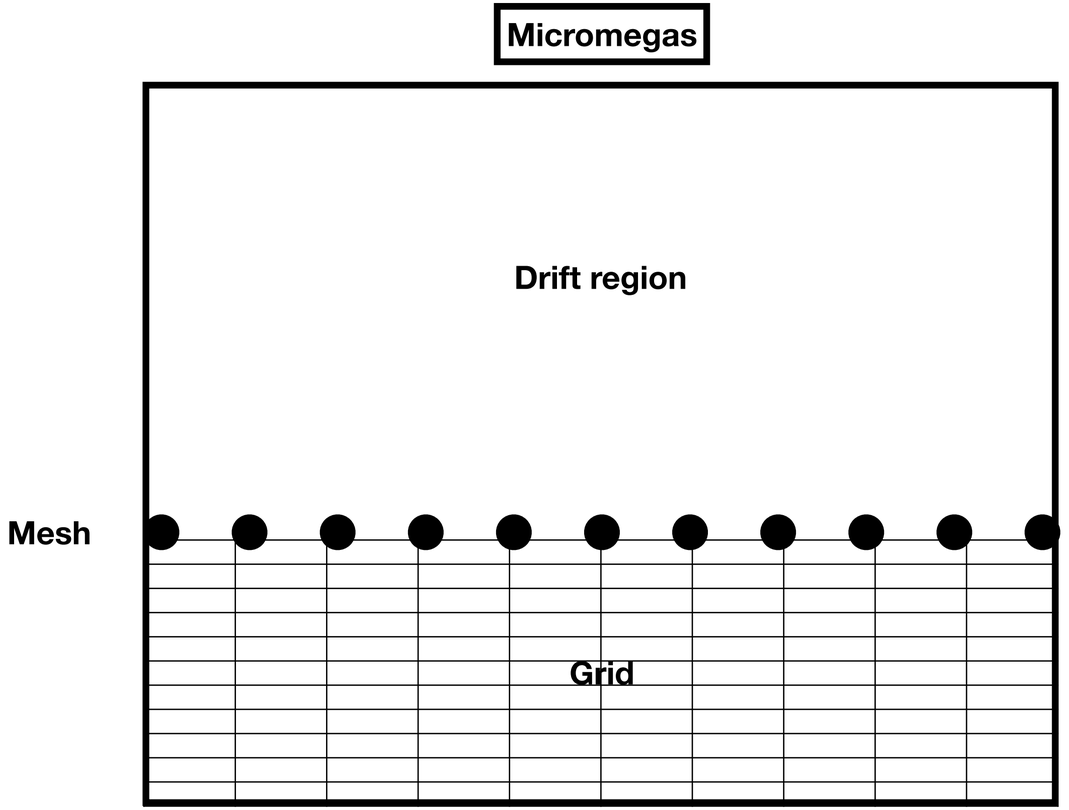
Space charge effects in simulations of large avalanche dynamics

Garfield++

Tom Szwarcer | University of Oxford | DRD1

The aim

- Use space charge effect in microscopic/MC avalanche simulations
- Allow the user to define regions where the effect should be included
- Include both electrons and ions
- No mirror charges in a general geometry - so use free space field as a first approximation



The grid

- Snap particles to a 2D (z, r) grid
- Compute the field due to axisymmetric rings of charge at a given (z, r)
- This prevents divergences that would be present in point charge fields.

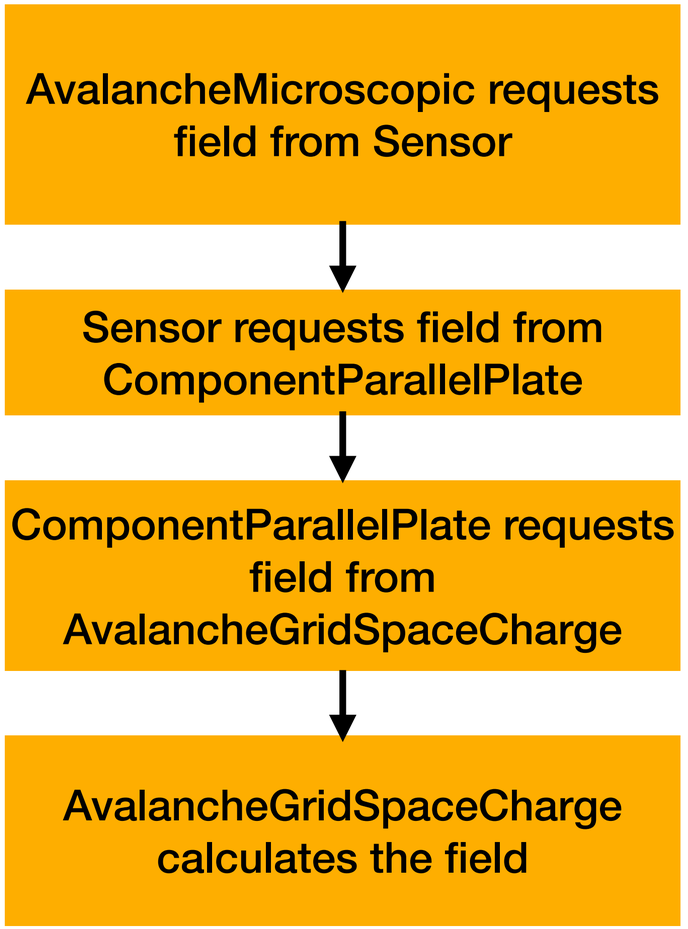
Making use of the RPC code

- We set up a ‘fake’ parallel plate system that exists at grid limits
- We set dV=0 so the only field present is the space charge field
- Tell Sensor that ComponentParallelPlate has its own field
- Sensor automatically adds the contribution to the field from all components
- This lets us reuse existing code for the space charge effect in RPCs

```
ComponentParallelPlate cmp;
if (enableSpaceCharge) cmp.EnableSpaceCharge();
sensor.AddComponent(&cmp);
```

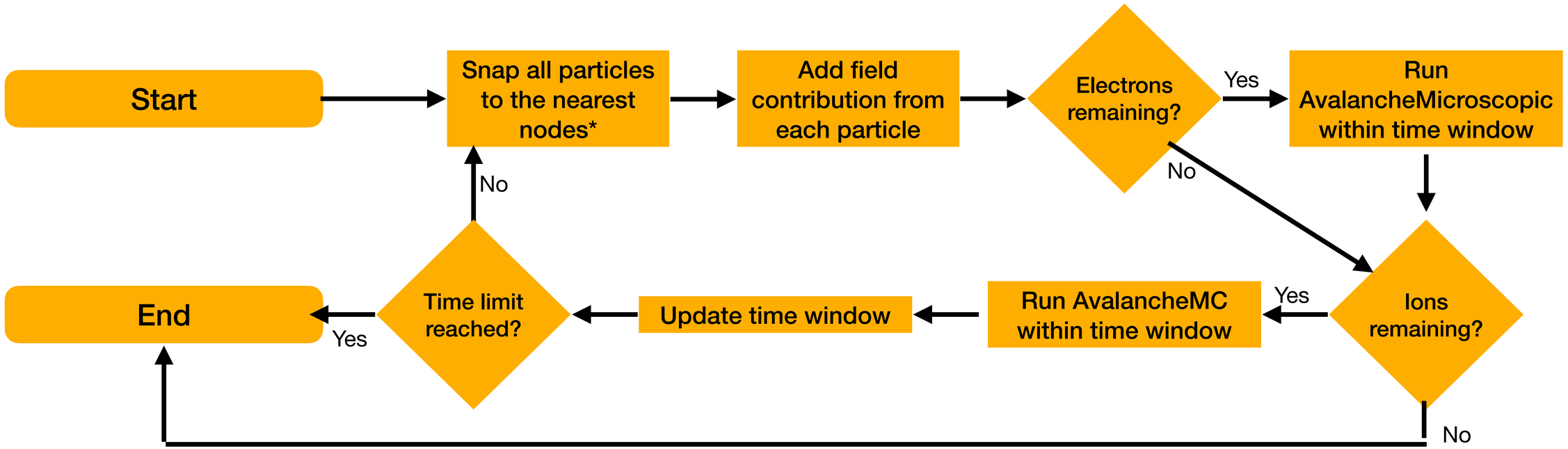
Communicating the field

- In order to get the field into sensor we have to go through ComponentParallelPlate
- This allows a lot of code to be reused



Electron & ion simulation

- As far as I know, it is not possible to simultaneously run AvalancheMC and AvalancheMicroscopic
- I do a leapfrog approach where ions and electrons are simulated alternately.



* the particles do not move in physical space - only on the grid, in order to calculate the field

Updating the field during simulation

- New charged particles may be produced during ionisations, penning transfer, or attachment
- These occur in real time during the simulation
- However the field is only updated once per timestep…
- How do we update the field mid-timestep?

User handles

- Functions that Garfield++ calls every time an ‘event’ happens.
- In practice the total charge at a given location doesn’t change in these events
- As new particles are produced at the same location, they will be snapped to the same grid point
- Therefore the field won’t change, and doesn’t need to be updated

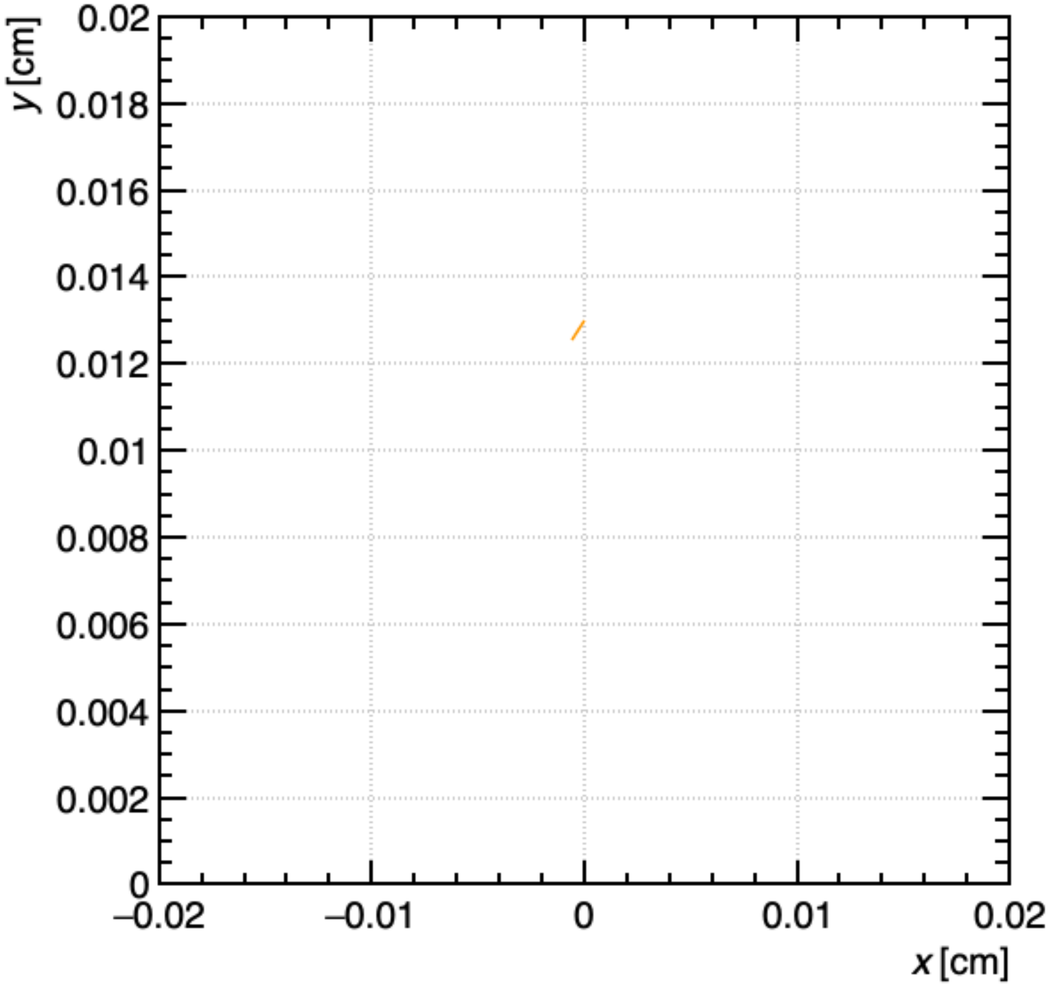
- However user handles provide a convenient way to pass on any new ions created to AvalancheMC
- New electrons are already accounted for in AvalancheMicroscopic

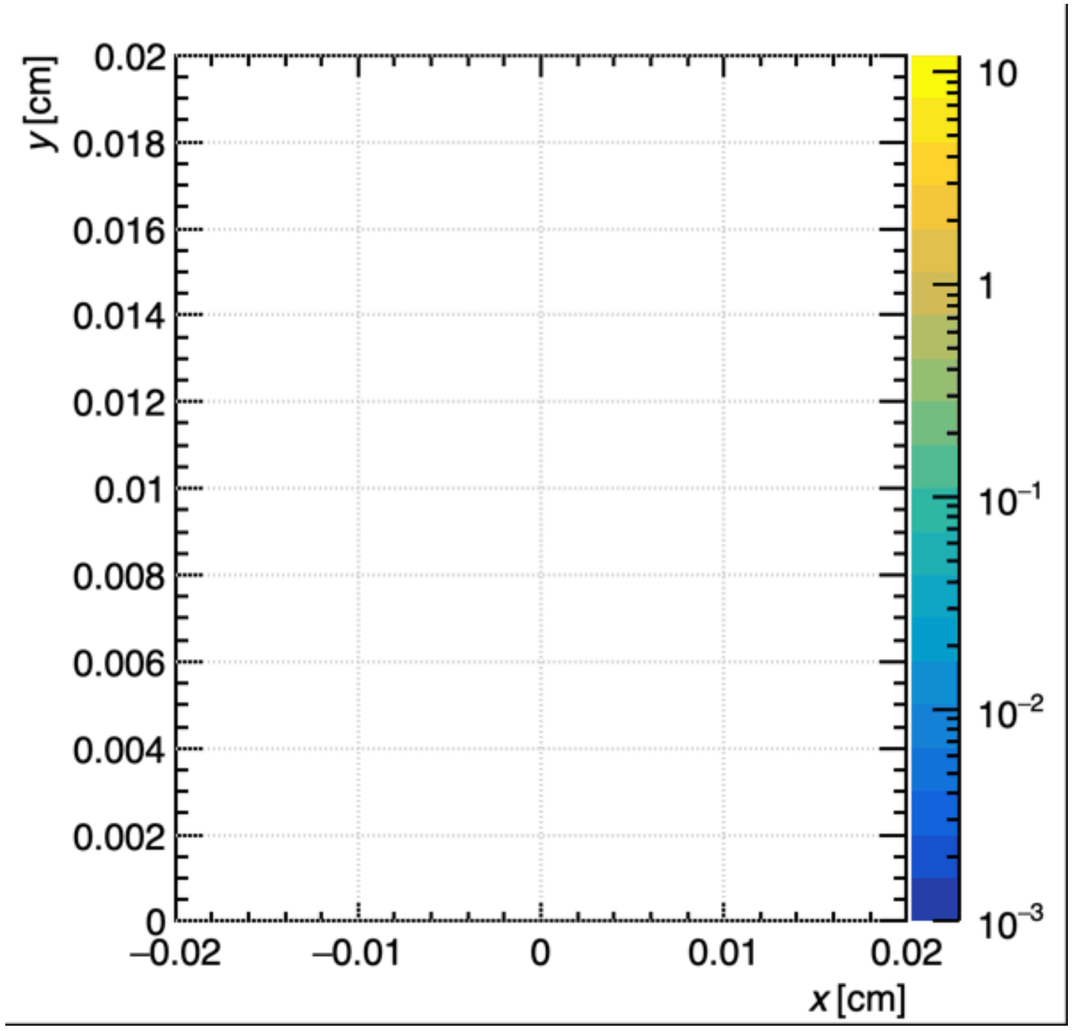
Dynamic mean position

- Avalanches will not always be centred on zero
- We need a way to centre the field on the mean position of the particles in the avalanche
- If there are particles in the avalanche, find the mean position and set as the zero of the radial coordinate

Visualising the field

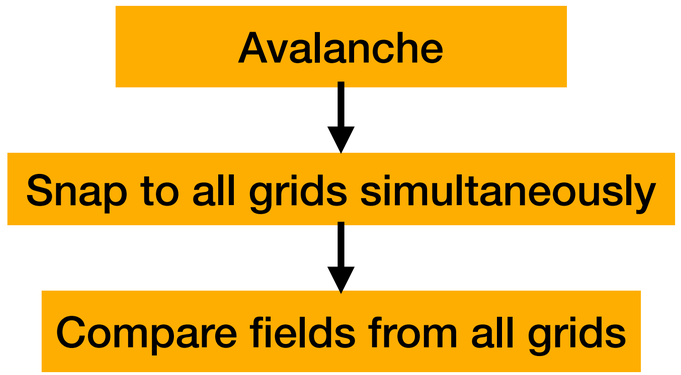
GIFs viewable at https://github.com/tomszwarcer/spacecharge-sim





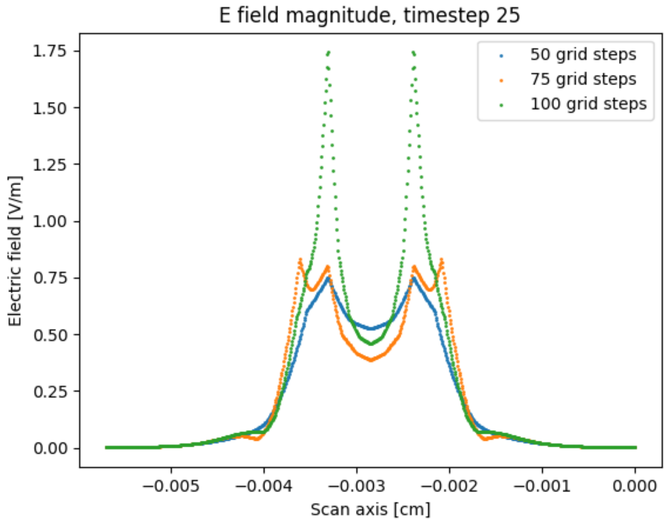
Investigating optimal grid spacing

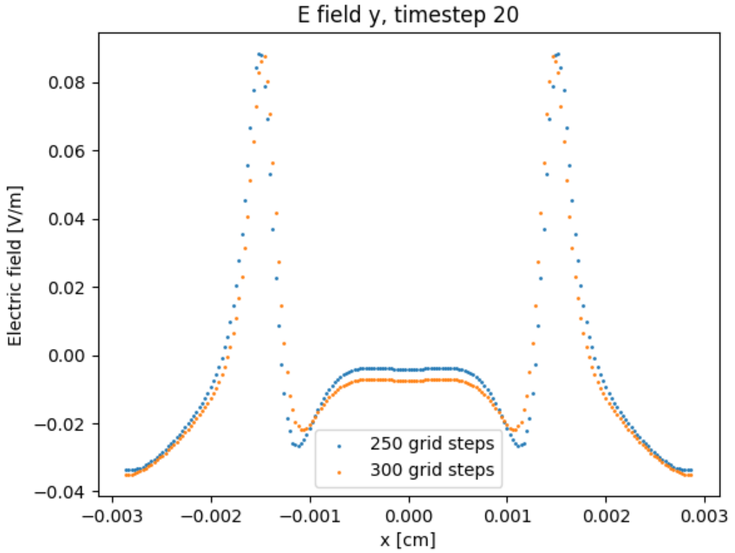
- Simulate an avalanche with multiple grids simultaneously
- Decide at what point it has converged ‘enough’ to justify a shorter computation time

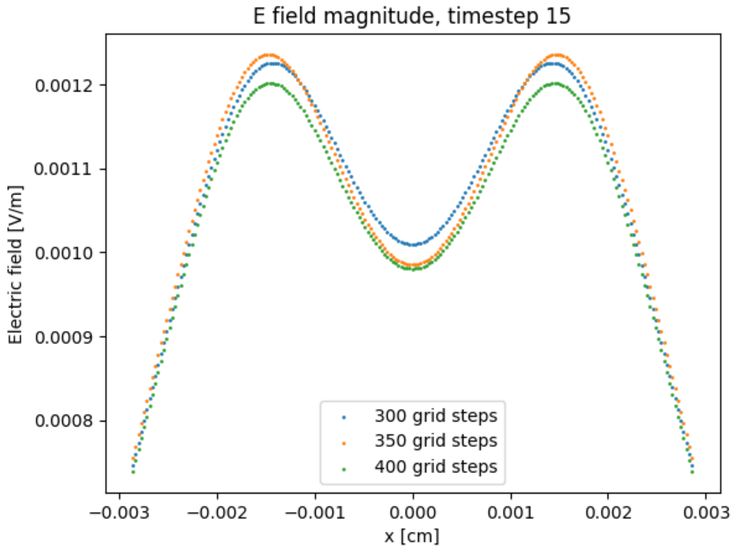


Investigating optimal grid spacing

- At low grid spacings, the field is a strong function of the grid spacing
- This is because electrons are being snapped a non-negligible proportion of the region of interest
- Above about 300, they start to converge





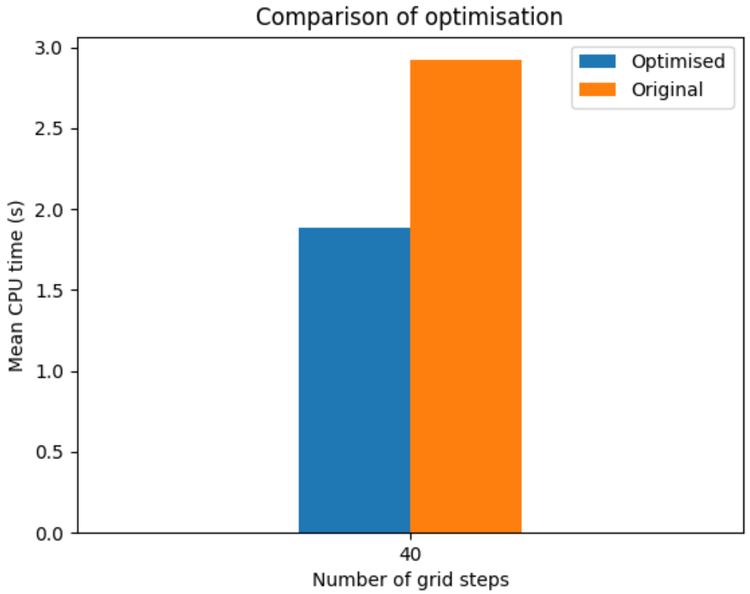


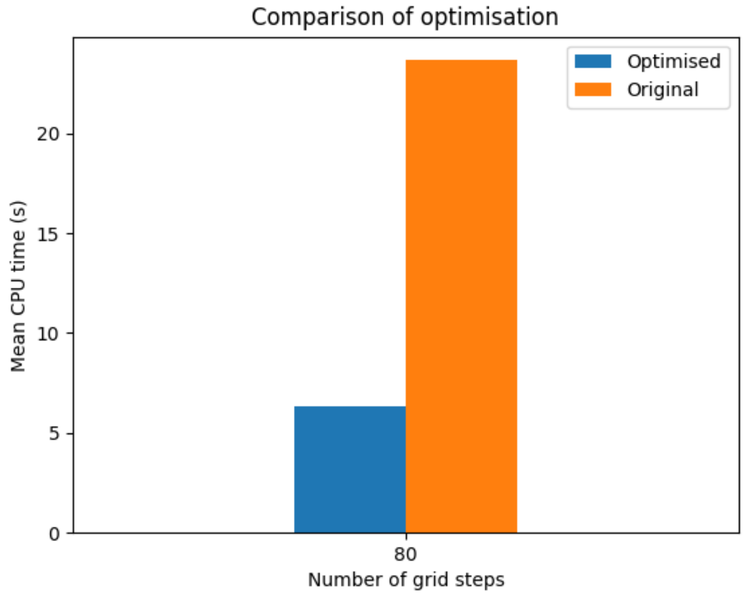
(These plots are from separate runs)

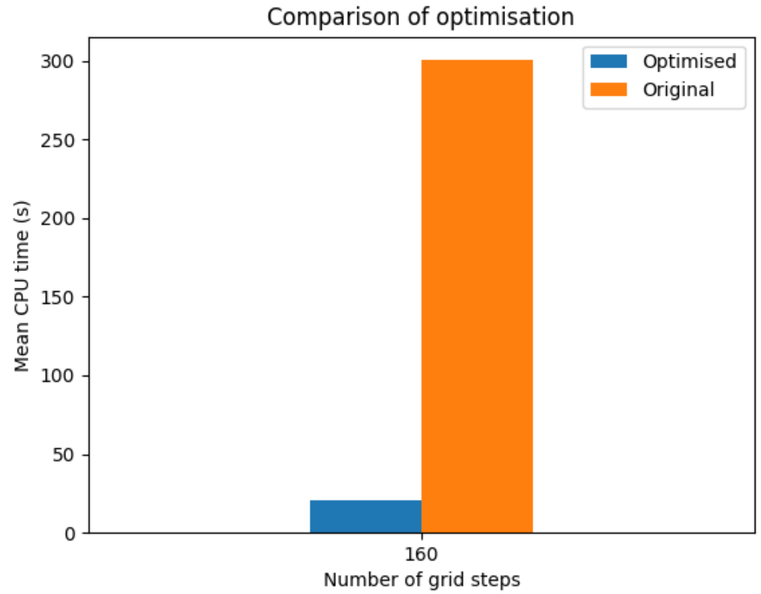
This plot took many hours to make

Optimisation

- Every time the field is requested, all grid nodes are looped over
- To avoid this we can just store a list of ‘active’ nodes and iterate over that
- For small gains (~10^2) this gives a major speedup and makes fine grids more accessible

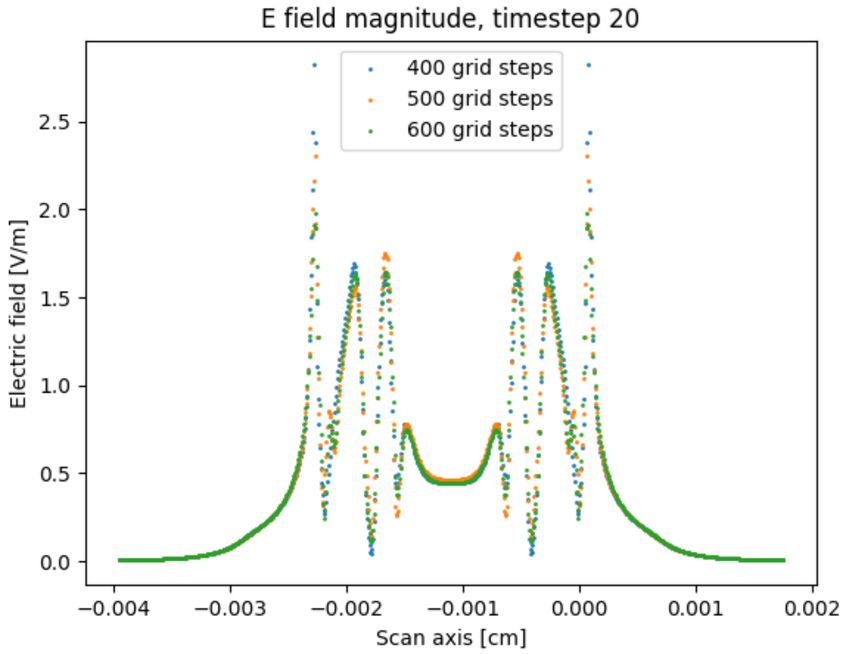


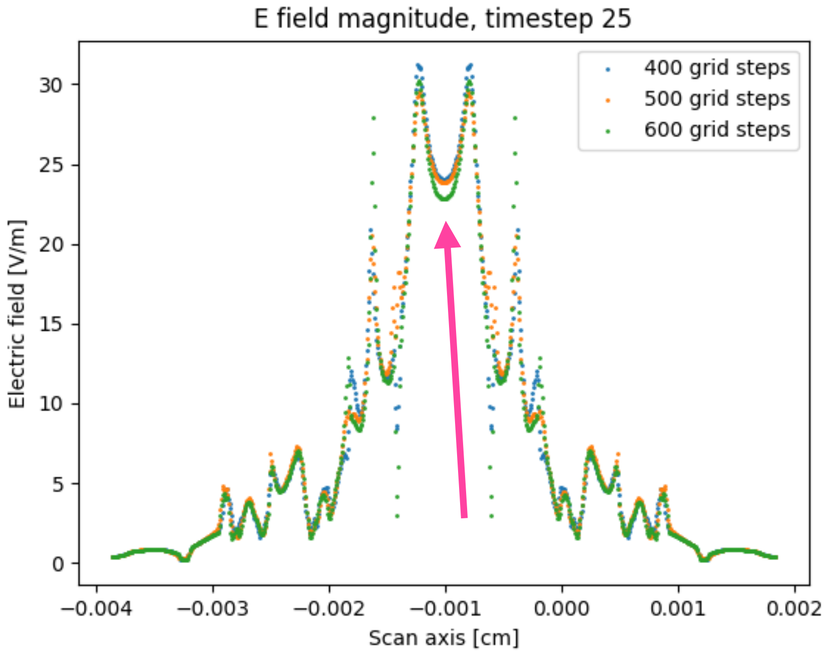


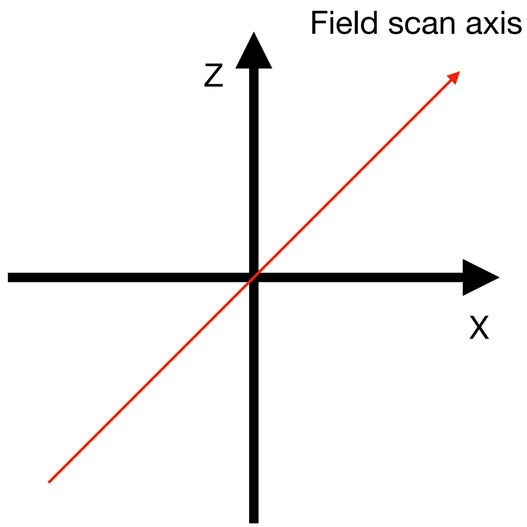


Convergence

- At higher grid spacings there is better convergence but still a small disagreement in certain regions







These plots are centred on the mean position of all particles

Future work

- Investigate using a point charge field instead of axisymmetric rings
- Run many avalanches with and without space charge effect and investigate the impact
- We are still working in a free space approximation: try adding a parallel plate boundary condition and see what the effect is
- Optimise further…