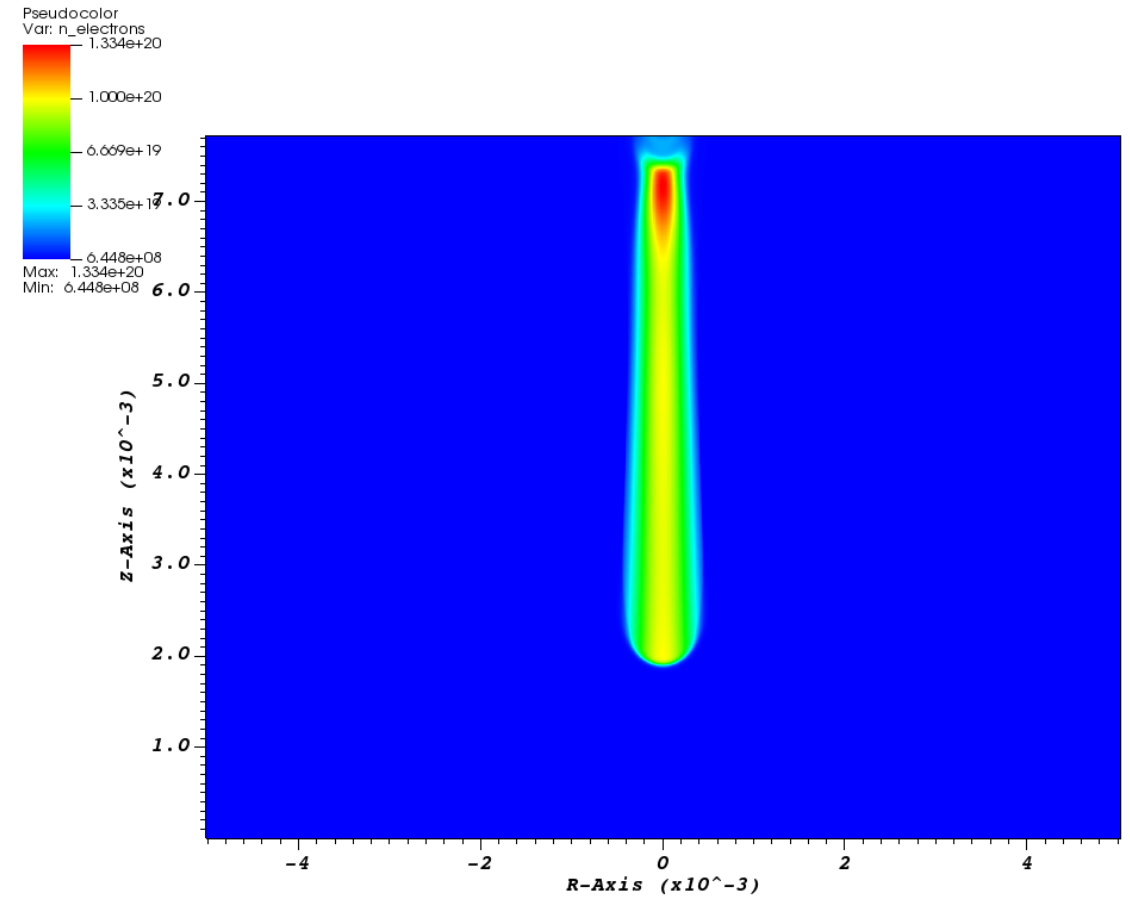
Afivo-streamer

DB: streamer_cyl_000032.silo
Cycle: 32 Time: 8e-09



AMReX-streamer

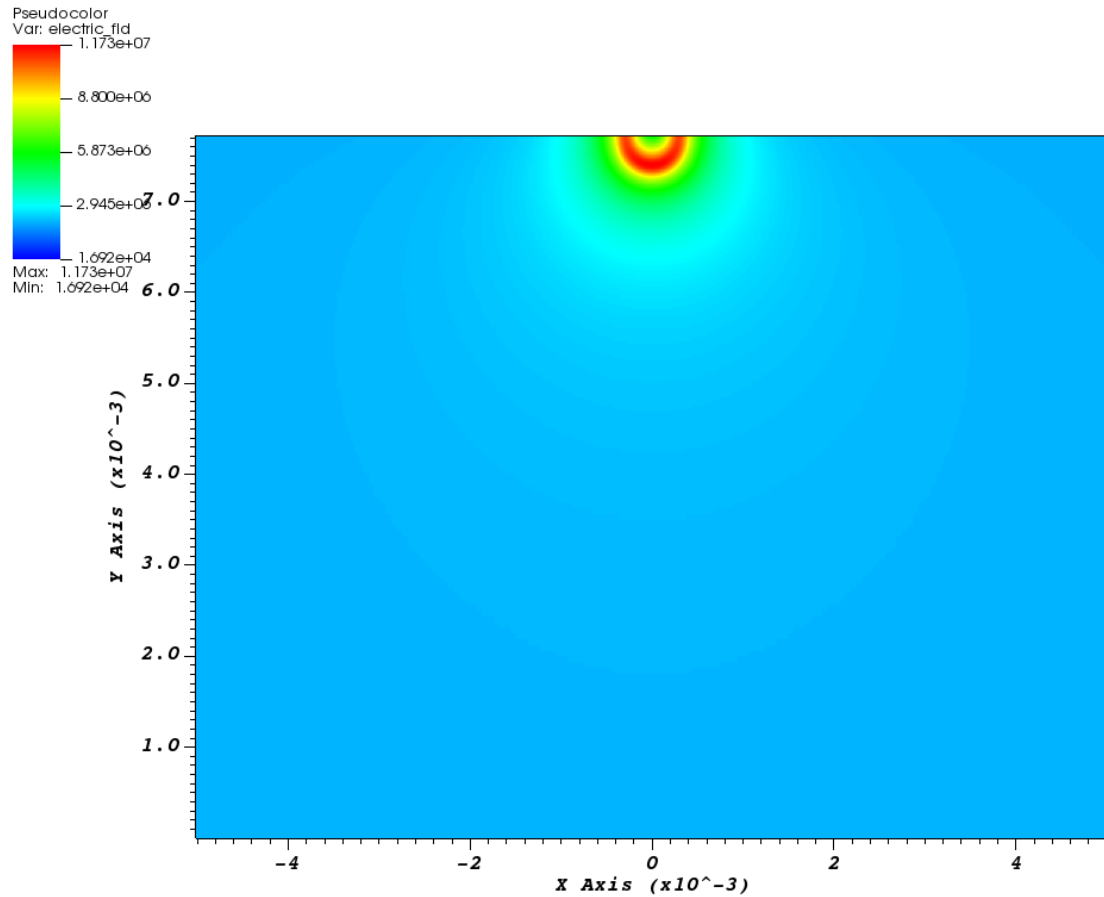
DB: Header
Cycle: 32 Time: 8e-09



Electric field intensity at $t = 0.25$ ns

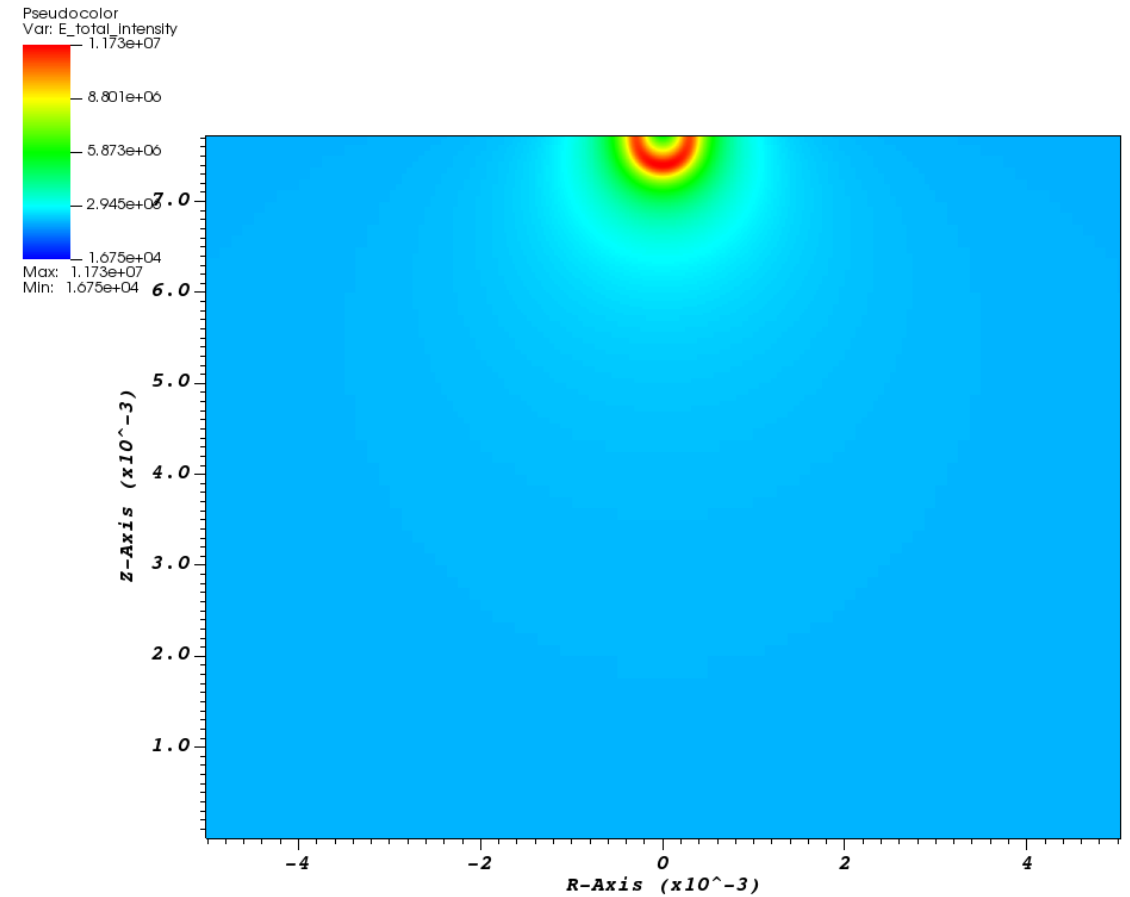
Afivo-streamer

DB: streamer_cyl_000001.silo
Cycle: 1 Time: 2.5e-10



AMReX-streamer

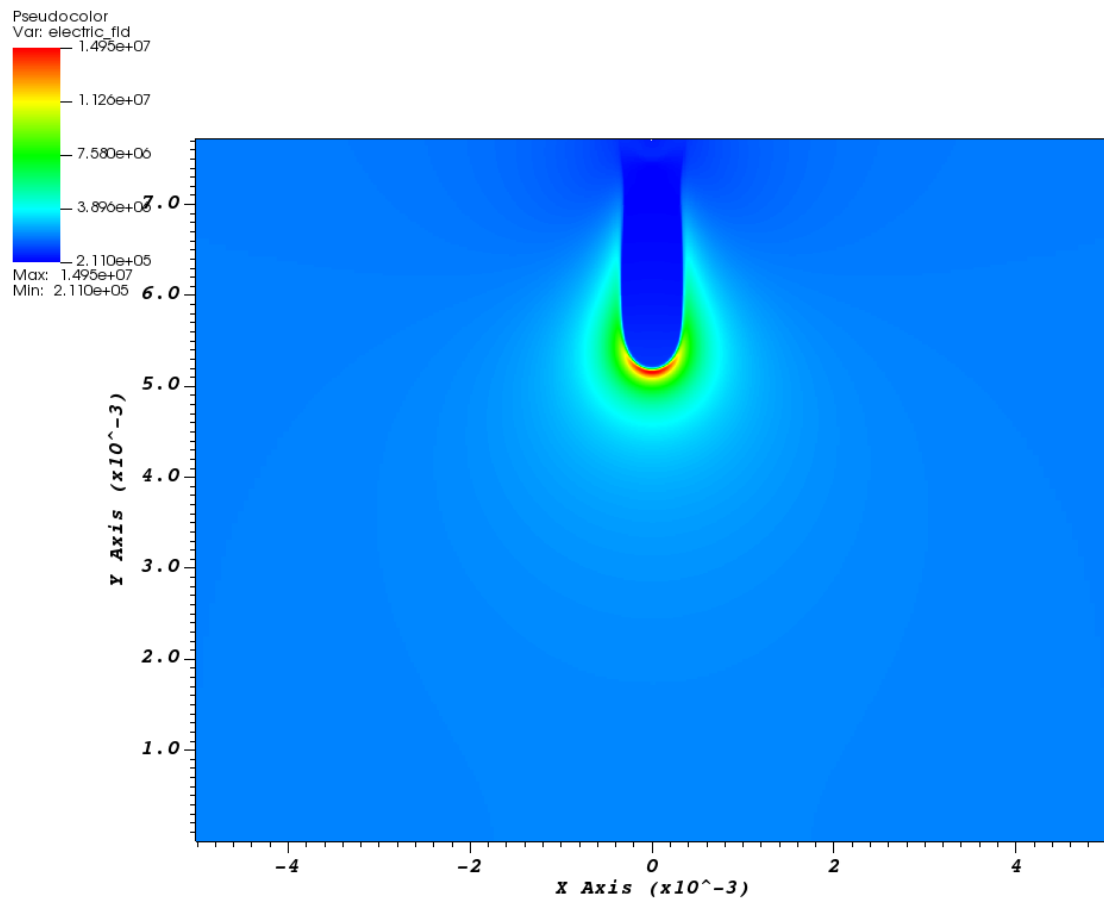
DB: Header
Cycle: 1 Time: 2.5e-10



Number density of electrons at $t = 4 \text{ ns}$

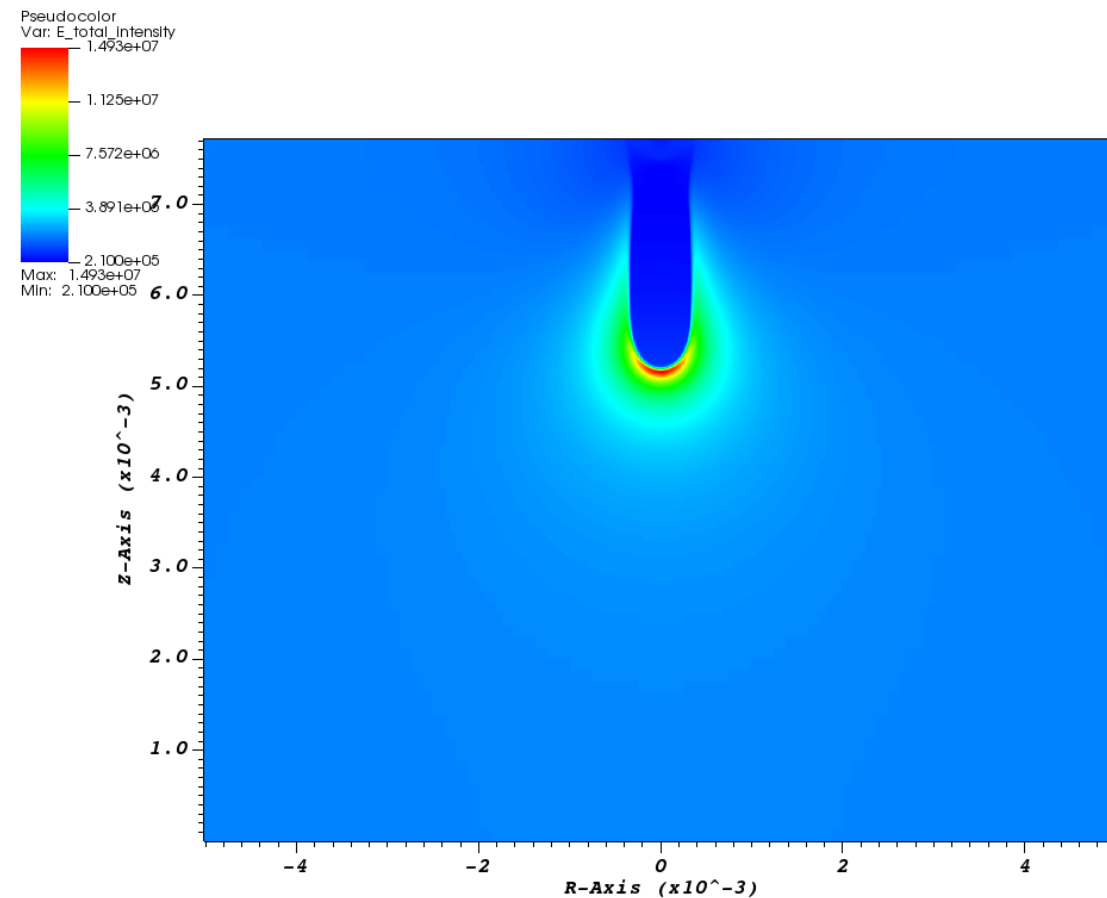
Afivo-streamer

DB: streamer_cyl_000016.silo
Cycle: 16 Time: 4e-09



AMReX-streamer

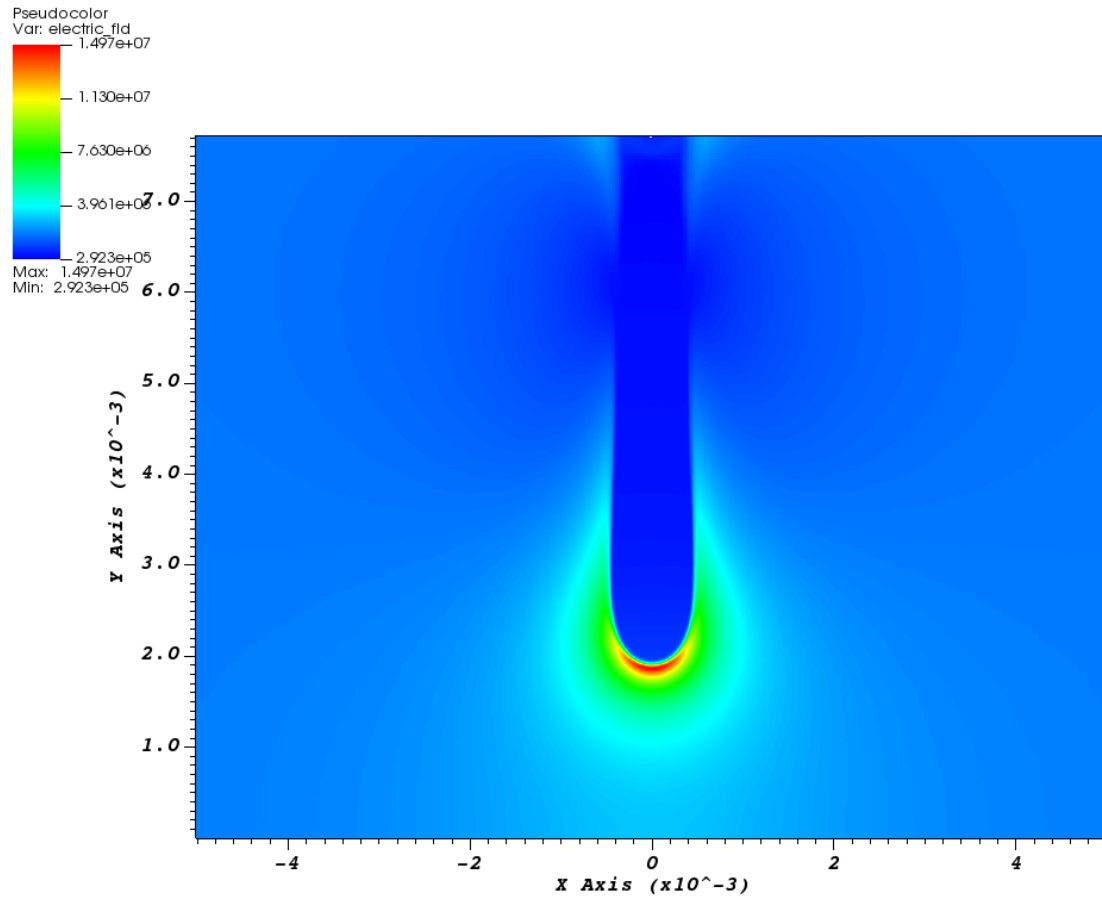
DB: Header
Cycle: 16 Time: 4e-09



Number density of electrons at $t = 8$ ns

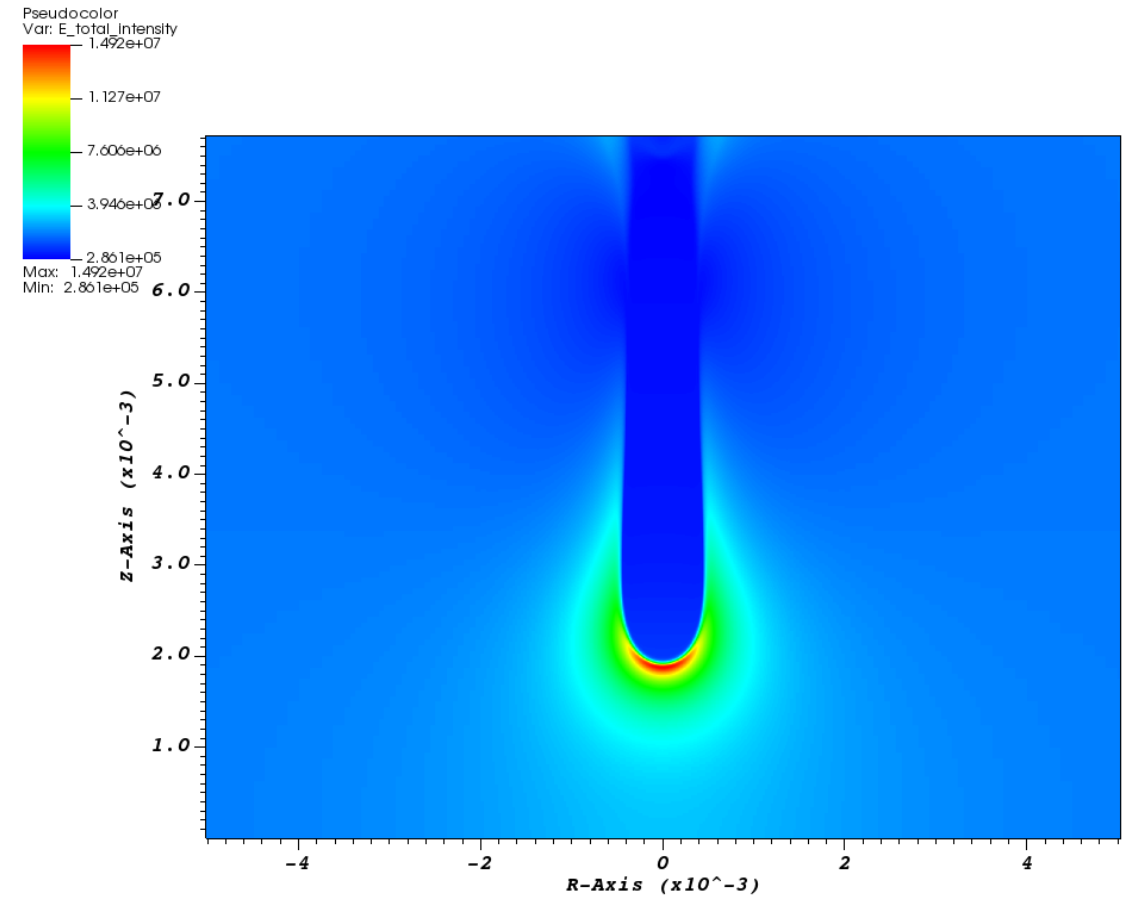
Afivo-streamer

DB: streamer_cyl_000032.silo
Cycle: 32 Time: 8e-09



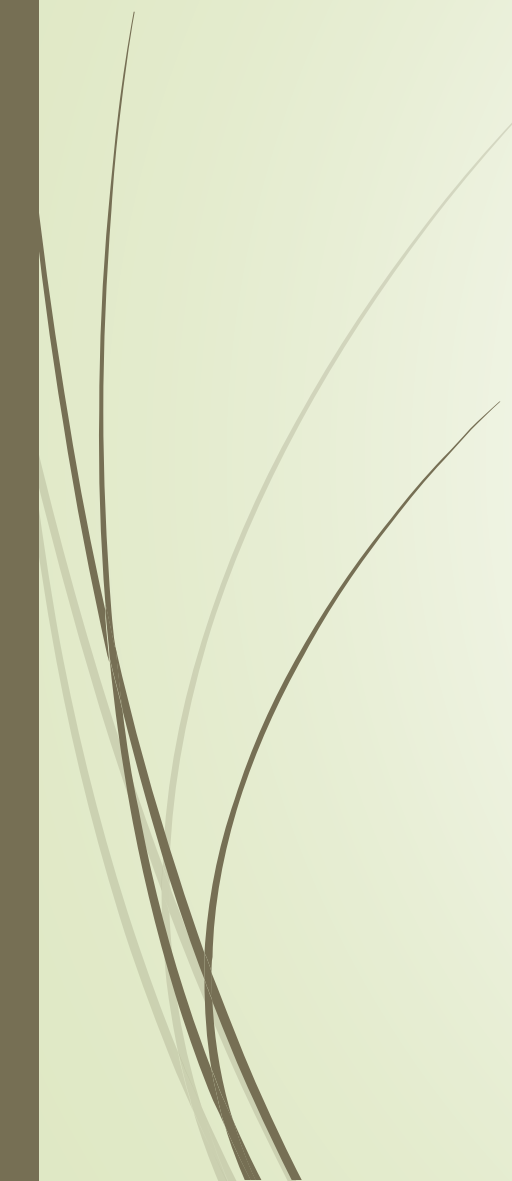
AMReX-streamer

DB: Header
Cycle: 32 Time: 8e-09





Conclusions

- We have implemented an axisymmetric fluid model in the AMReX library
 - Our model is based on the first-order fluid model with local field approximation
 - Photoionization is implemented by solving a system of Helmholtz equations
 - The good agreement between the results of our program and the Afivo-streamer open-source program confirms the validity of our code.
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➤ Acknowledgements

- This research was supported by the Science Fund of the Republic of Serbia, Grant No. 7749560, project EGWIn.
- The authors acknowledge support from the MESTD of Serbia.

THANK YOU

