



GEM simulations: a flexible and efficient microscopic model based on Garfield++



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- **Experimental Context**
- **Simulation tools**
- **Models/Approaches**
- **Results**



SuperBigbite Spectrometer for HallA @ JLab

Physics:

Nucleon Form Factors at high Q^2

Neutron spin and TMD's, high statistics, high x_B

Pion structure functions

... an experimental tool for hadron structure investigation

- Large luminosity
- Moderate acceptance
- Forward angles
- Reconfigurable detectors

Expected Background:

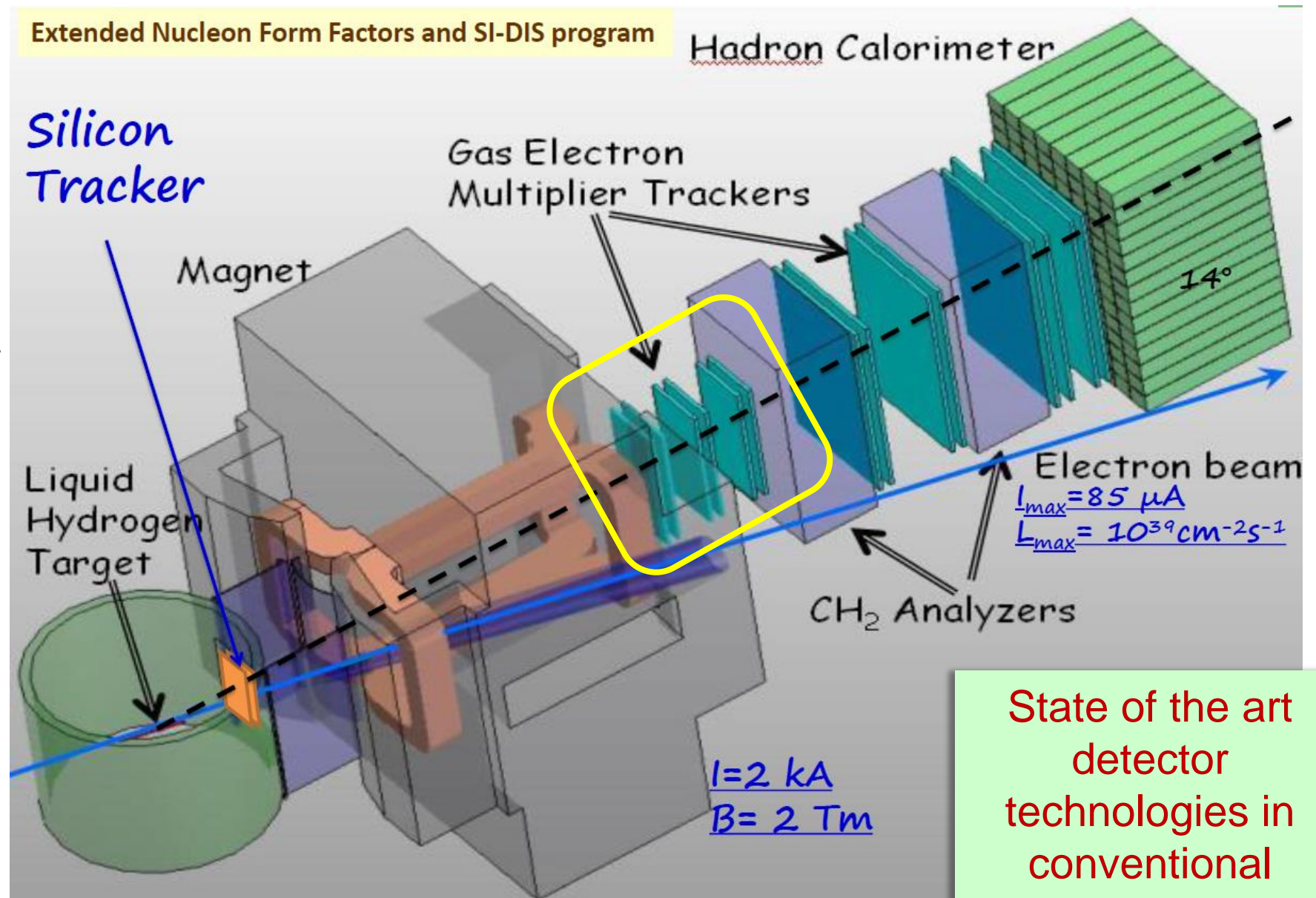
photon (250 MHz/cm²)

charged (160 kHz/cm²)

Kondo GNANVO talk (tomorrow on SBS GEMs Commissioning)

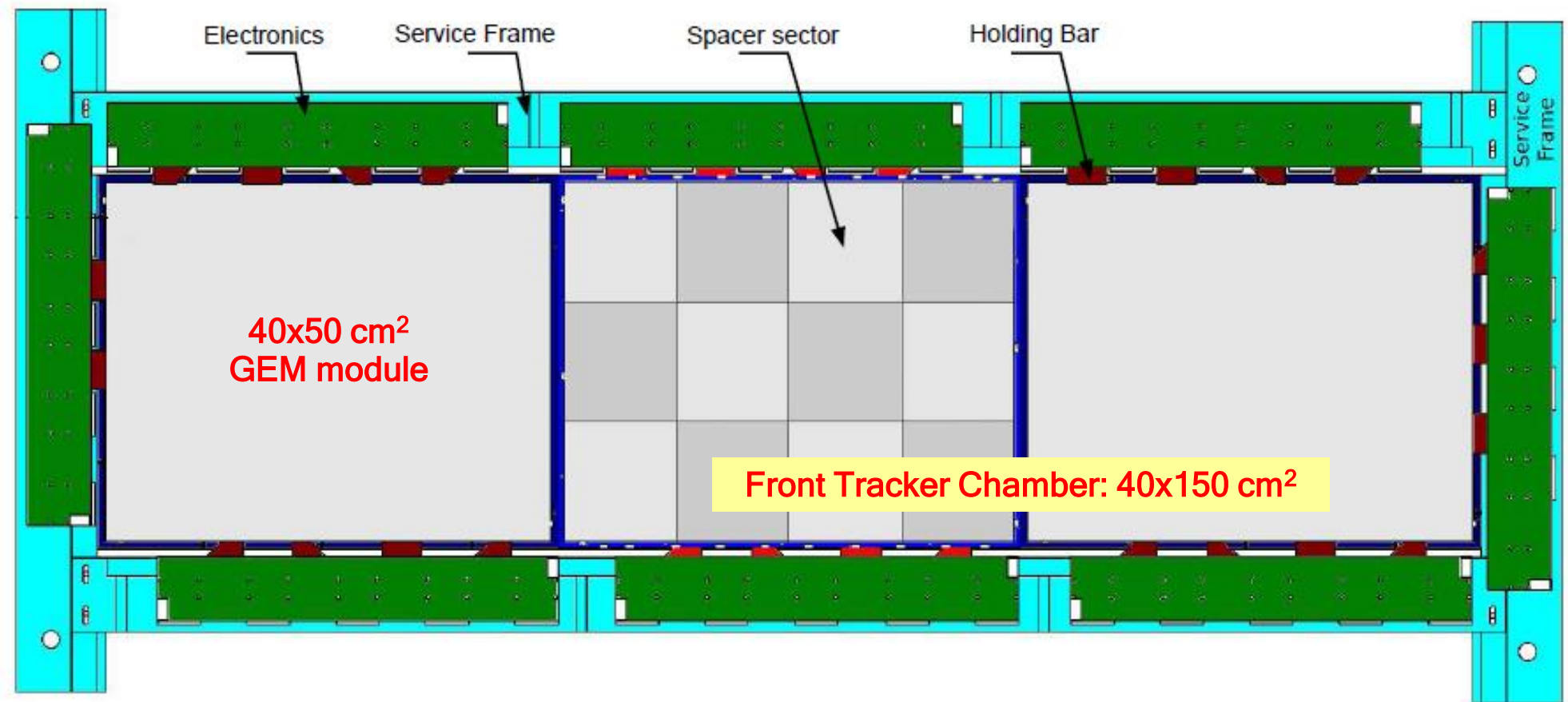
Extended Nucleon Form Factors and SI-DIS program

Hadron Calorimeter



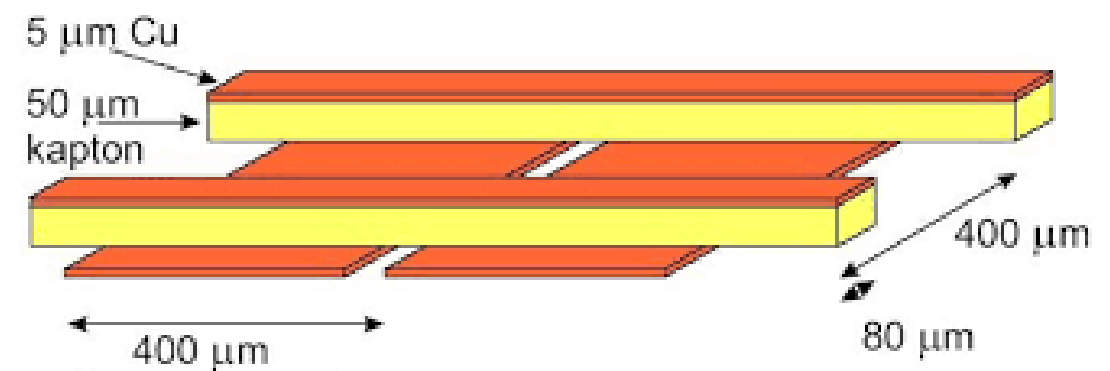
State of the art detector technologies in conventional configuration

Front Tracker GEM for SBS



3xGEM: single mask GEM foil production
x/y Strip Readout Plane (a la COMPASS)

Double strips with 90° angle



Main Tools / Libraries

Two years ago we decided to implement a «microscopic» GEM simulator to better model the real behavior of the tracker and to prepare for the physics experiments

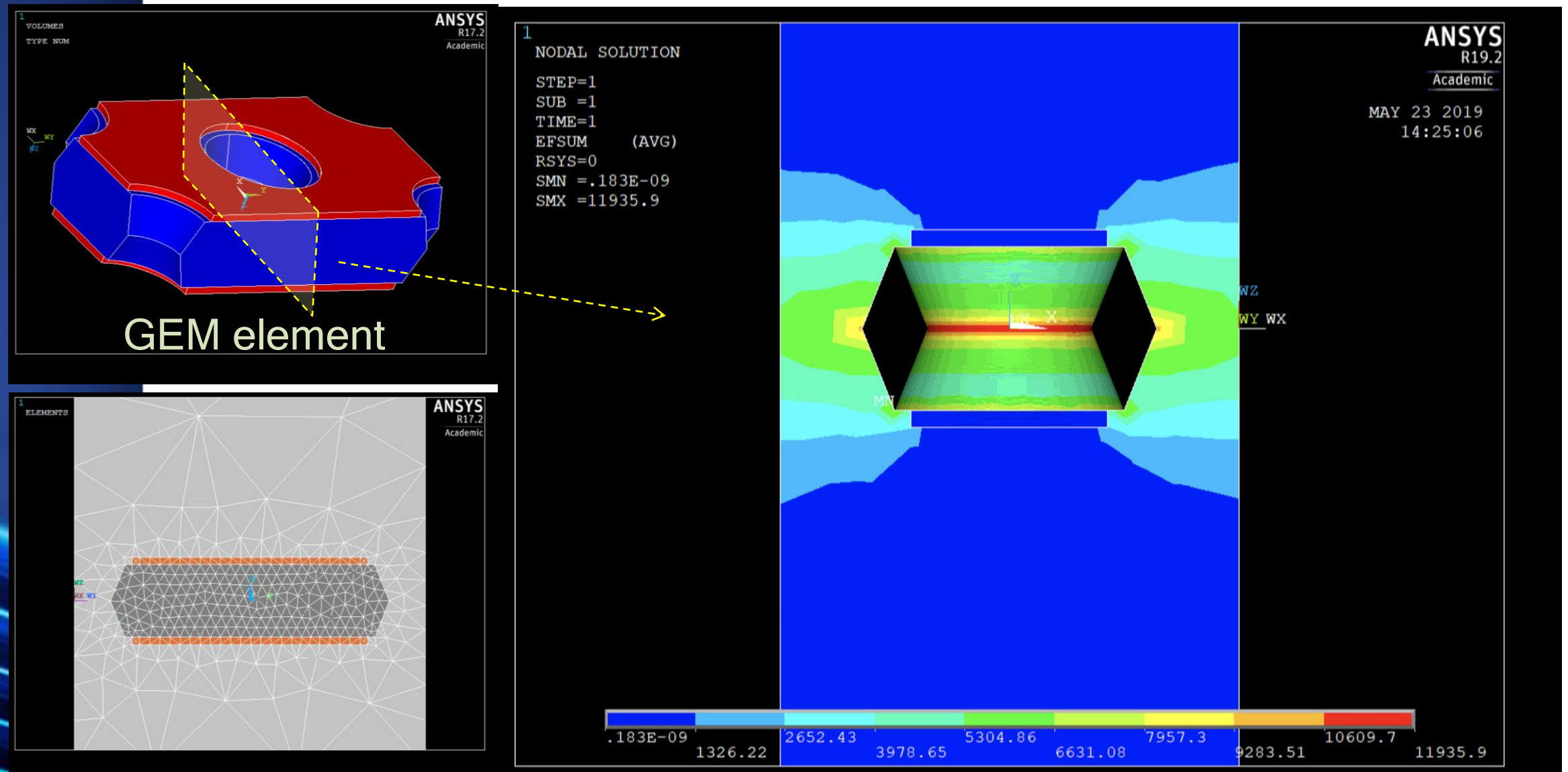
Geometric Model Definition	GMSH (3.05 and 4.0.7) OpenCascade	ANSYS Multiphysics (R17-19)
Mesh Generation		
Electrostatic Field Solution	ELMER (2018)	
Gas properties	Magboltz	
High Energy particle Photoabsorption ionization	HEED	
Microscopic Gas Simulator	Garfield++ (v3)	
Automate and parallelize multiple simulations	Bash scripts	
Analysis/Visualization	ROOT	
Simulation Platforms	Xeon/16 cores and i7/16 cores	

All steps are implemented programmatically

ANSYS Mechanical APDL software

Engineering software useful to create complex geometries, assign materials to different volumes and create electrostatic field solutions.

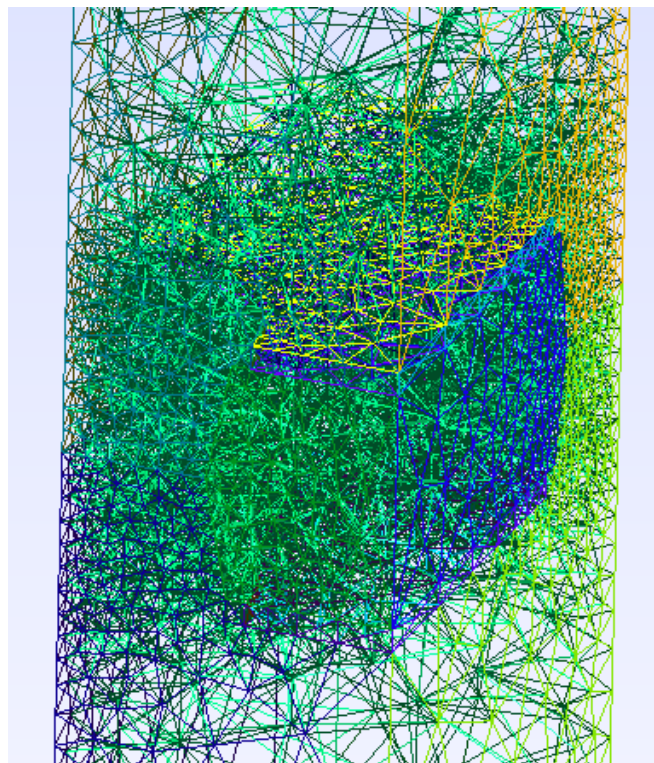
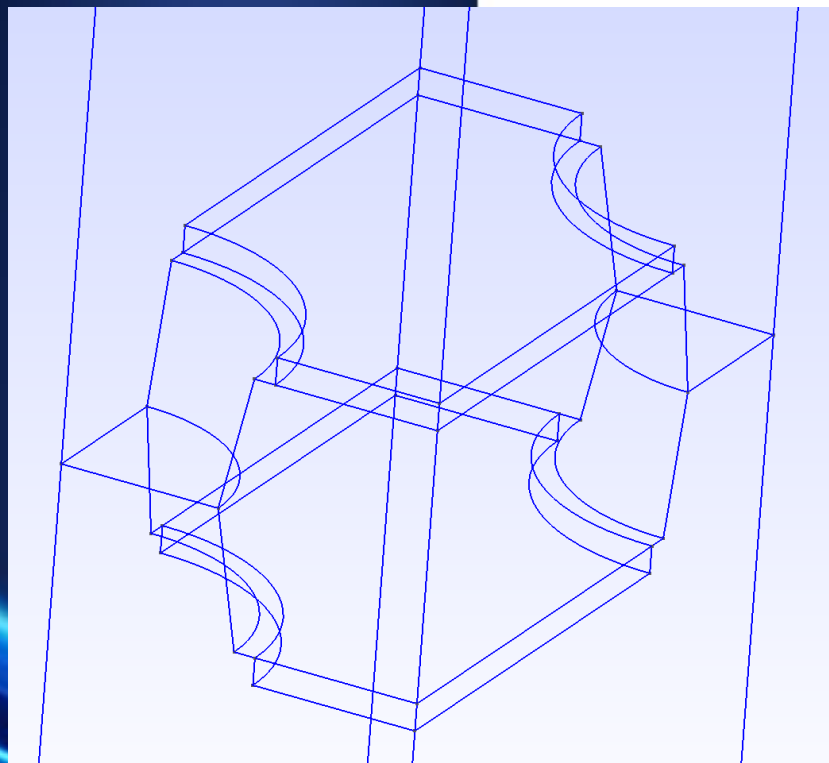
Model design, mesh generation and electrostatic field within a reasonably coherent framework; mesh generation is well optimized; better visualization tools; **need license**



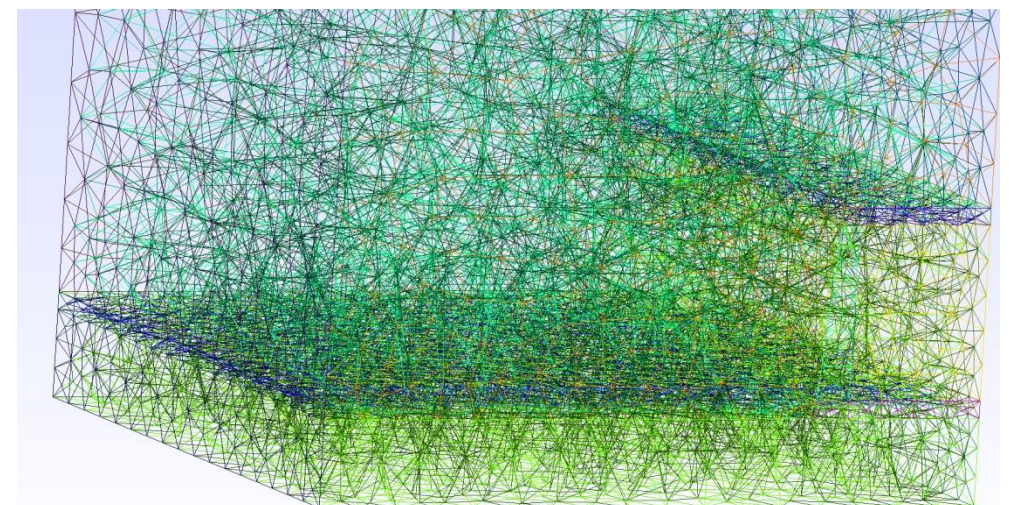
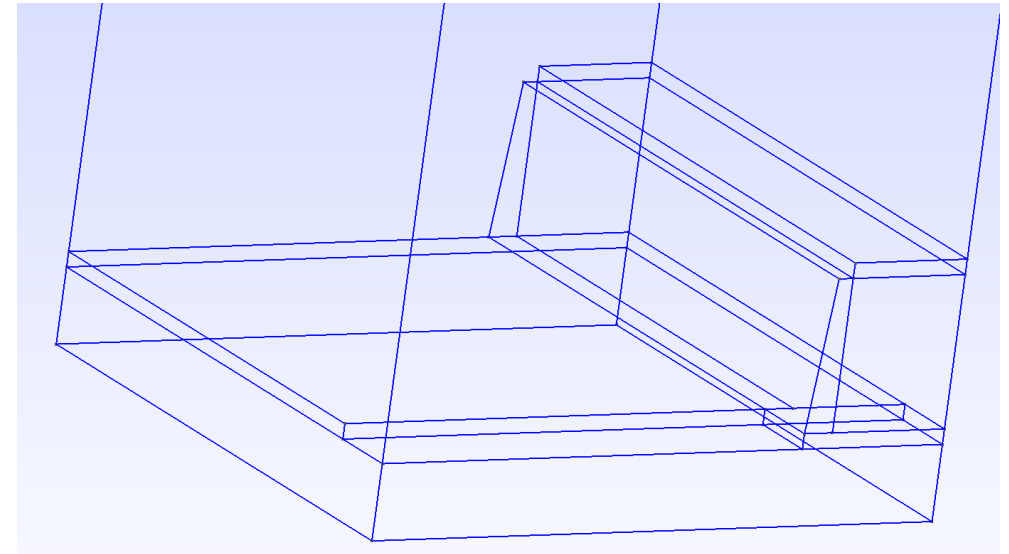
GMSH + ELMER

Open source software; reasonably well documented; somehow more flexible scripting language (c/c++ oriented) than in ANSYS
GMSH versions incompatibility, decent mesh generation (and field solution) is not straightforward as in ANSYS

GEM foil element



x/y strip readout



Denser mesh nodes near foil parts

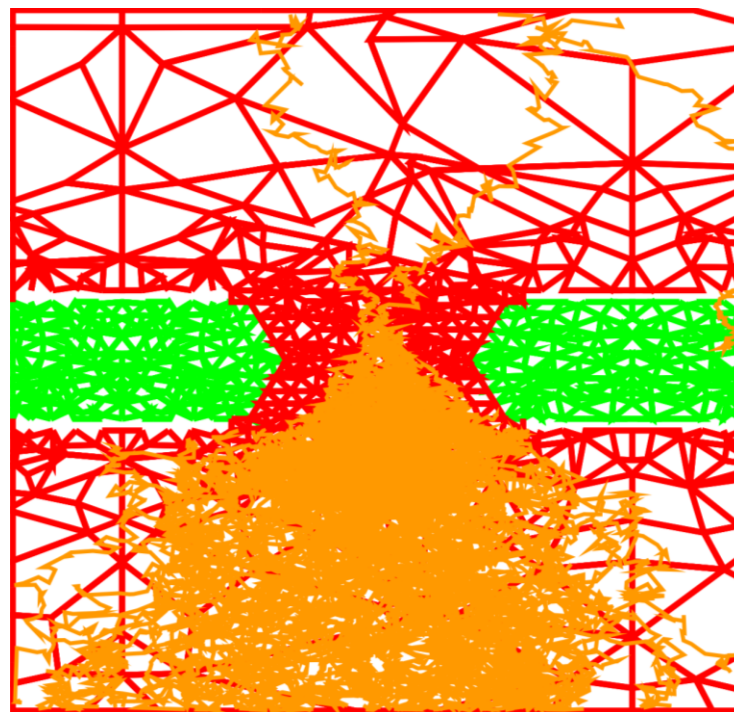
Garfield++ tool

HEED: simulate the photoabsorption ionization of the high energy particles (GeV electrons and protons)

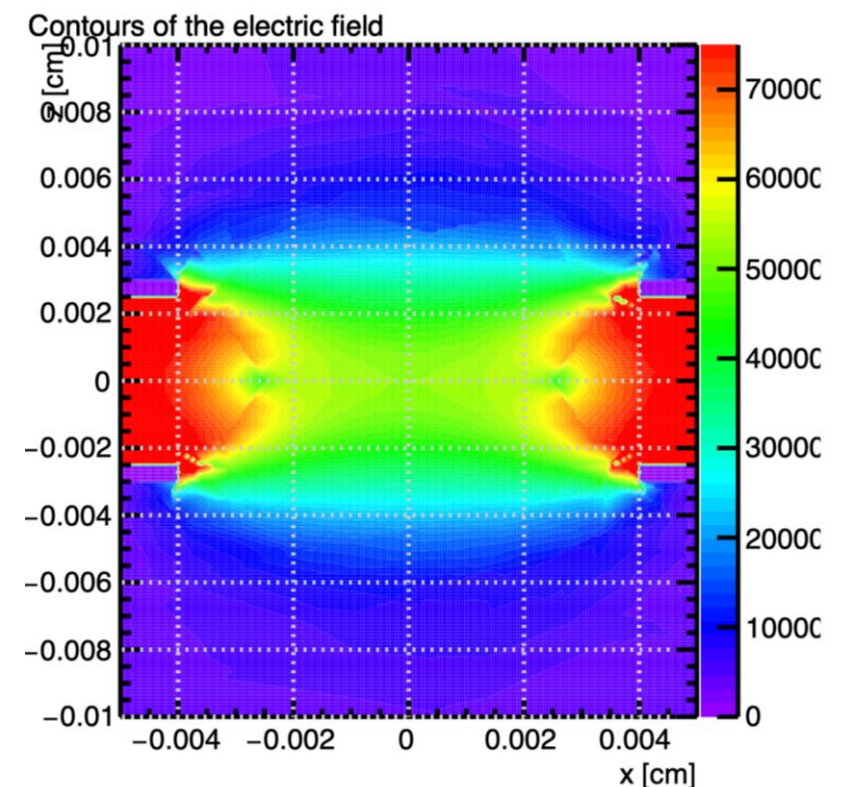
Magboltz: define the atomic and thermodynamic characteristic of the gas:
Mixture **70% Argon** and **30% CO₂** (at STP); Penning Transfer ON

Garfield++: AvalancheMicroscopic for e- and AvalancheMC for ion

ROOT: for analysis and visualization



Avalanche creation and its propagation

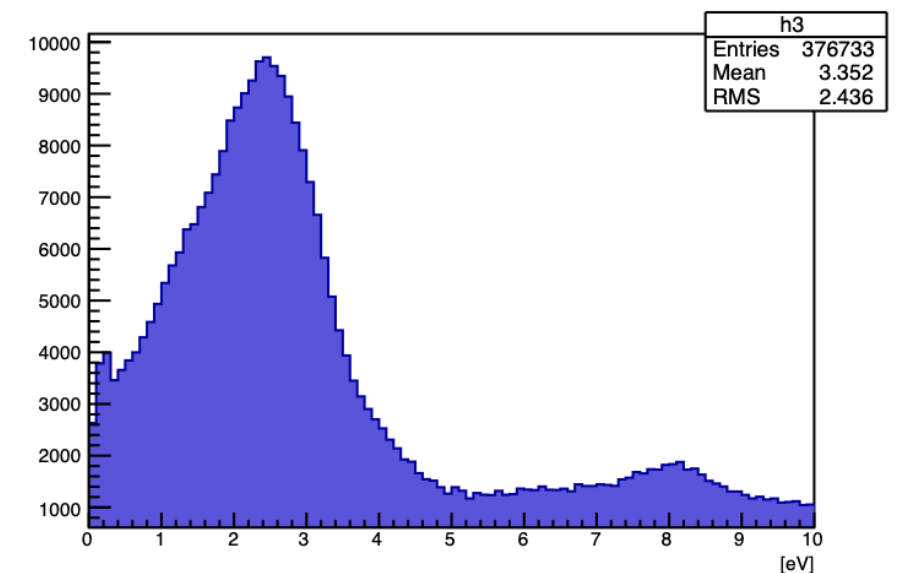
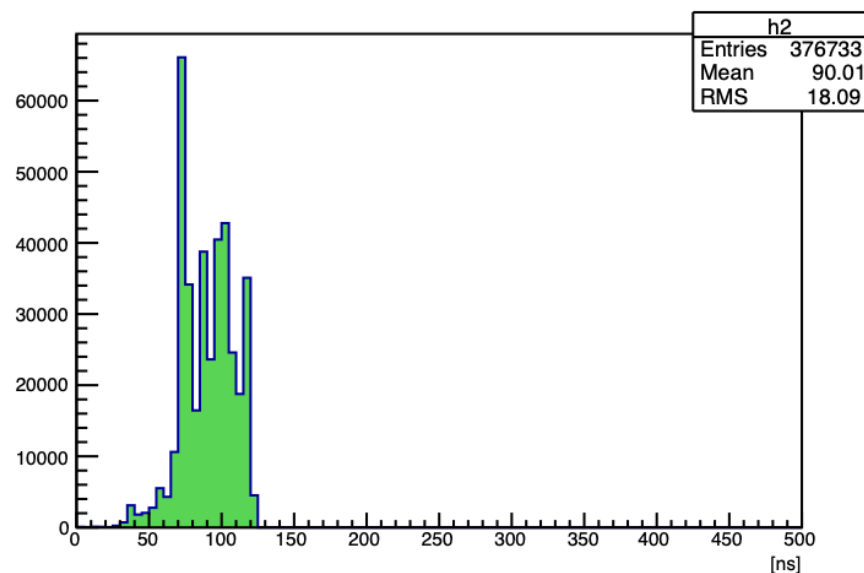
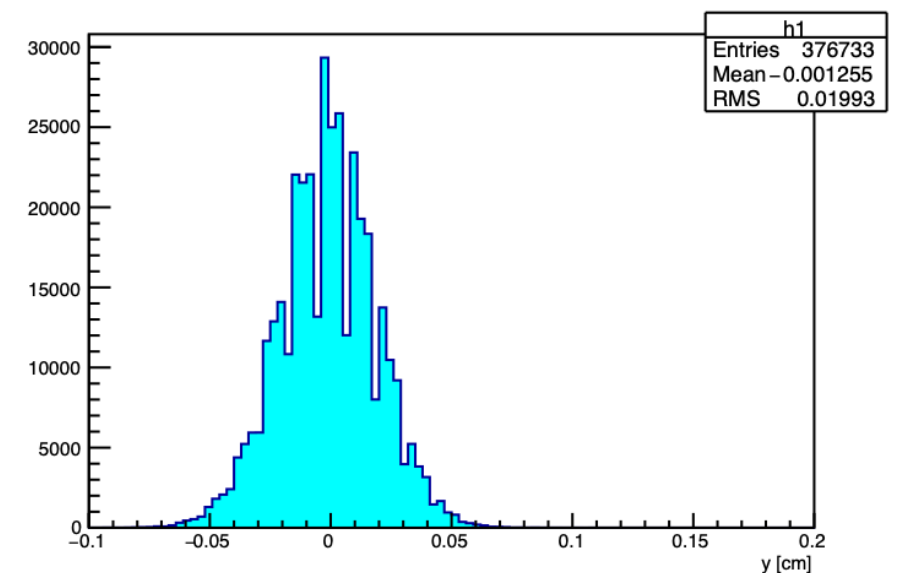
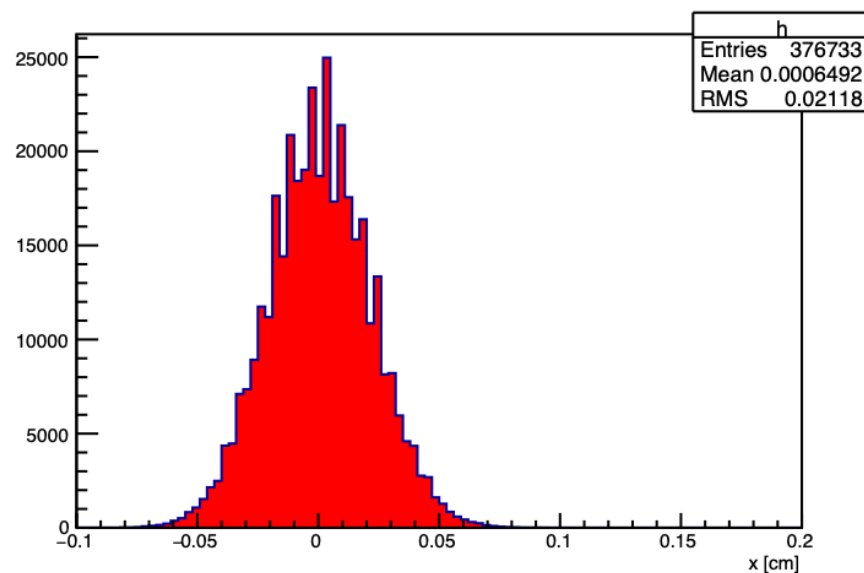


Electric Field map

Microscopic Simulations

Systematic study to verify the consistency of the simulations:

- Number of secondary electrons reaching the readout plane (strips)
- Spatial distribution of the avalanche (x and y axes)
- Secondary electrons drift time
- Energy of the particles in the readout plane



Simulations models

Single GEM foil

Simulate 2.8 GeV protons passing through a single GEM chamber with ideal readout plane. Used to compare GMSH+ELMER and ANSYS models

Multistep 3xGEM+RD

Microscopic simulations are carried on each GEM layer in a sequential way: simulation outcome of the previous layer is the input of the next.

It's a flexible multistep model that easily allows to simulate different schemes, imperfections, foil misalignment, by decomposing the 3xGEM+Readout chamber in 4 adjacent layers.

Full 3xGEM

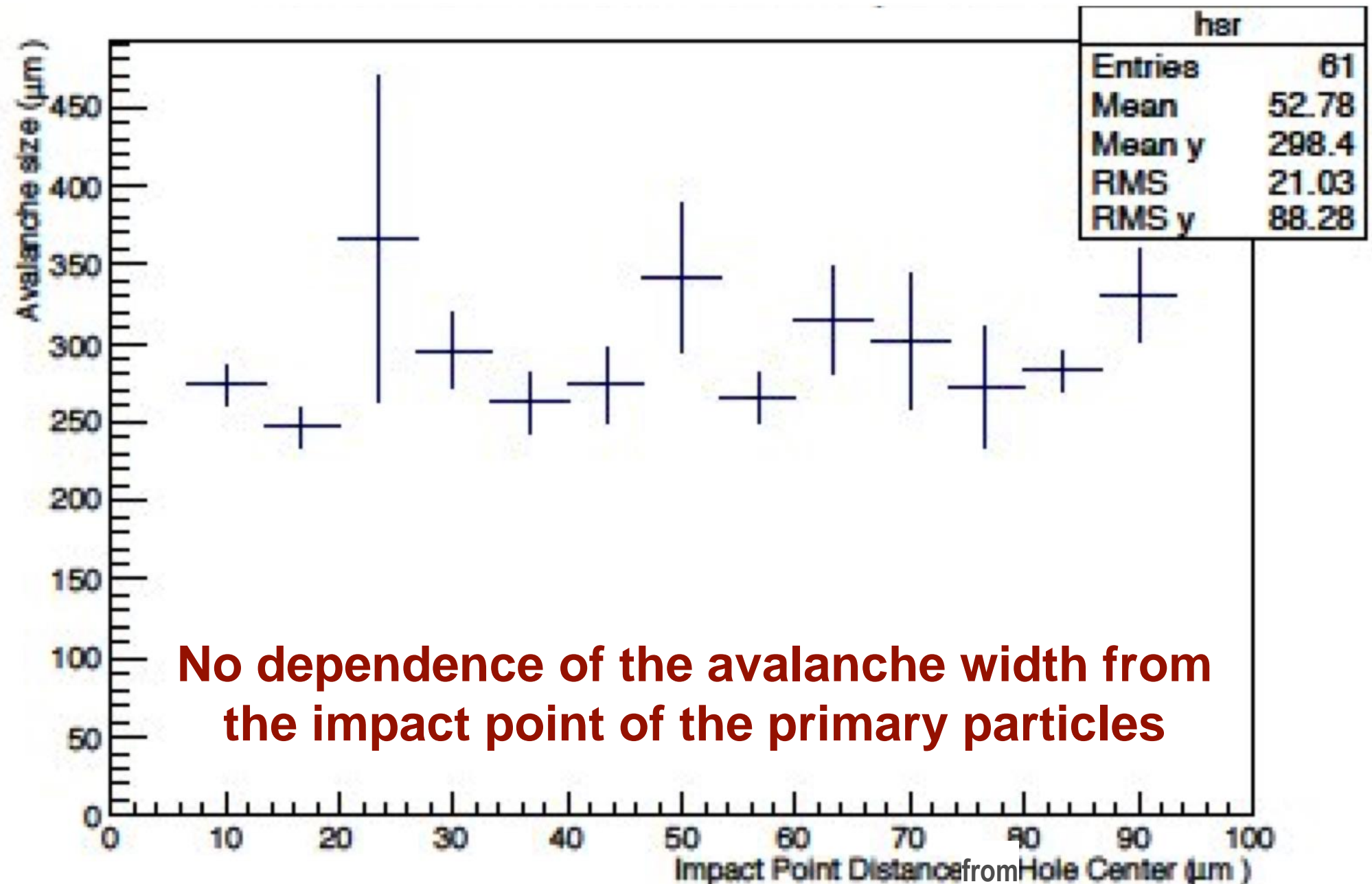
Microscopic simulations are carried from the drift plane, through the 3 GEM foils, to an ideal readout plane.

Segmentation of the readout plane is performed on the collected electrons (no readout structure in the simulation)

Somehow used as reference and comparison for the multistep model.

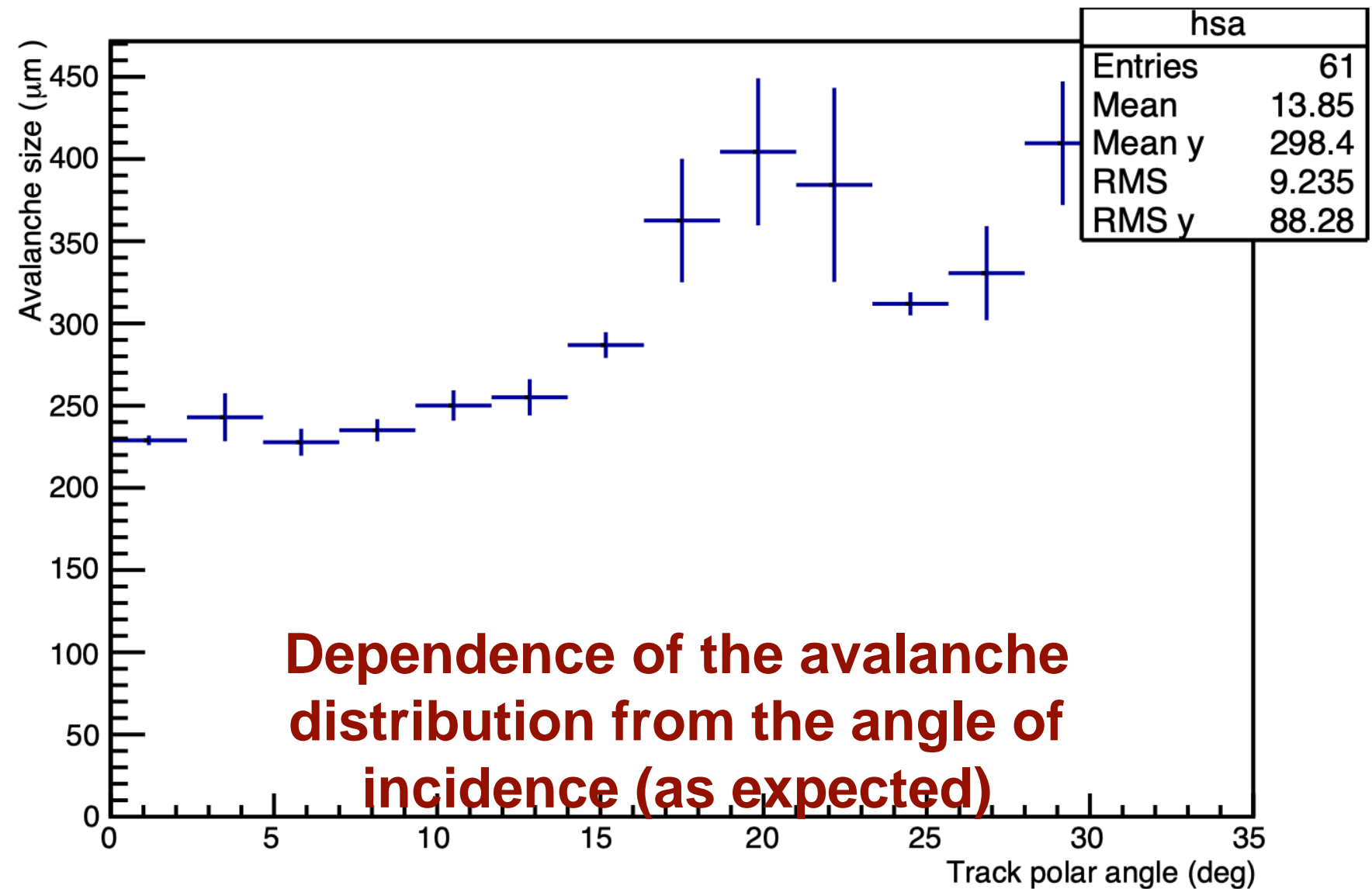
Simulation - some consistency checks

Avalanche width VS impact point of primary particles



Simulation - some consistency checks

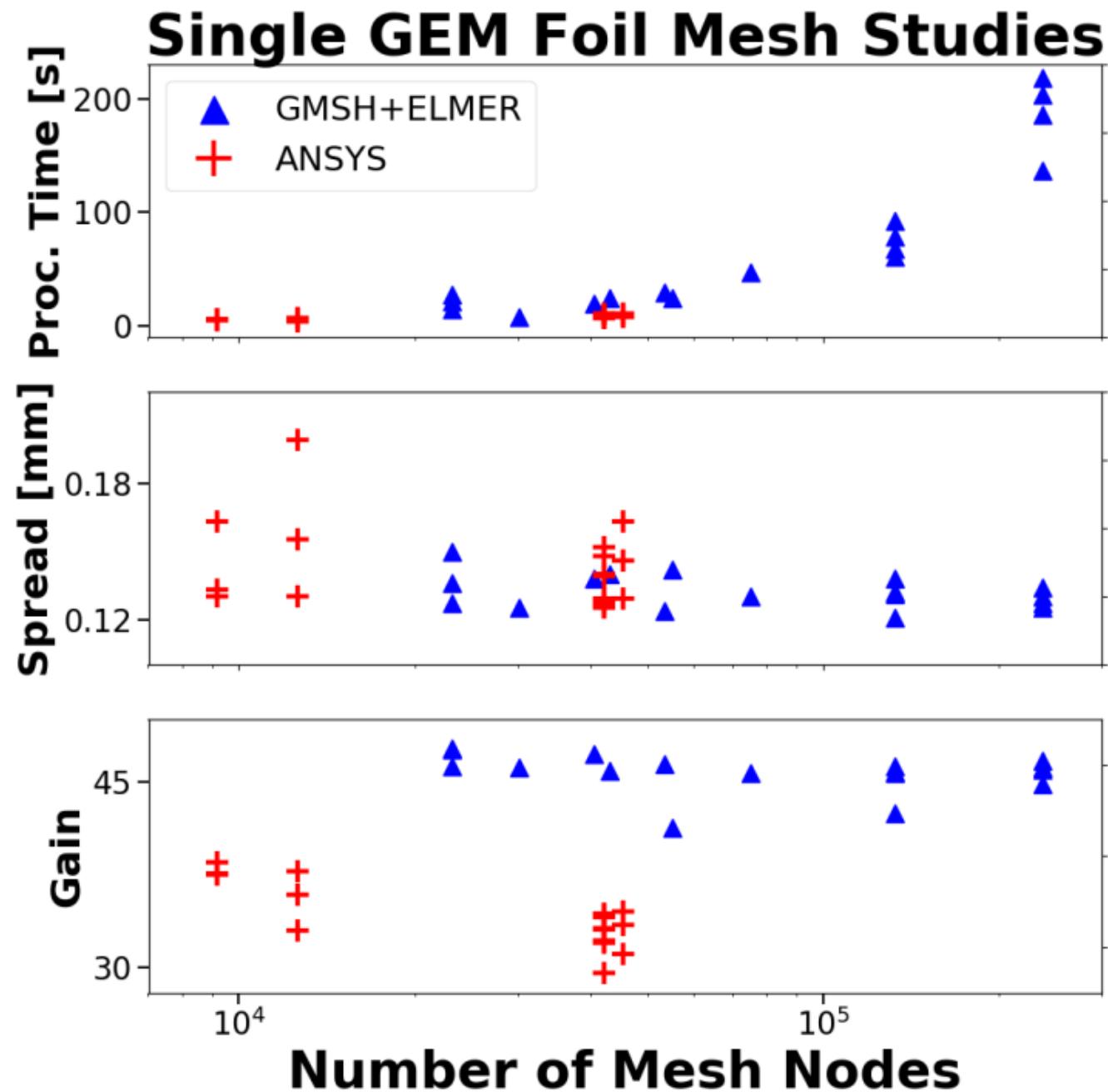
Avalanche width VS incidence angle of primary particles



Incidence Angle	0°	10°	30°
RMS	233 μm	295 μm	381 μm

Single GEM: GMSH+ELMER vs ANSYS

We compared response of single GEM foil simulation from both modellers/solvers at different mesh sizes



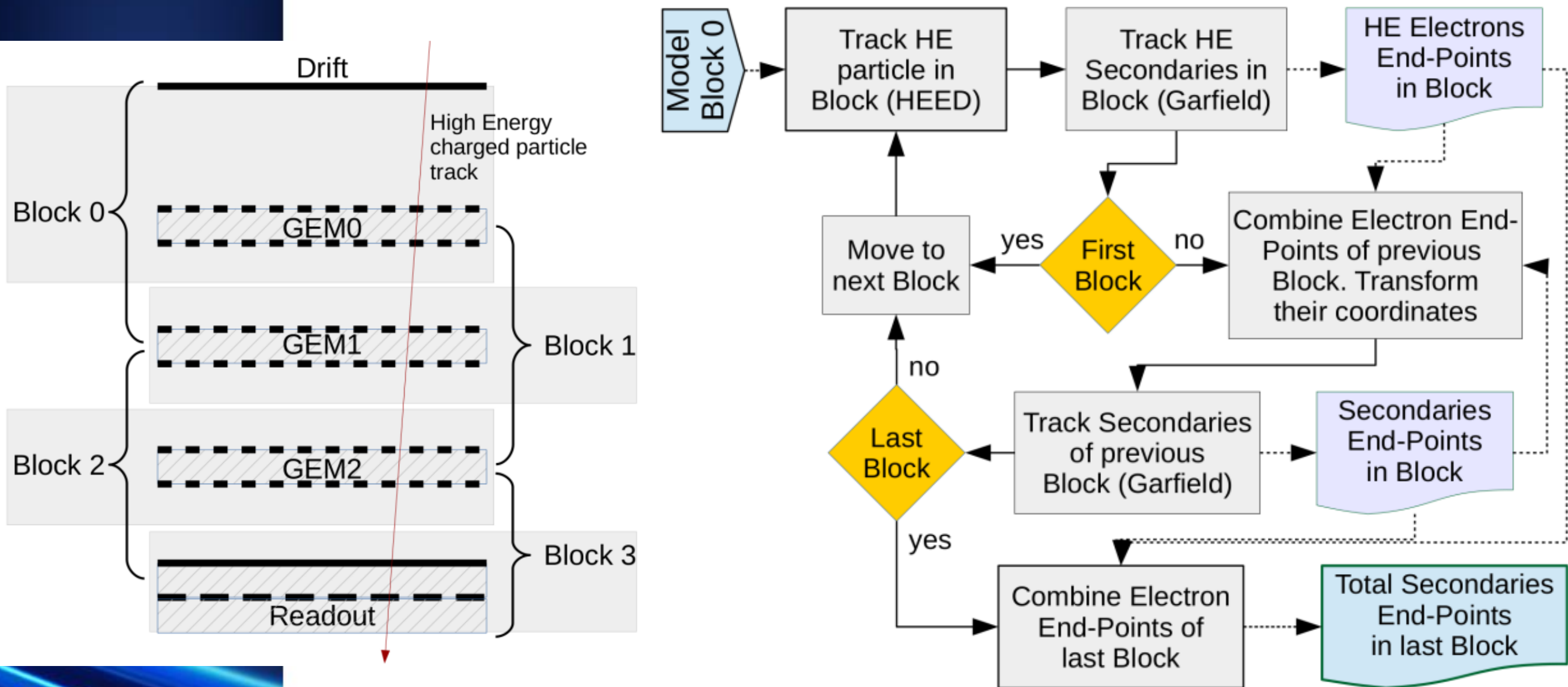
2.8 GeV protons, 0 incidence angle on Drift [-1836 V] +3 mm + [-1116 V] GEM [-721 V] + 2 mm + [0 V] ideal readout

- Processing times and Avalanche spread consistent between G+E and A
- Avalanche spreads (x/y show similar spreads) reasonably constant over mesh size, but larger meshes produce more stable results
- Gains discrepancy between G+E and A; G+E based simulations look more stable over mesh size

Gain = average number of electrons collected after the GEM foil on an ideal readout plane, for each primary ionized electron

Multistep - 3xGEM+RD approach

Assume approximate independence of each GEM foil response:
the 3xGEM+Readout is decomposed into 4 overlapping simulation blocks



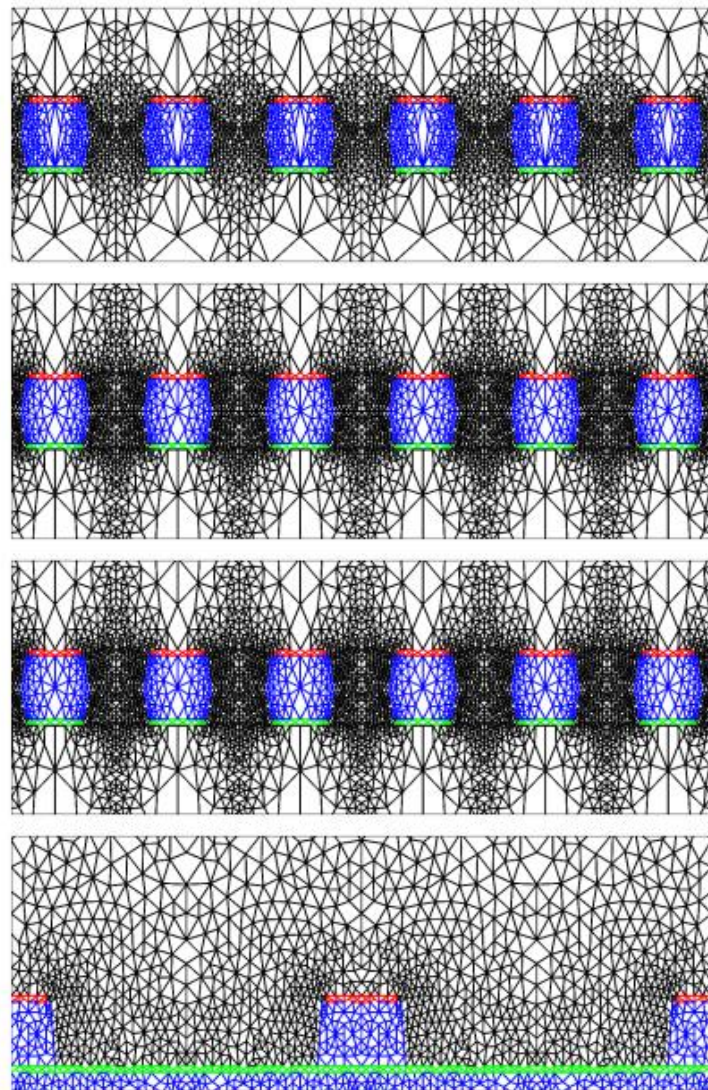
Flexible approach, easy integration of readout plane, allows to simulate different schemes, imperfections, foil misalignment

DO NOT CONSISTENTLY HANDLE ions drift on different blocks (YET)

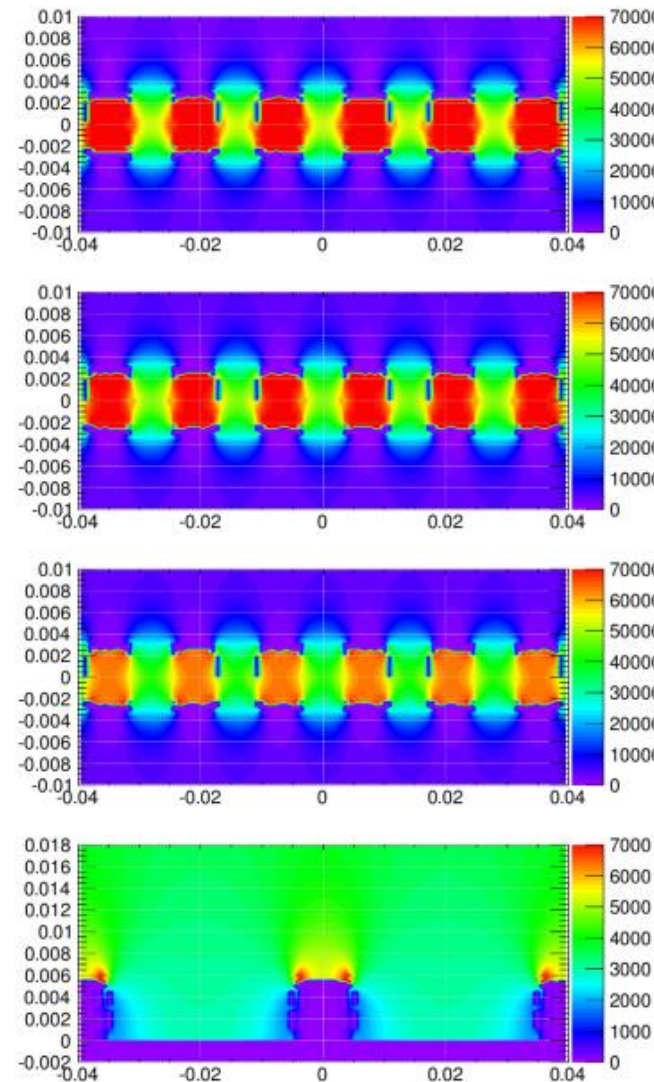
Multistep – 3xGEM+RD model and simulation

GMSH + ELMER models simulated

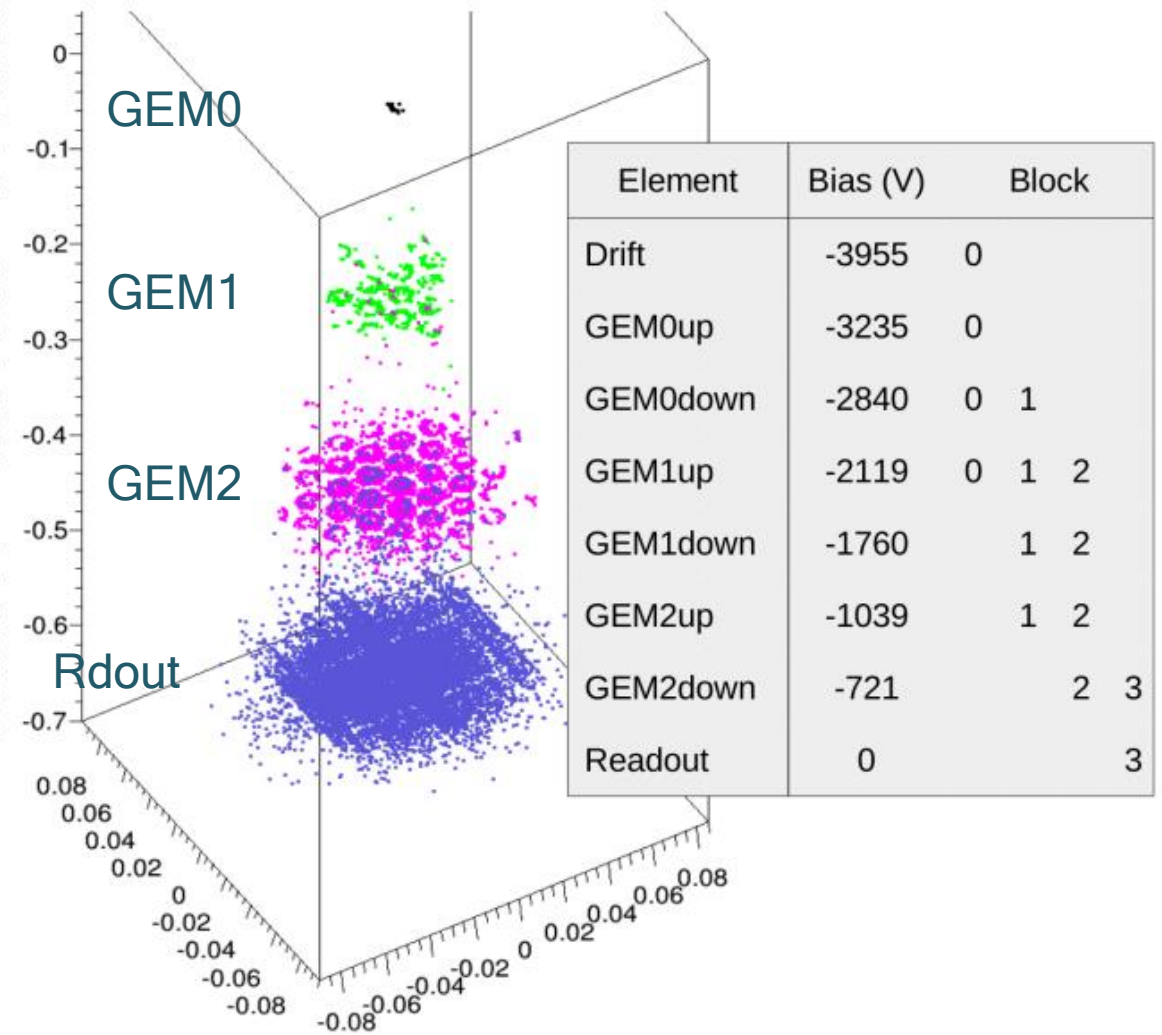
Mesh



Electrostatic Field



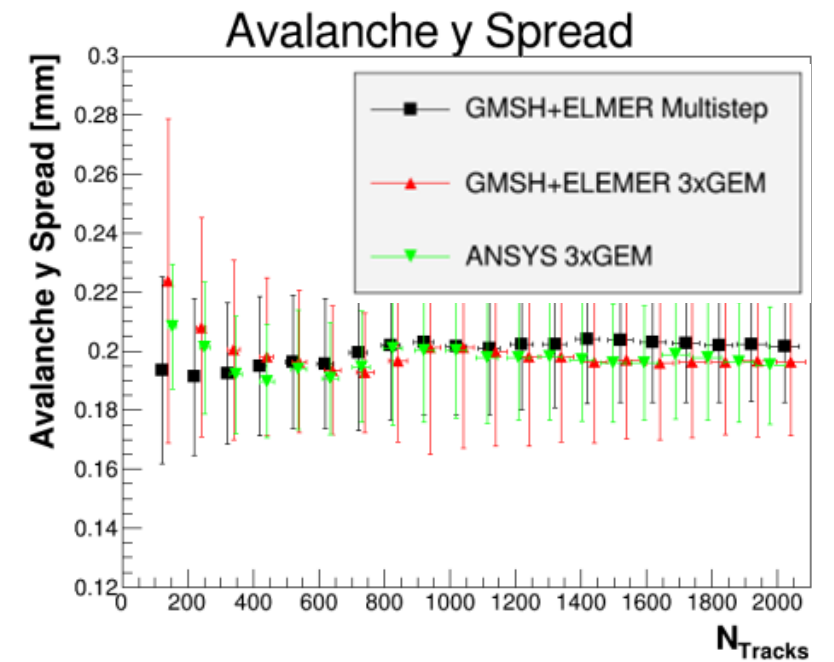
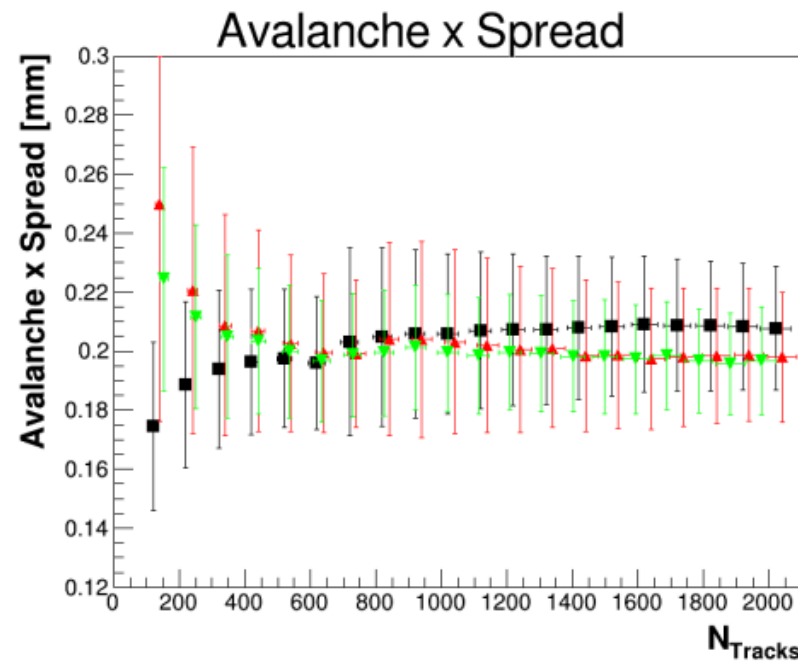
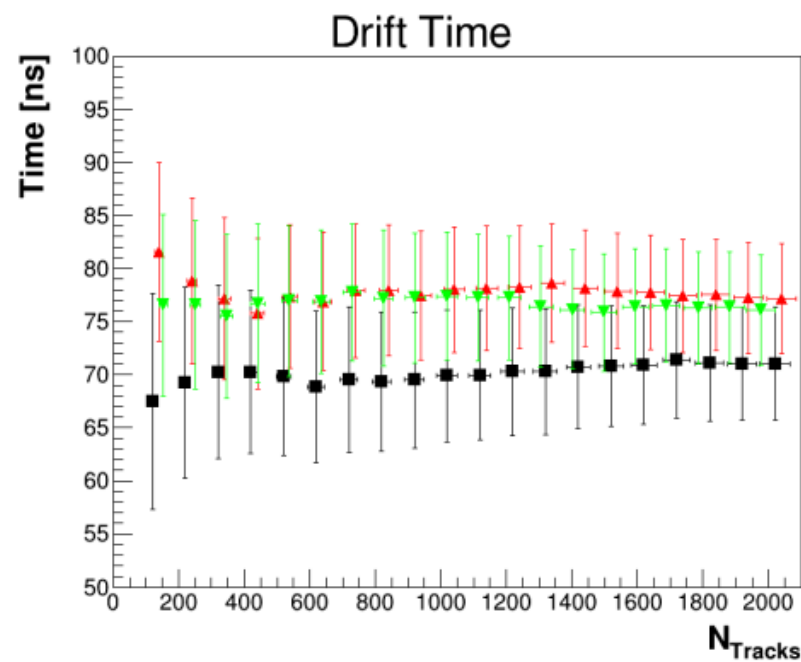
Garfield e- endpoints



Each block is simulated in sequence (following avalanche development)

Different approaches comparison

GMSH+ELMER multistep approach and full 3xGEM
ANSYS on full 3xGEM

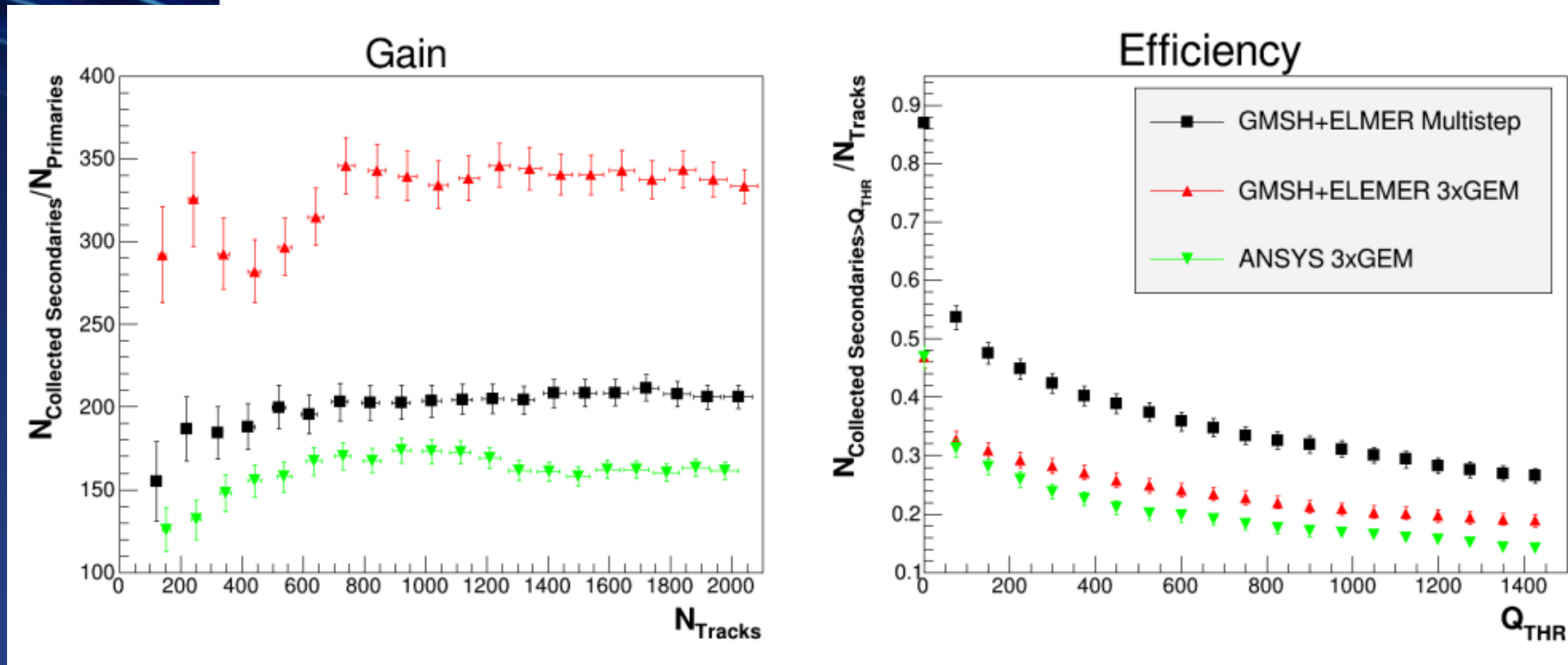


Cumulated plots (increasing number of simulated tracks)

All quantities tend to reach rather stable estimates when averaged above ~ 800 tracks

e- drift time and avalanche x/y spreads consistent within statistical uncertainties

Different approaches comparison



Efficiency = fraction of high energy tracks producing a number of secondaries collected on the readout plane larger than the given threshold Q_{THR}

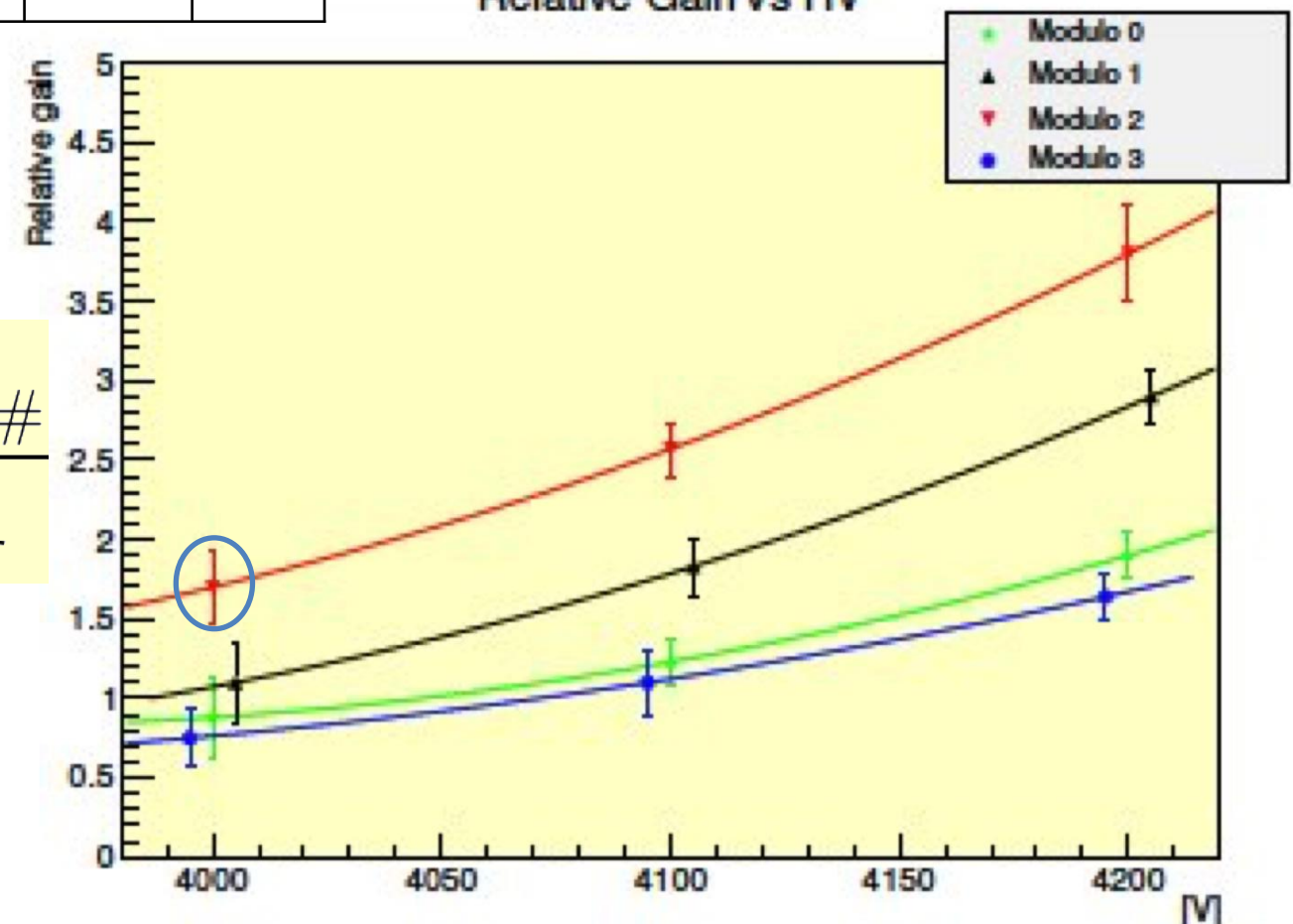
Gain and Efficiency show discrepancies between models
Gain differences between G+E and A qualitatively consistent with single foil results
→ Discrepancy most likely related to the different meshes

Preliminary comparison with real data

We have quite detailed test-beam (2.8 GeV p) data of 5 GEM modules

Module	HV Divider Resistors [MΩ]							
	R ₁	R ₂	R ₃	R ₄	R ₅	R ₆	R ₇	R ₈
0	0.441	7.2	3.76	7.2	3.63	7.2	2.98	7.2
1	0.441	7.2	3.92	7.2	3.57	7.2	3.57	7.2
2	0.441	7.2	3.92	7.2	3.92	7.2	3.92	7.2
3	0.441	7.2	3.76	7.2	3.66	7.2	2.98	7.2
Reference	0.441	4.8	2.66	4.8	2.66	4.8	2.27	4.8

Relative Gain vs HV



$$RelativeGain = \frac{[I_{HV}^{Beam} - I_{HV}^{NoBeam}]_{Mod\#}}{[I_{4100}^{Beam} - I_{4100}^{NoBeam}]_{Ref}}$$

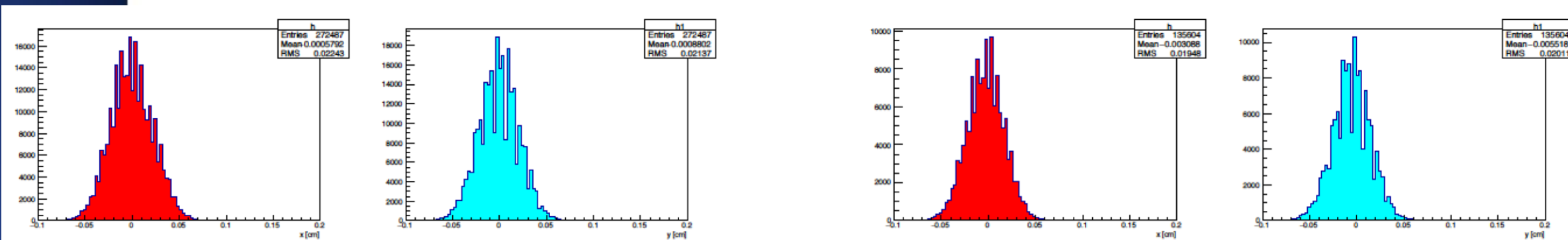
Preliminary Garfield++ prediction vs real data

ANSYS-Model

HV=4000 - Module 2 - Relative Gain

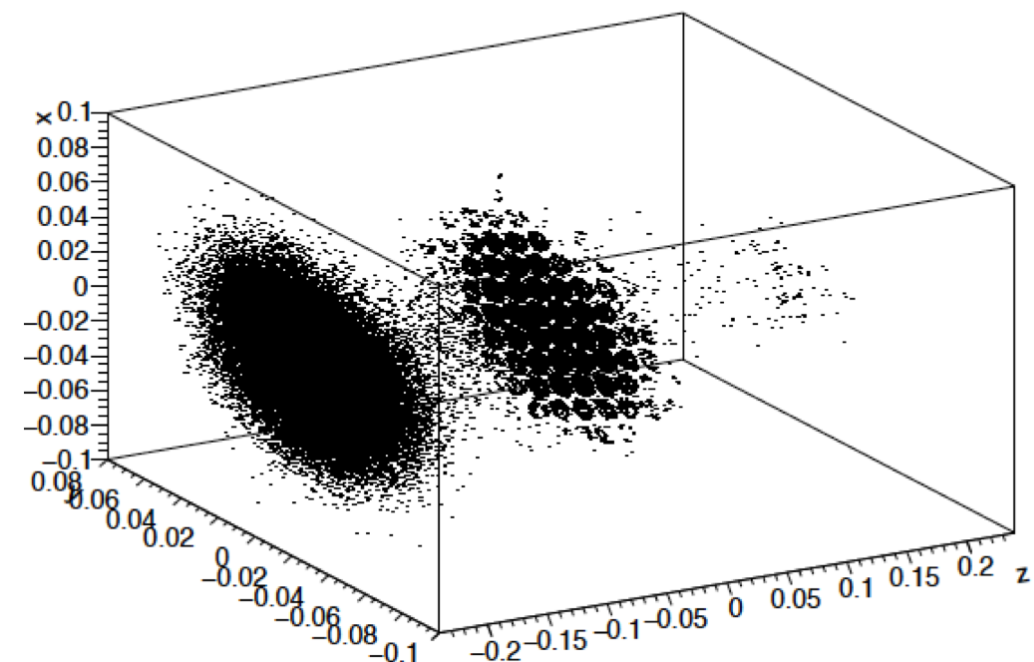
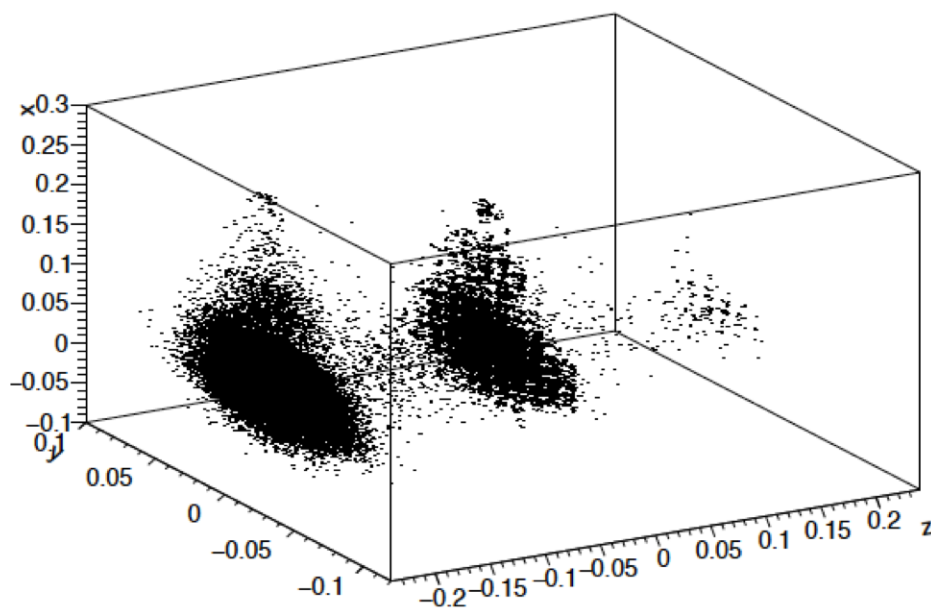
Predicted : 2.000 ± 0.013

Real data : 1.70 ± 0.25



Module 2 - Entries: 272487

Ref Module - Entries: 135604



Summary

Implemented GEM simulations based on:

- Garfield++ (& Heed & Magboltz)
- Programmatic models from both ANSYS Multiphysics and GMSH+ELMER

Compared ANSYS and G+E on single foil response vs mesh size → show different gain predictions (related to how the mesh is generated ?)

Implemented two different GEM chamber models:

- Multistep 3xGEM+x/y Readout: flexible, permit an easy integration of the readout plane; no consistent ion drift simulation yet
- Full 3xGEM: ideal readout only

Comparison shows:

- good agreements on drift time, avalanche spread
- discrepancy on gain prediction (but consistent with single foil)
- The multistep approach seems to work similar to the Full GEM; discrepancy likely related to mesh generation (and/or field solution) (?)

Very preliminary, limited, comparison of relative gain with real data of the 3xGEM model is encouraging