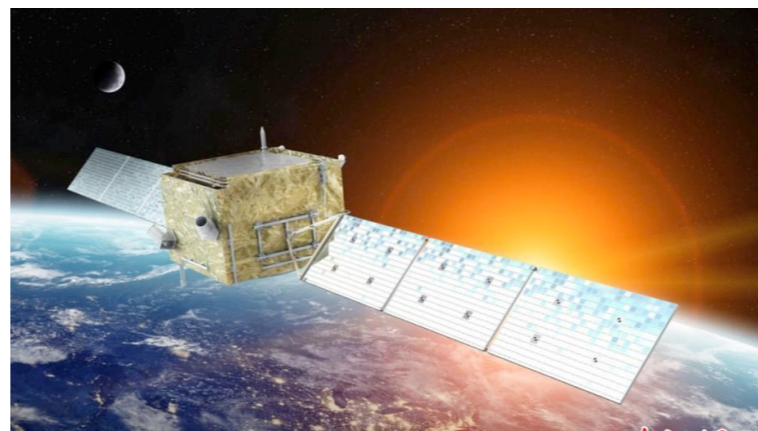


High-energy simulation experience from DAMPE

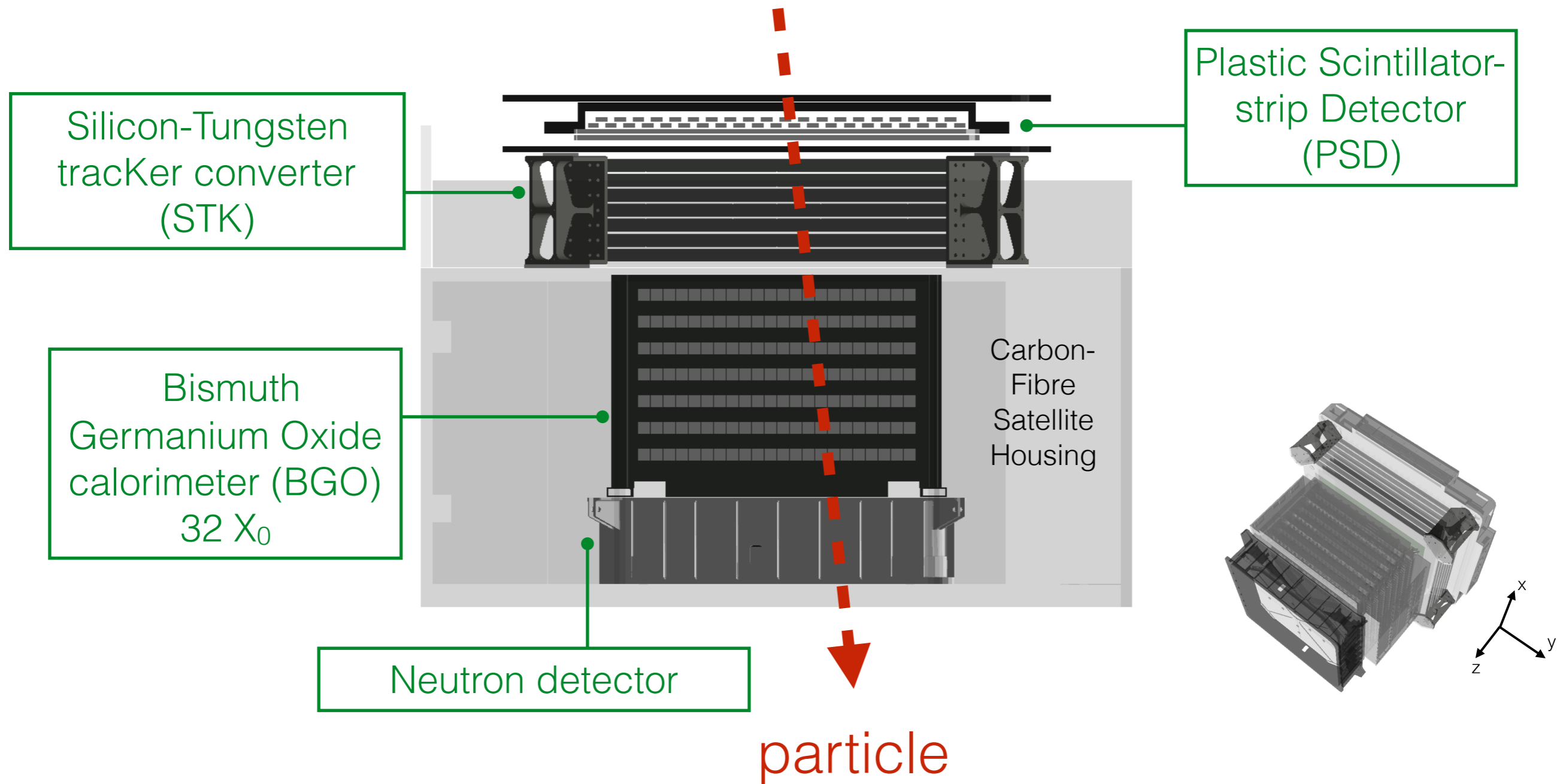
Andrii Tykhonov
University of Geneva



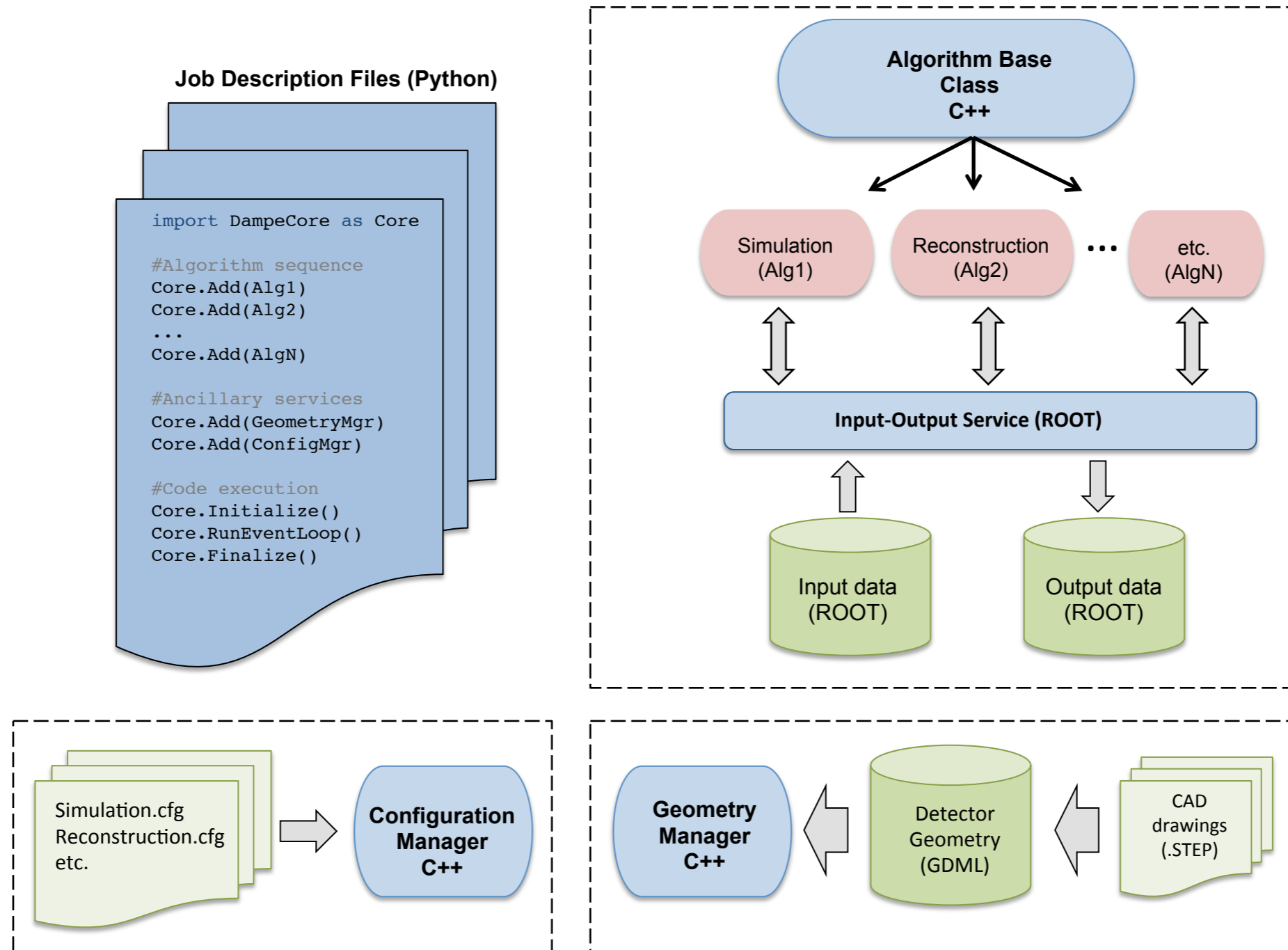
Outline

- DAMPE & Simulation:
 - Cosmic-Ray (CR) proton/ion analyses
 - High-energies (>100 TeV)
 - Cosmic-Ray (CR) electron analysis

The DAMPE detector

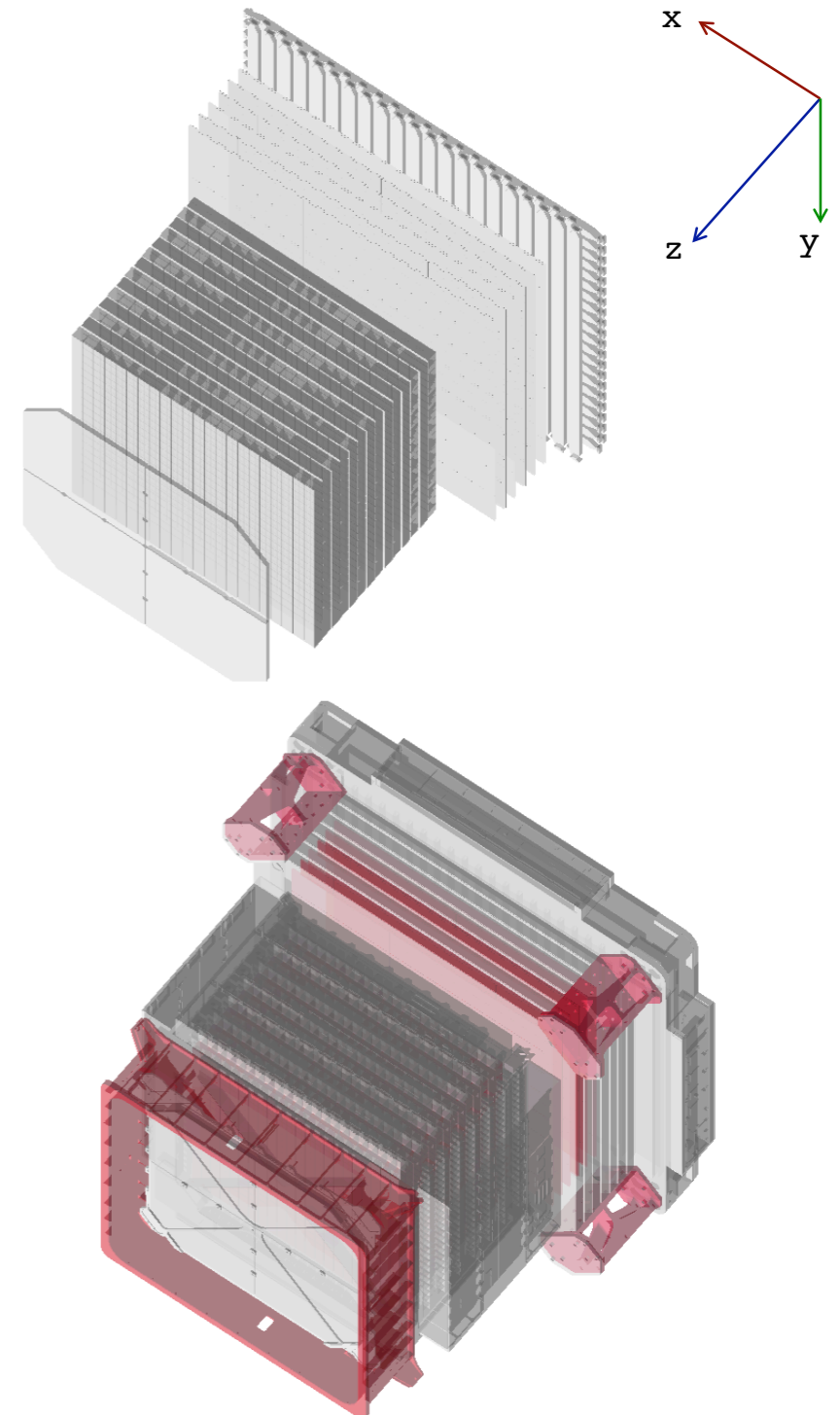
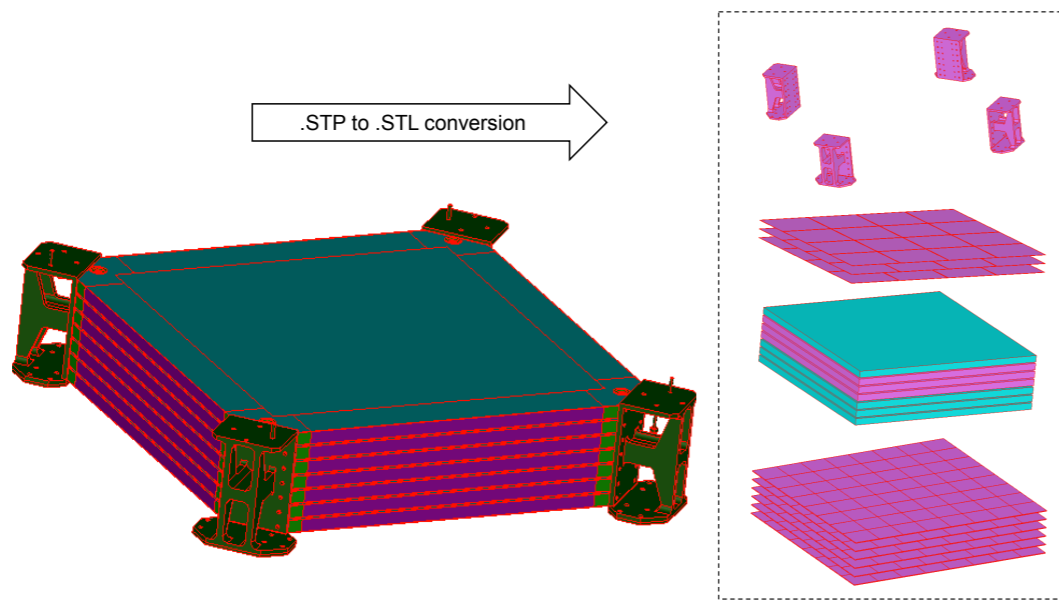


DAMPE Software framework



Geometry

- Geometry model in Geant4 is created from the CAD drawings
 - Each CAD geometry element is saved as a tessellated solid in an STL file
 - STL converted to GDML with material assignment using the conversion tool*



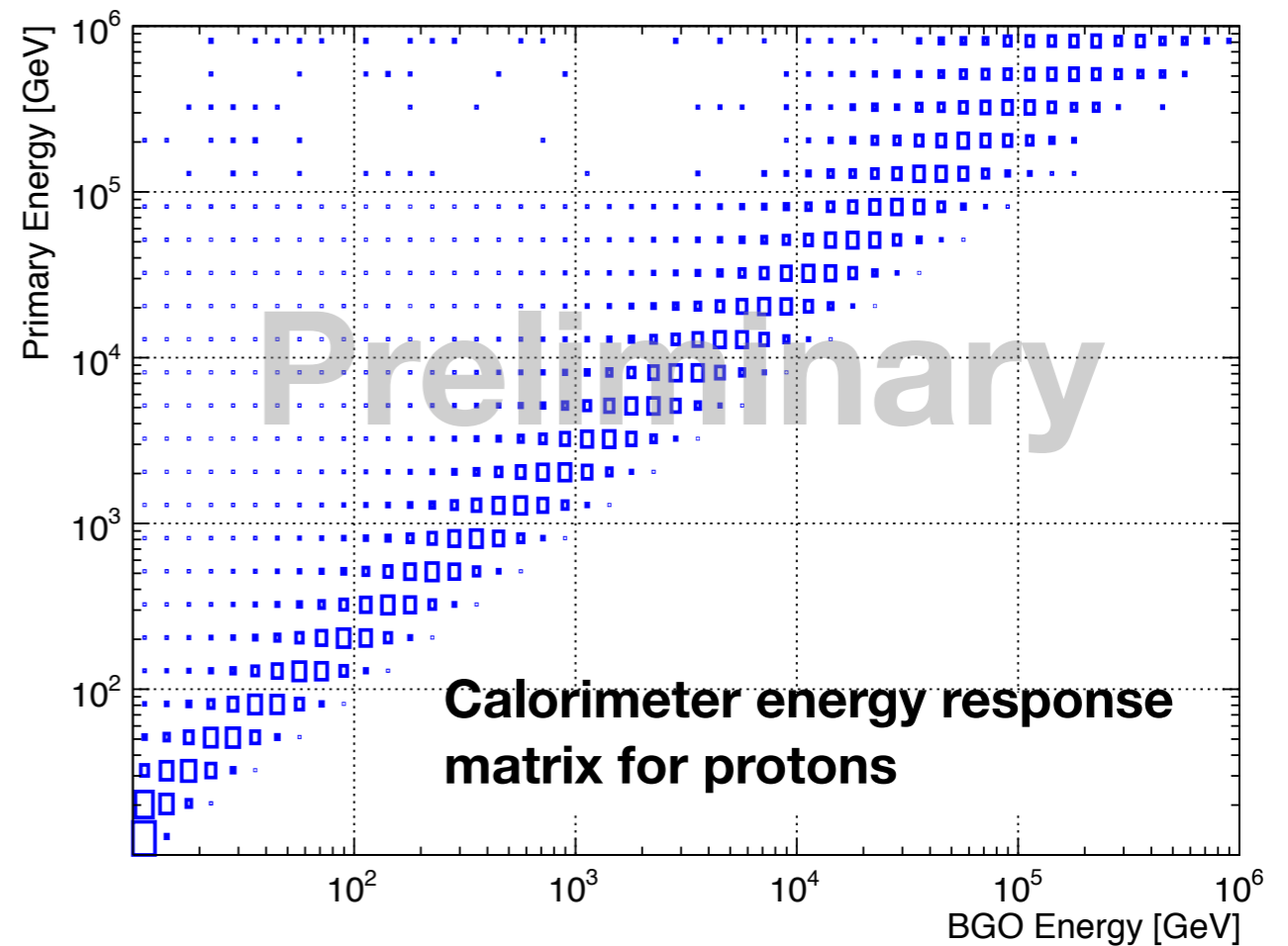
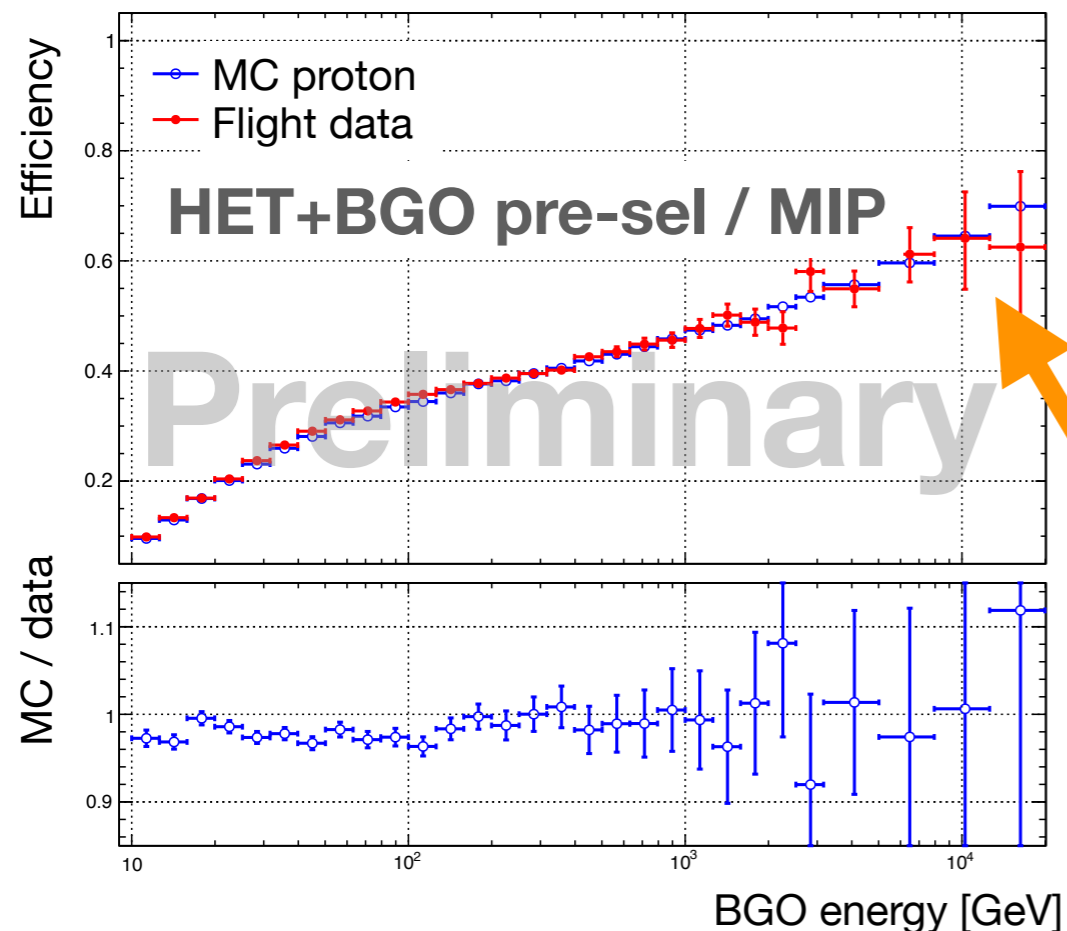
* <https://github.com/tihonav/cad-to-geant4-converter>

CR proton/ion analyses

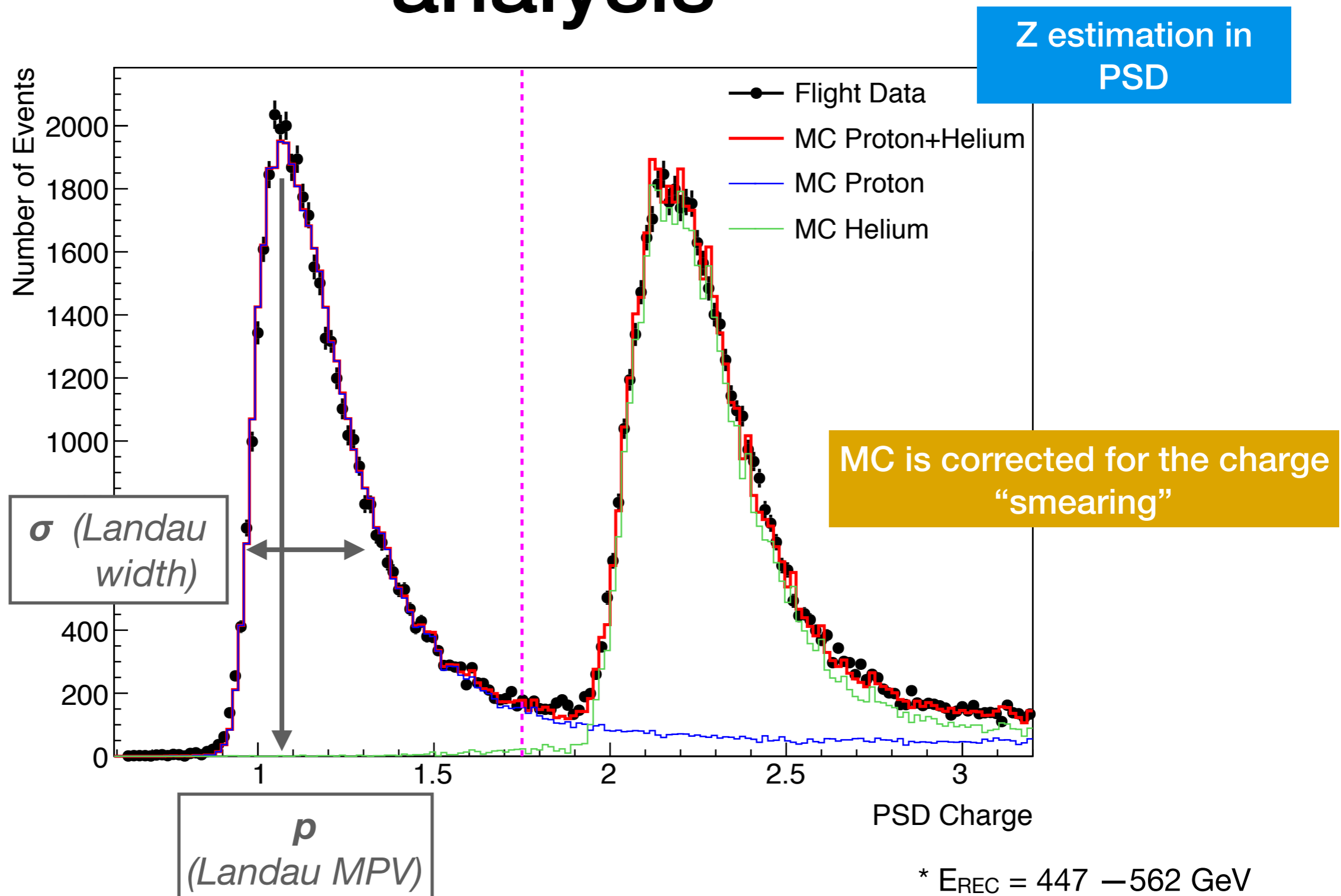
Simulation challenges

Simulation challenges in CR analysis

- High-Energy Trigger (HET) — largely determines the proton acceptance
- After the HET selection, average deposited energy fraction for protons ranges from ~50% at 100GeV to ~30% at 100 TeV:
 - **Energy unfolding is needed — relies on MC!**

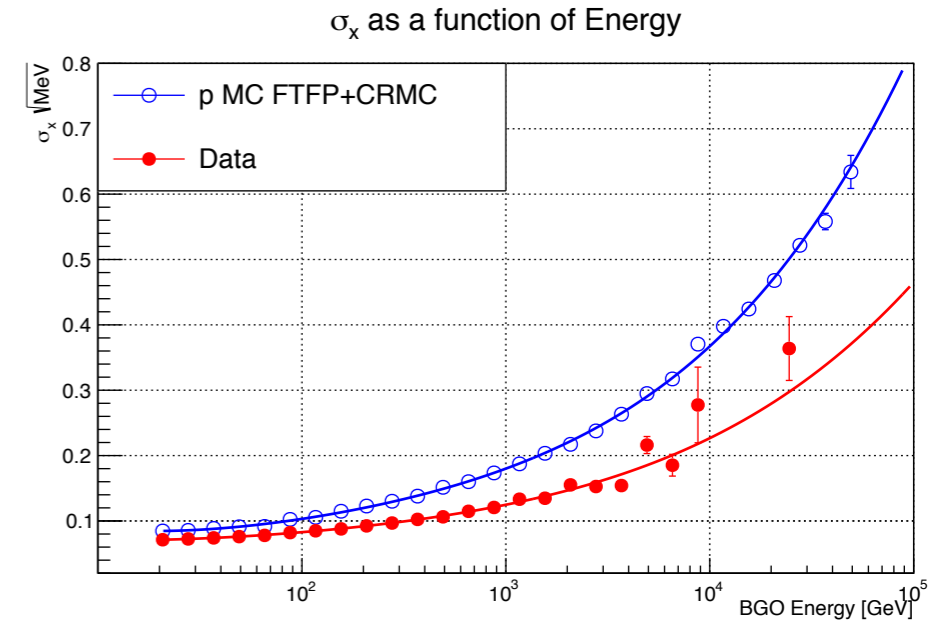
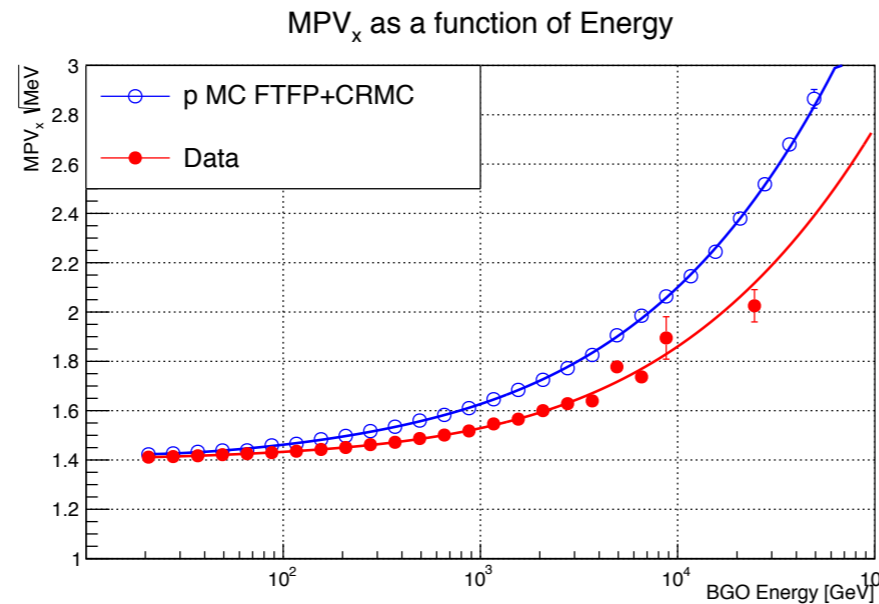


Simulation challenges in CR analysis



Simulation challenges in CR analysis

Charge
“smearing”
in MC



$$p_0 = f_{\text{MPV Data } x(y)}(E_{\text{rec}})$$

$$s_0 = f_{\sigma \text{ Data } x(y)}(E_{\text{rec}})$$

$$p_1 = f_{\text{MPV pMC } x(y)}(E_{\text{rec}})$$

$$s_1 = f_{\sigma \text{ pMC } x(y)}(E_{\text{rec}})$$

Data

$$\Rightarrow \text{PSD}_{x(y)}(\text{corr}) = (\text{PSD}_{x(y)} - p_1)s_0/s_1 + p_0.$$

MC

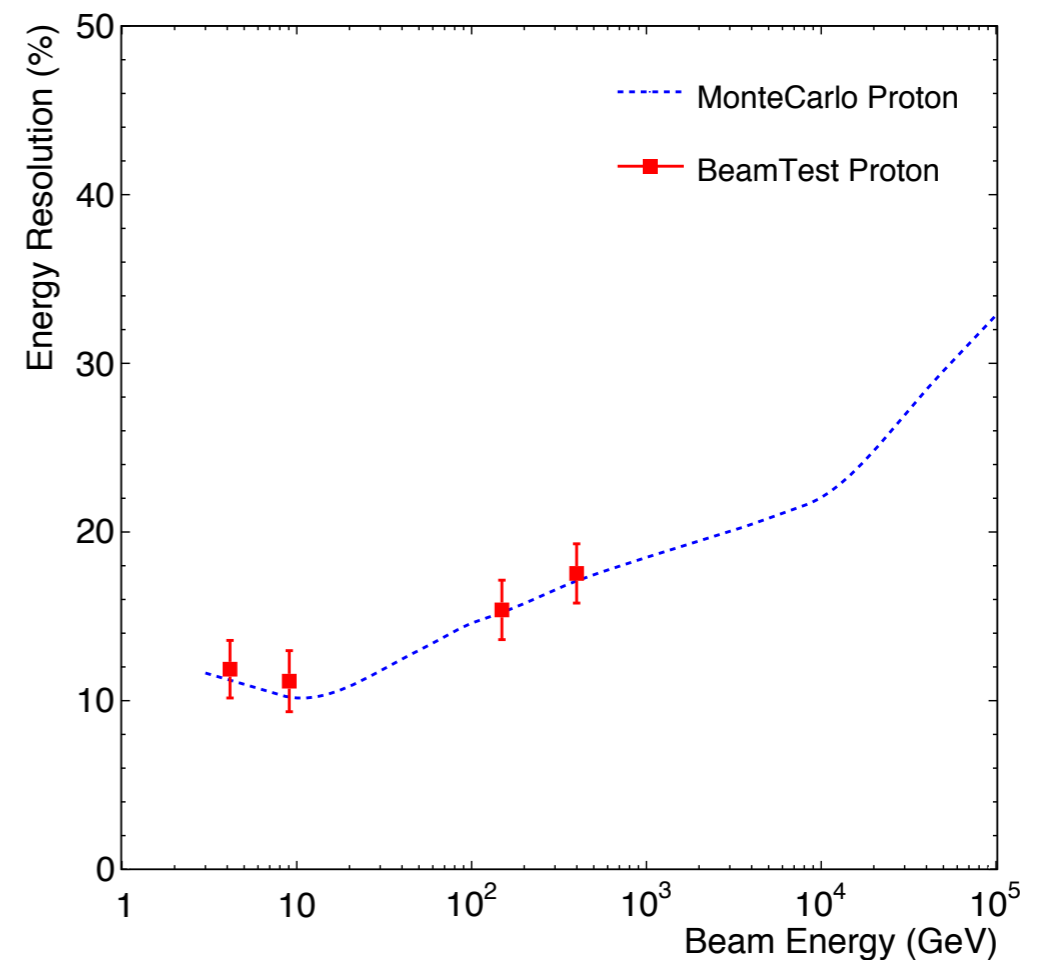
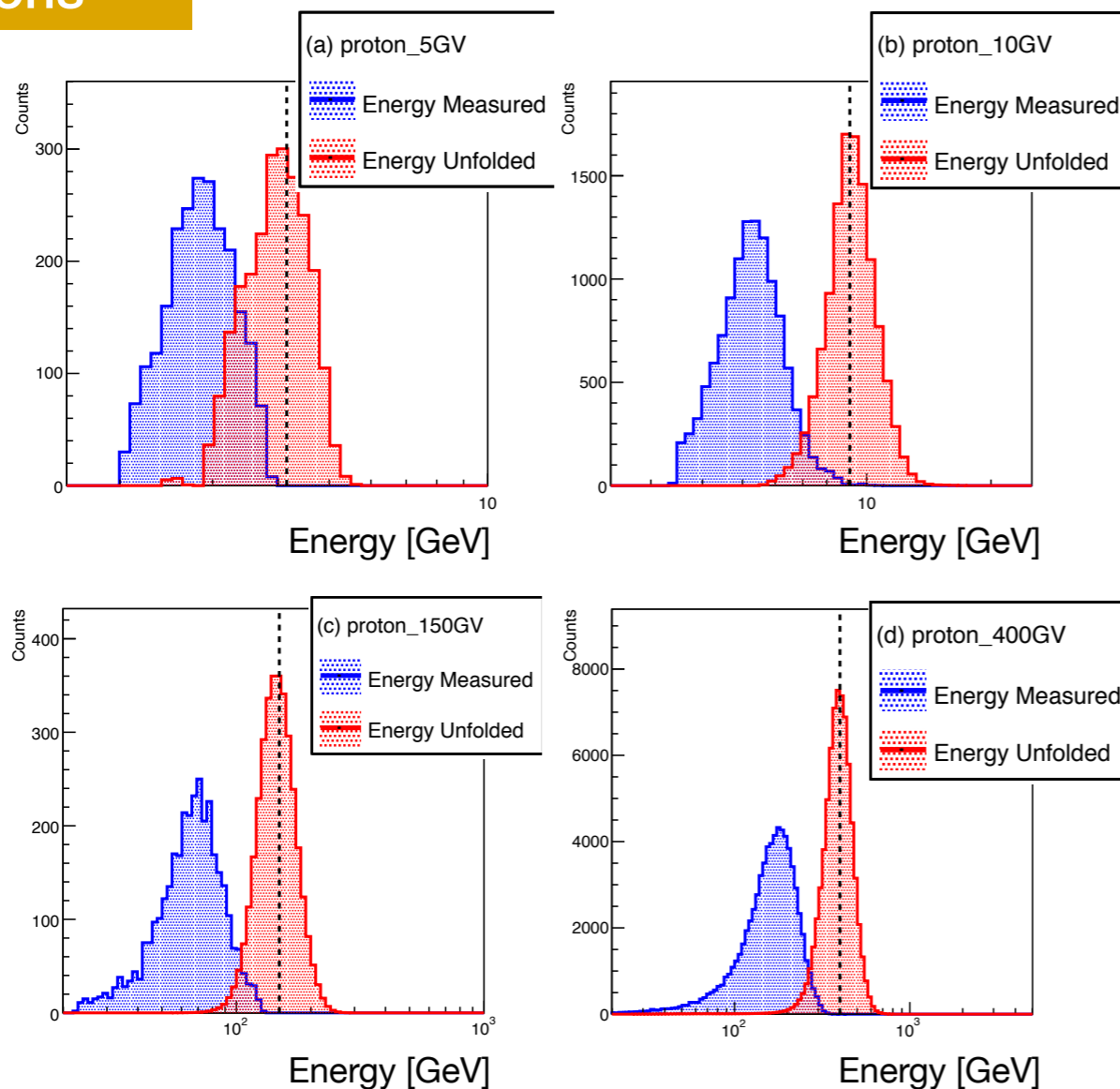
Simulation

$E > 100 \text{ TeV/nucl}$

Baseline simulation: Geant4

- Geant4 hadronic simulation shows very good agreement with beam-test data, at PS—SPS energies (up to 400 GeV for protons)

Protons

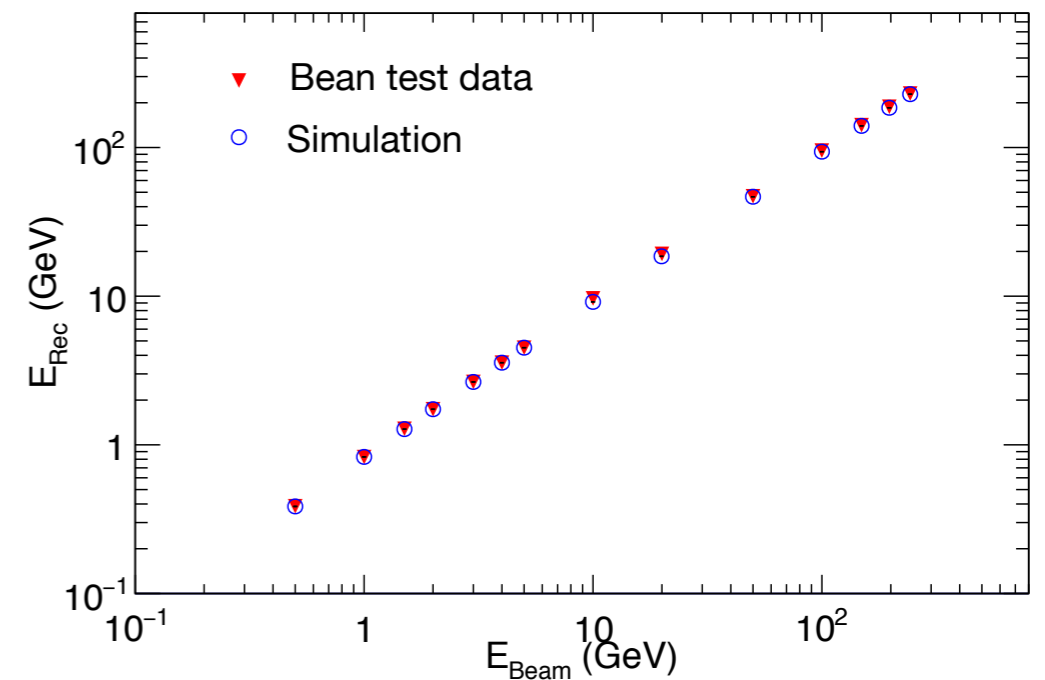
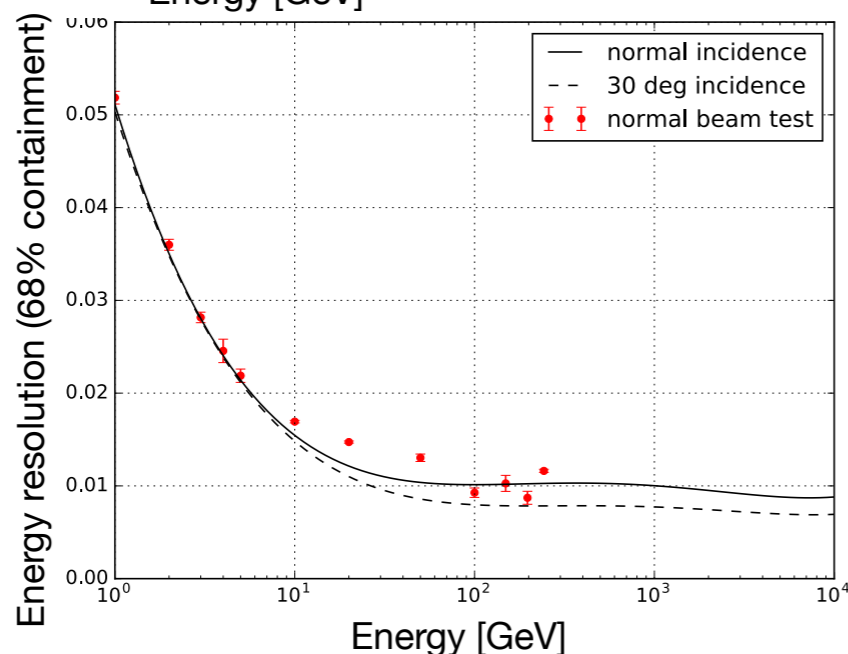
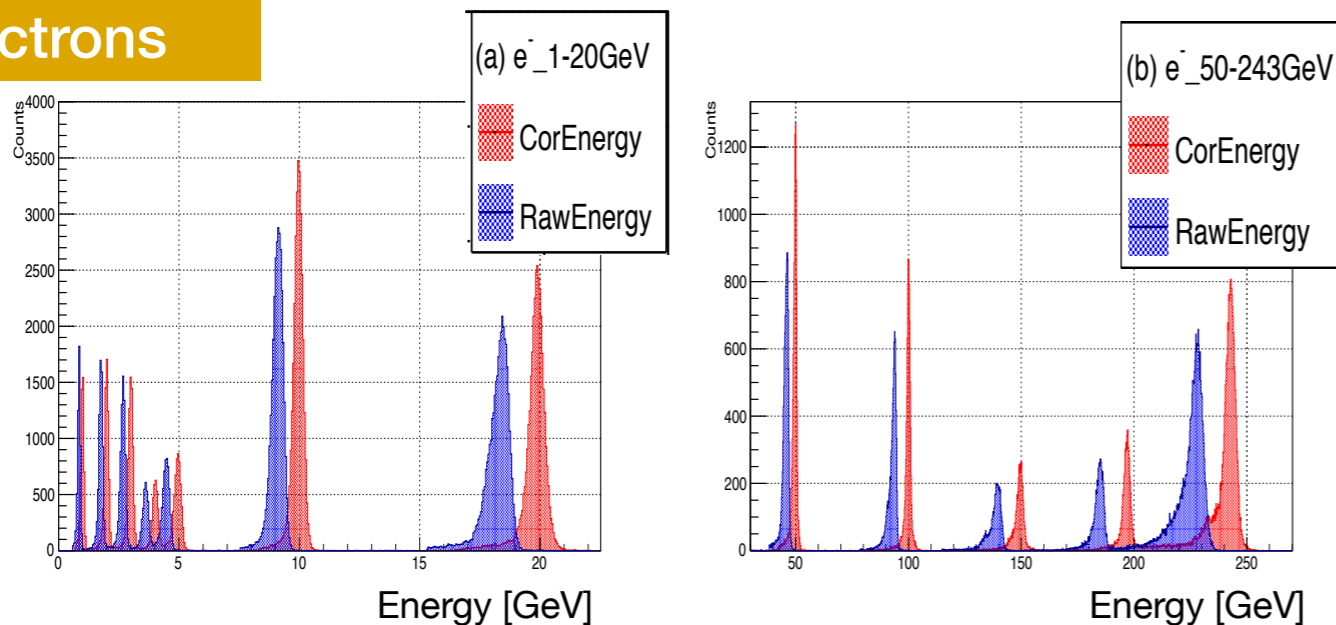


Geant4 is used as a baseline for DAMPE simulations up to 100 TeV

Baseline simulation: Geant4

- Geant4 hadronic simulation shows very good agreement with beam-test data, at PS—SPS energies (up to 243 GeV for electrons)

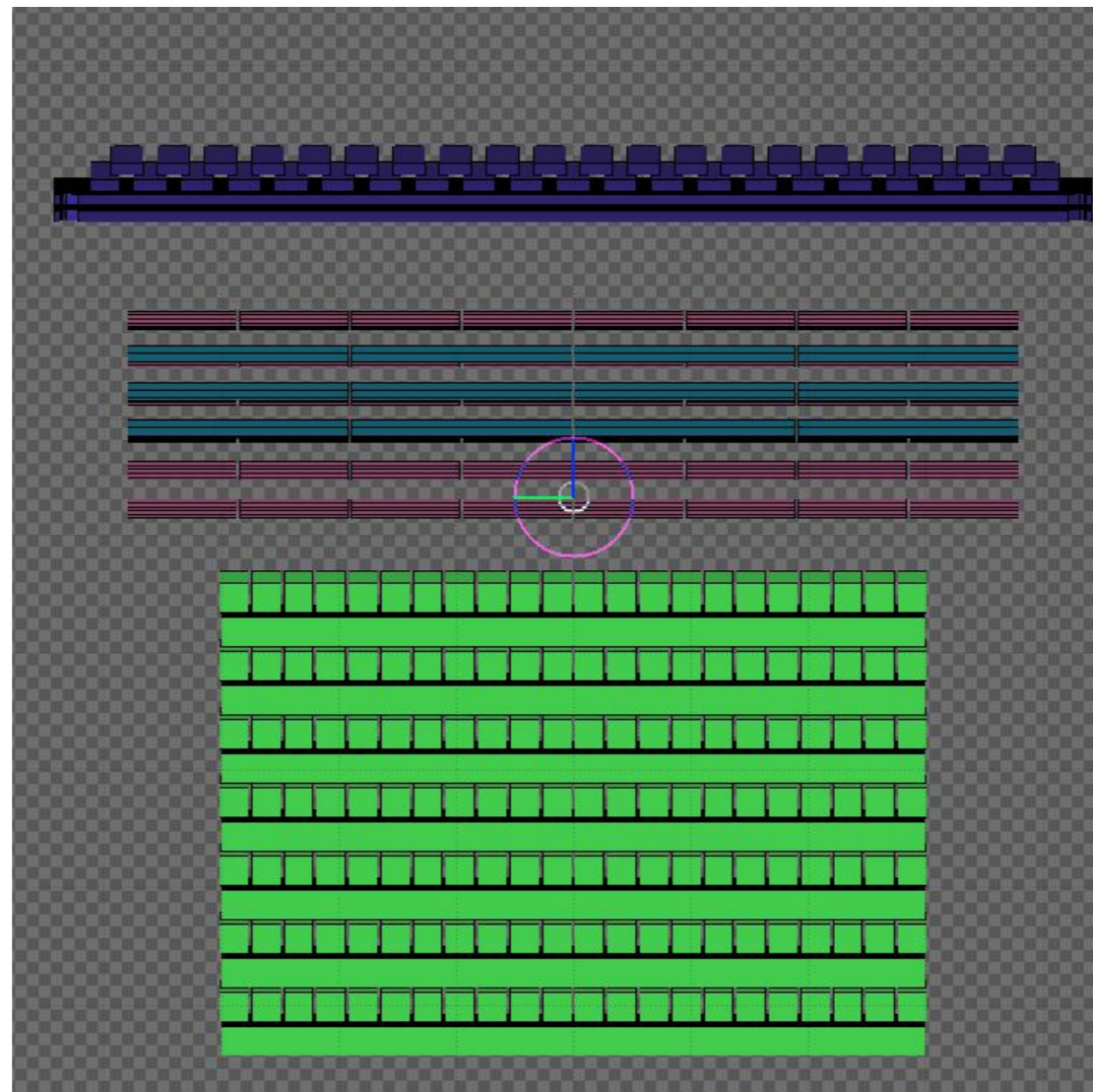
Electrons



Geant4 is used as a baseline for DAMPE simulations up to 100 TeV

FLUKA

- Alternative FLUKA simulation of DAMPE is implemented in parallel with the baseline Geant4 simulation



Geant4 – CRMC interface

- Running Geant4 simulations at > 100 TeV energies? → CRMC interface:
 - Allows to run simulations with using the DPMJET, EPOS and other models in Geant4:

```
# CRMC & Geant4 combined physics list:  
SimAlg.Set("HadronPhysicsList", "DPMJET3_FTFP_BERT")  
# ... or  
SimAlg.Set("HadronPhysicsList", "EPOS199_FTFP_BERT")  
  
# ... this one is optional:  
SimAlg.Set("CrmcEnergyThresholdGeV", "200")  
# ... default threshold = 300 GeV
```

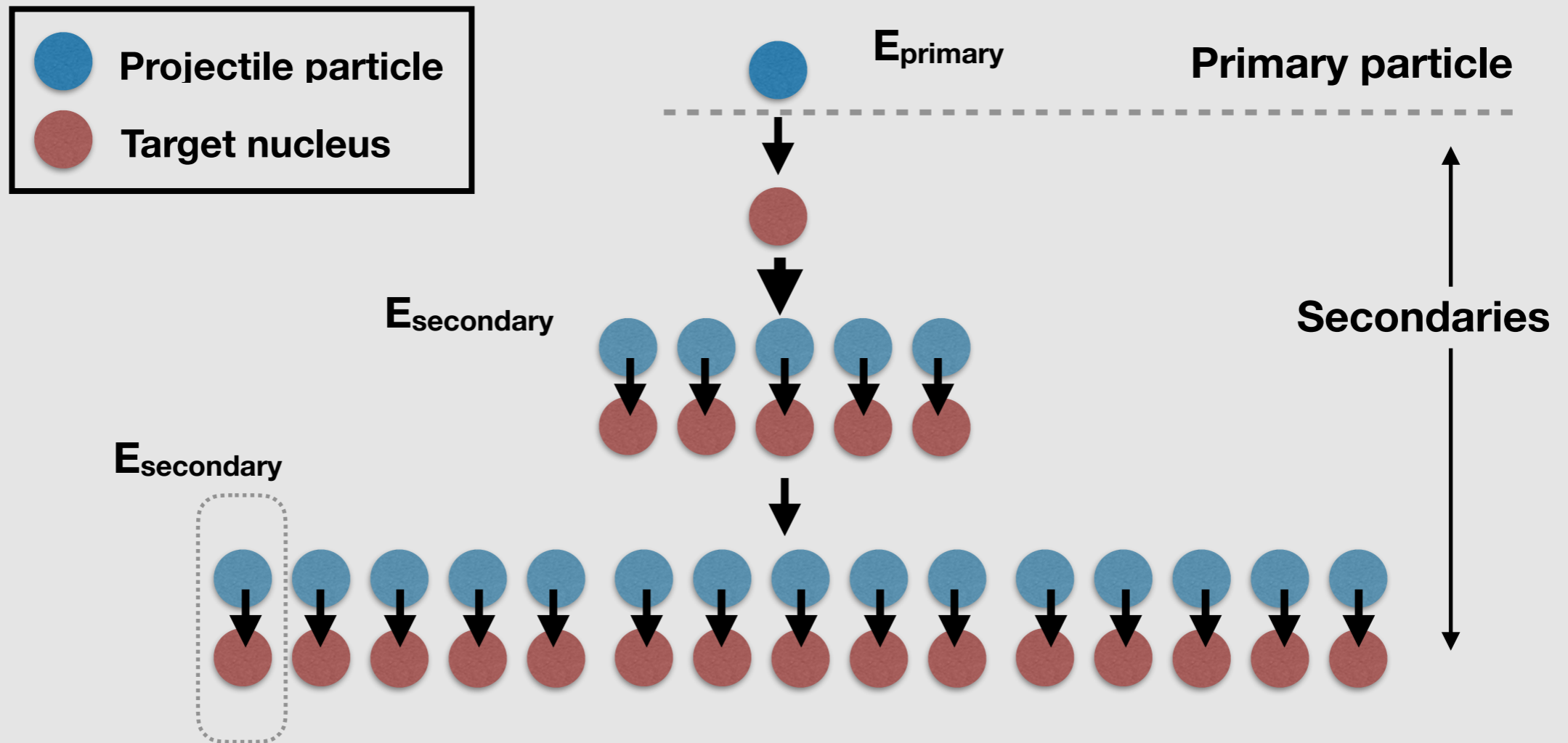


* <https://web.i kp.kit.edu/rulrich/crmc.html>

CRMC:

- DPMJET
- EPOS
- SYBILL
- ... other models

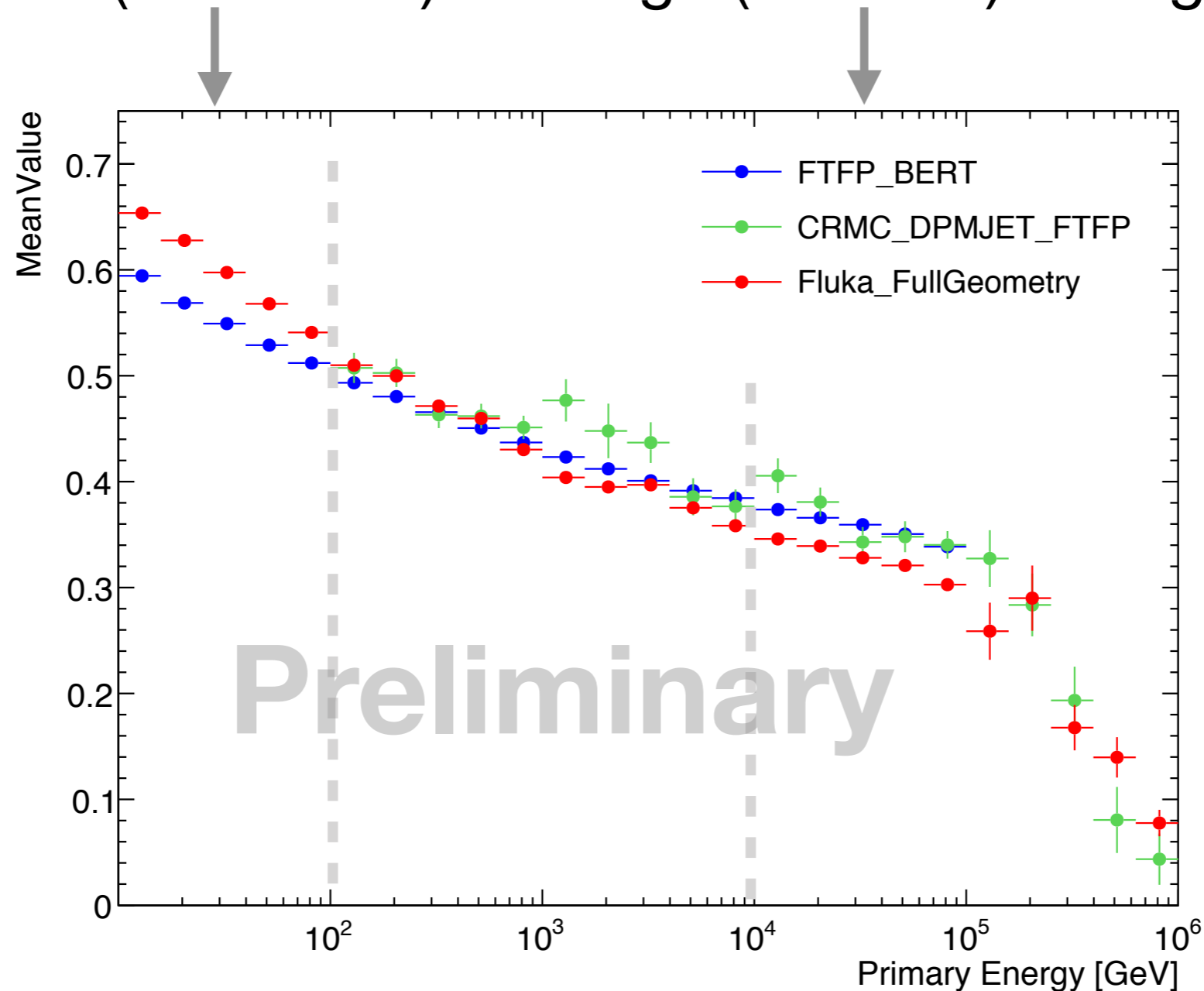
Geant4 – CRMC interface



```
If E < Ether  
    use FTFP (GEANT4)  
else  
    use CRMC (EPOS/DPMJET/..)
```

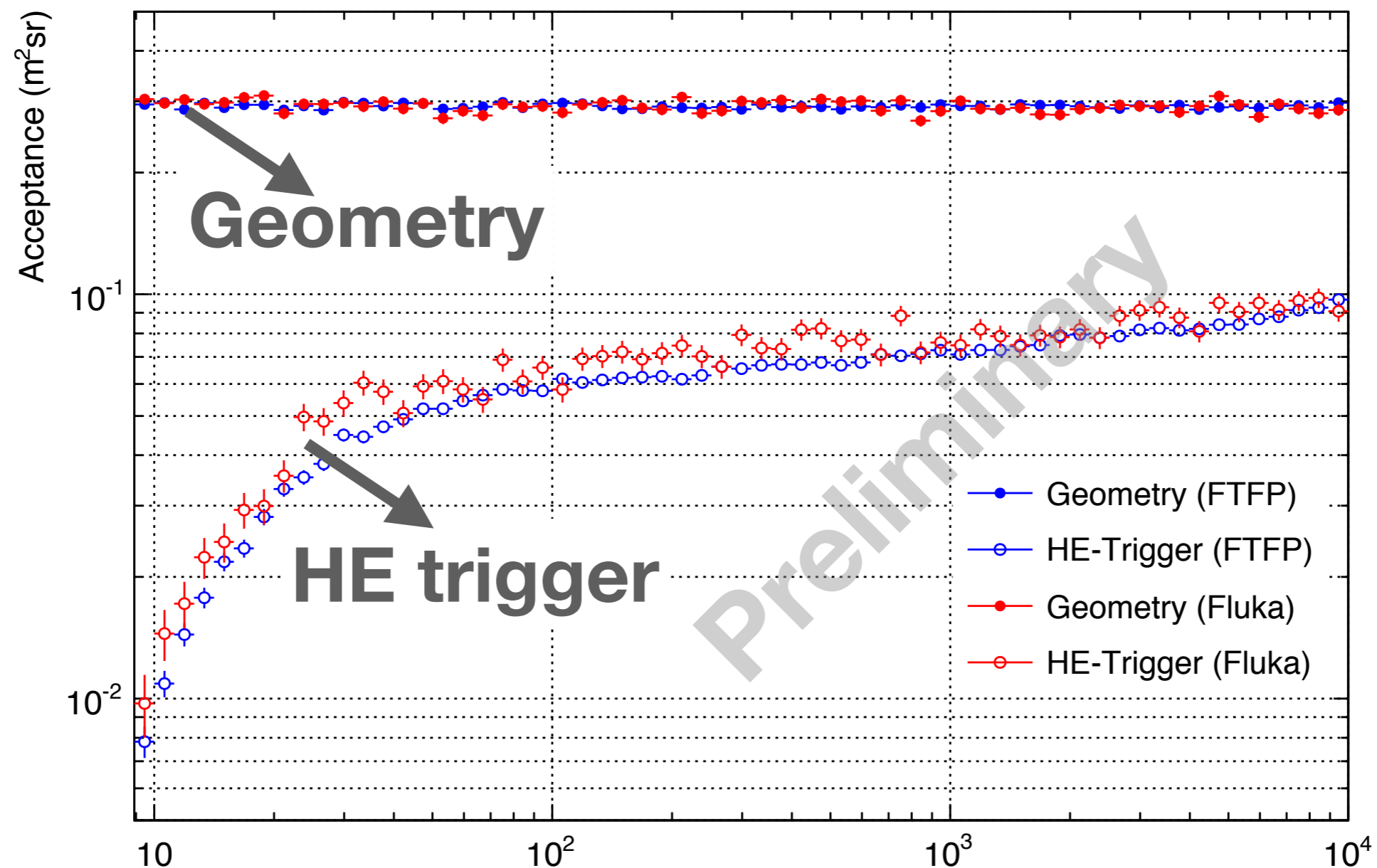
FLUKA & Geant4 – CRMC

- Geant4 and Fluka show discrepancy at low (<100 GeV) and high (>10 TeV) energies

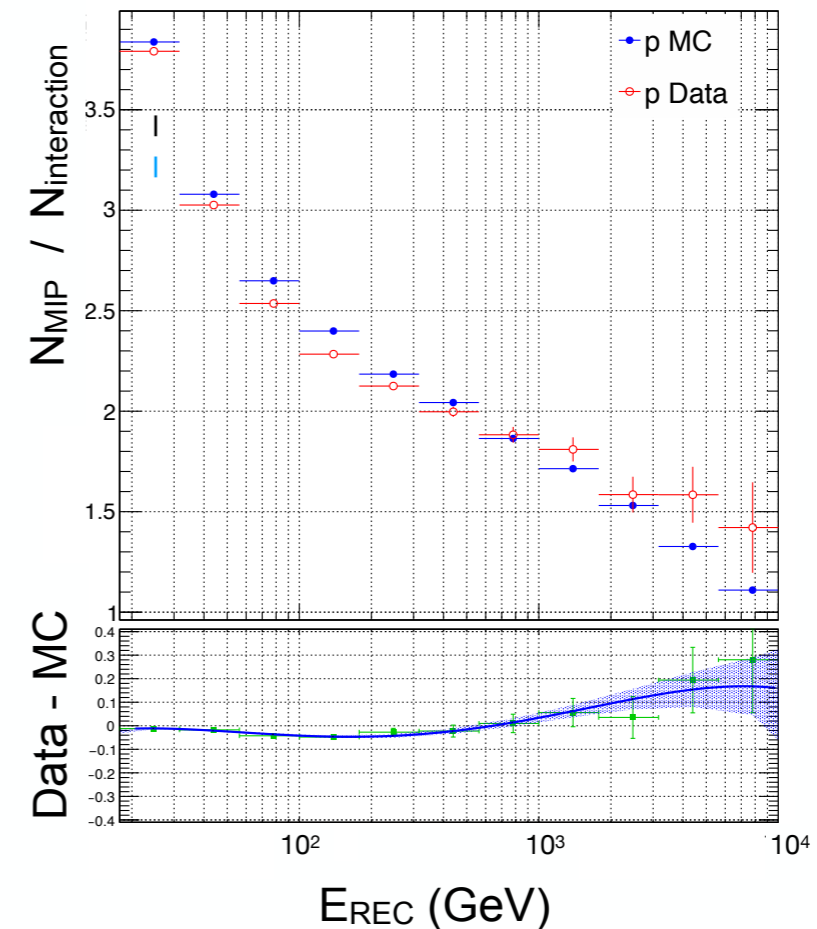
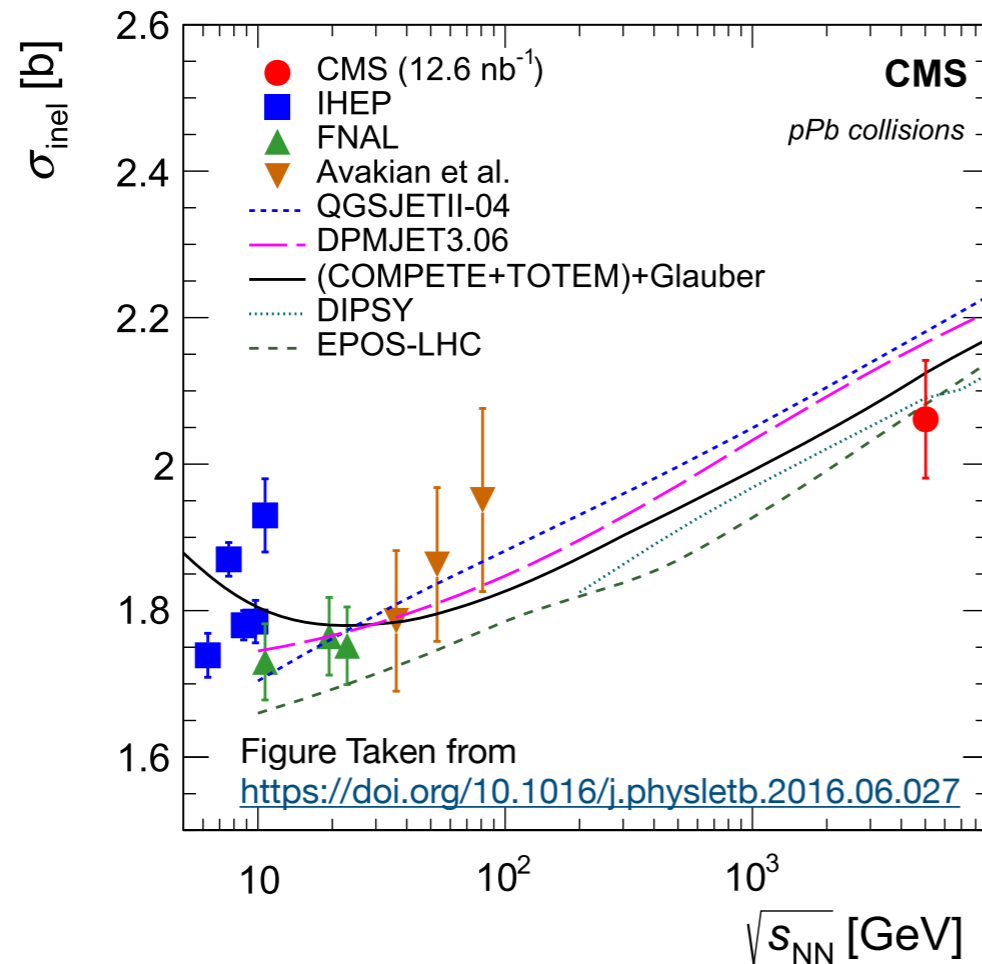


FLUKA & Geant4 – CRMC

- Geometrical acceptance is consistent between FLUKA and G4
- The difference observed for HE trigger efficiency at low energies (<100 GeV)



Hadronic cross-sections uncertainties



- Uncertainty of hadronic cross-sections in MC simulations → can be tested by varying cross-sections at Geant4 level:

```
# Modify elastic/inelastic cross-section by 10% down
SimAlg.Set("ModifyInelasticCrossSection", "0.9")

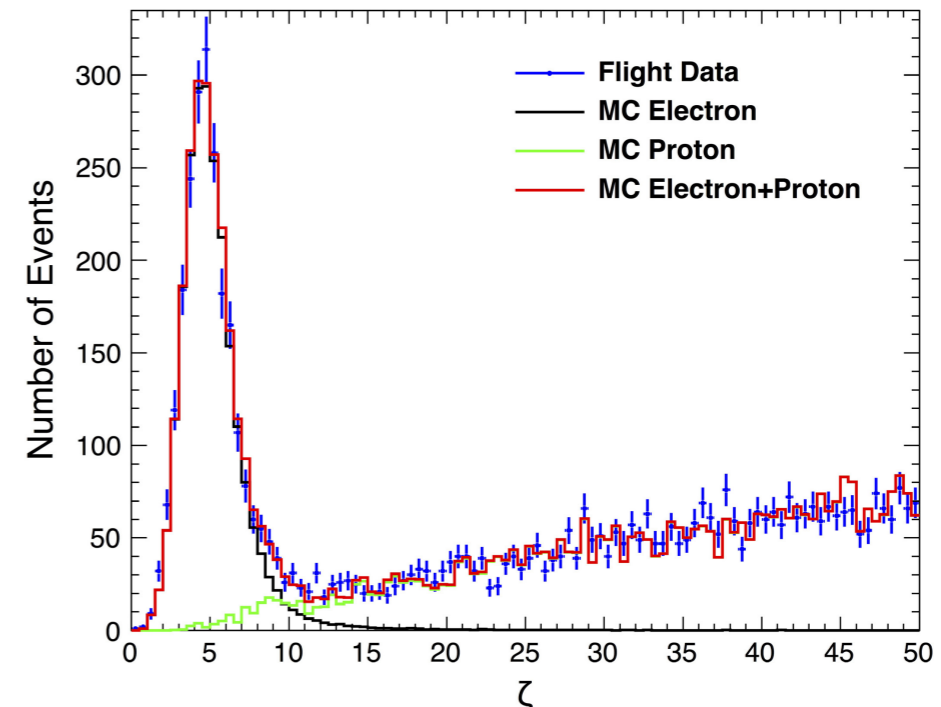
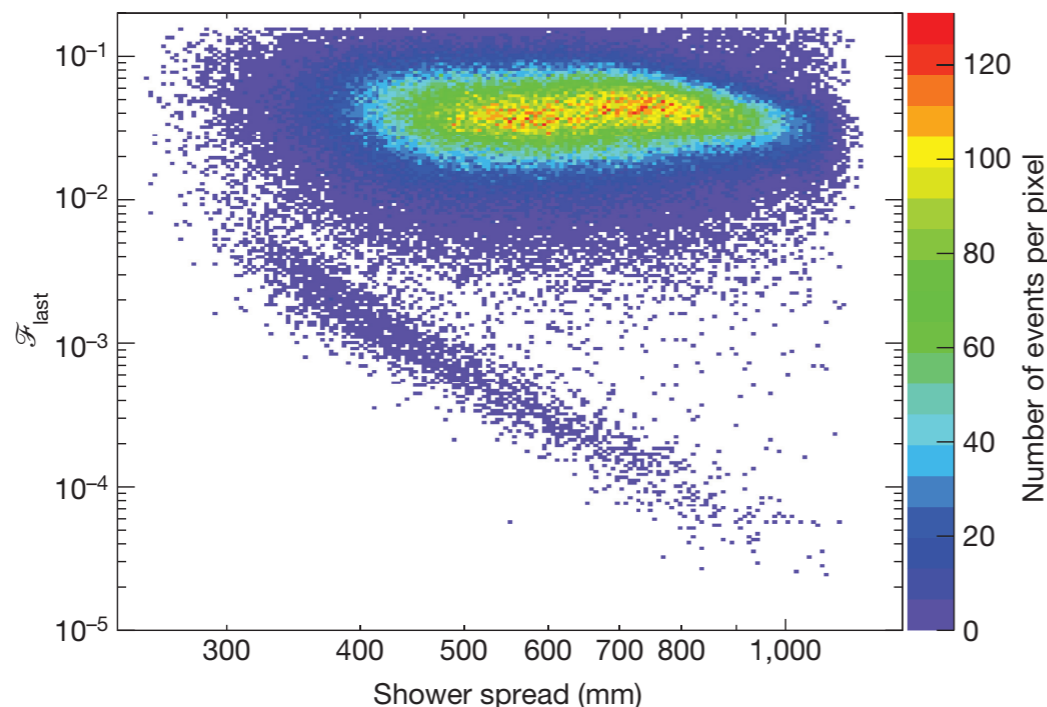
# Set minimal particle energy for which the cross-section is modified
SimAlg.Set("ModifyCrossSectionMinTotalEnergyGeV", "0")
```

CR electron analysis

Simulation challenges

Discrimination of e/p & Simulation

- Electron/proton discrimination in DAMPE is based on the shower-shape variables:
 - Sum of shower RMS in all 14 layers of the BGO calorimeter (RMS_i)
 - Fraction of energy in the last BGO layer (\mathcal{F}_{last})



$$RMS_i = \sqrt{\frac{\sum_j (x_{j,i} - x_{c,i})^2 E_{j,i}}{\sum_j E_{j,i}}}$$

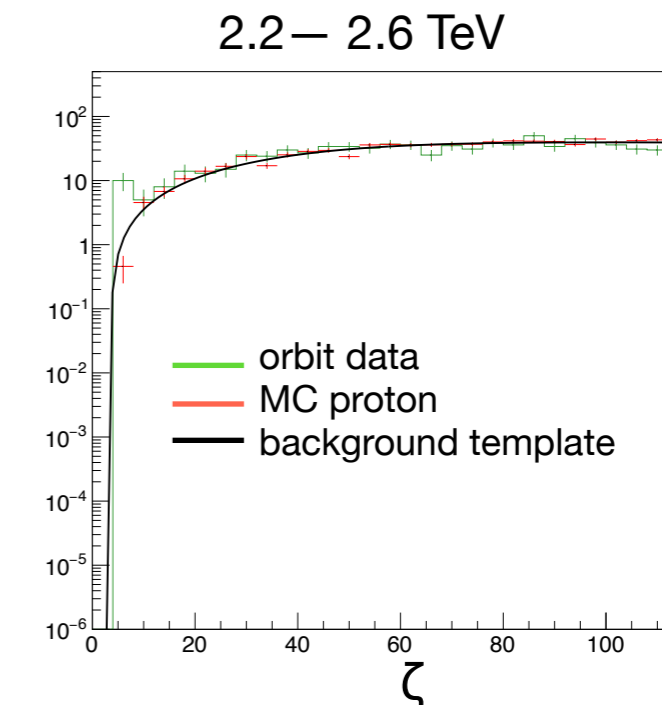
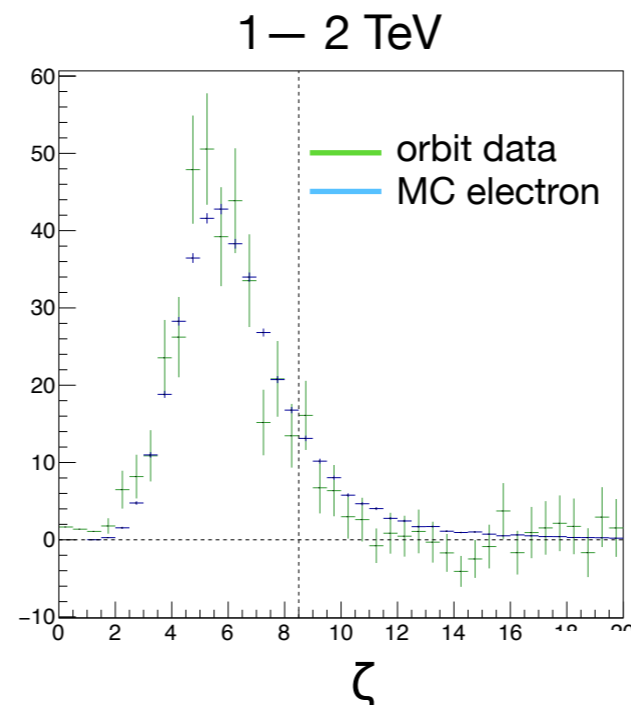
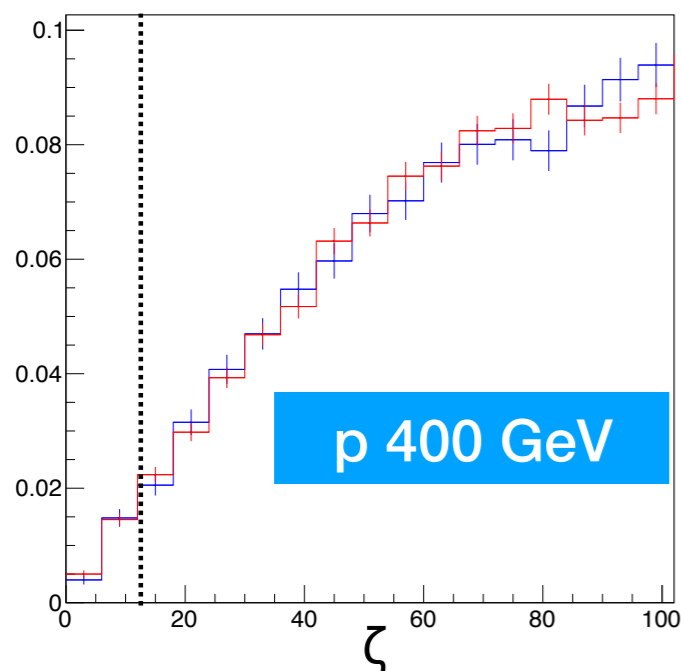
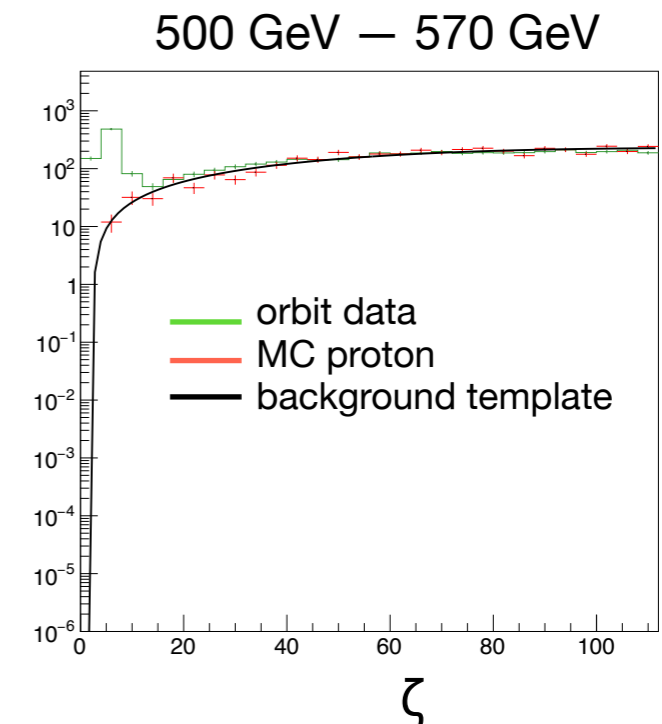
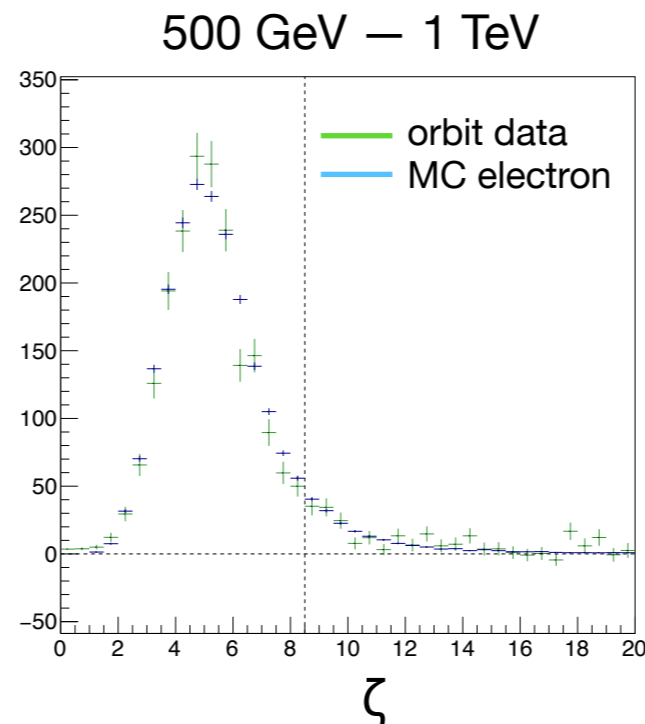
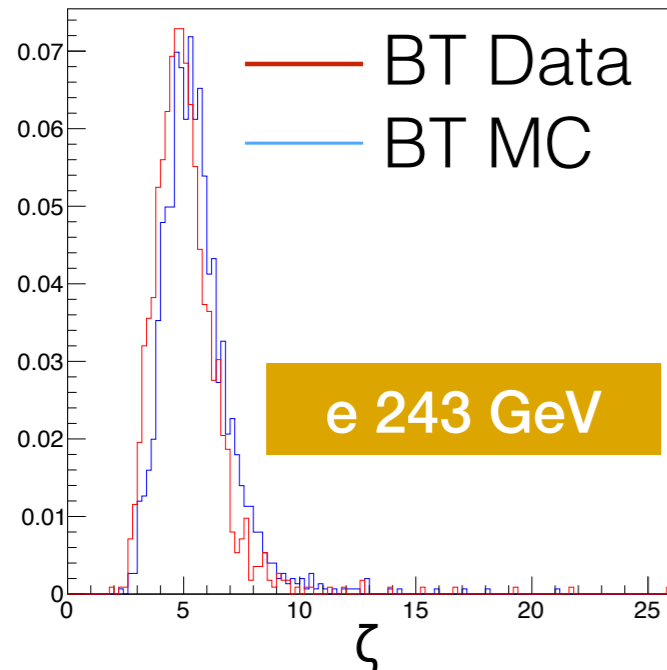


<https://doi.org/10.1038/nature24475>

$$\zeta = \mathcal{F}_{last} \times (\sum_i RMS_i / \text{mm})^4 / (8 \times 10^6)$$

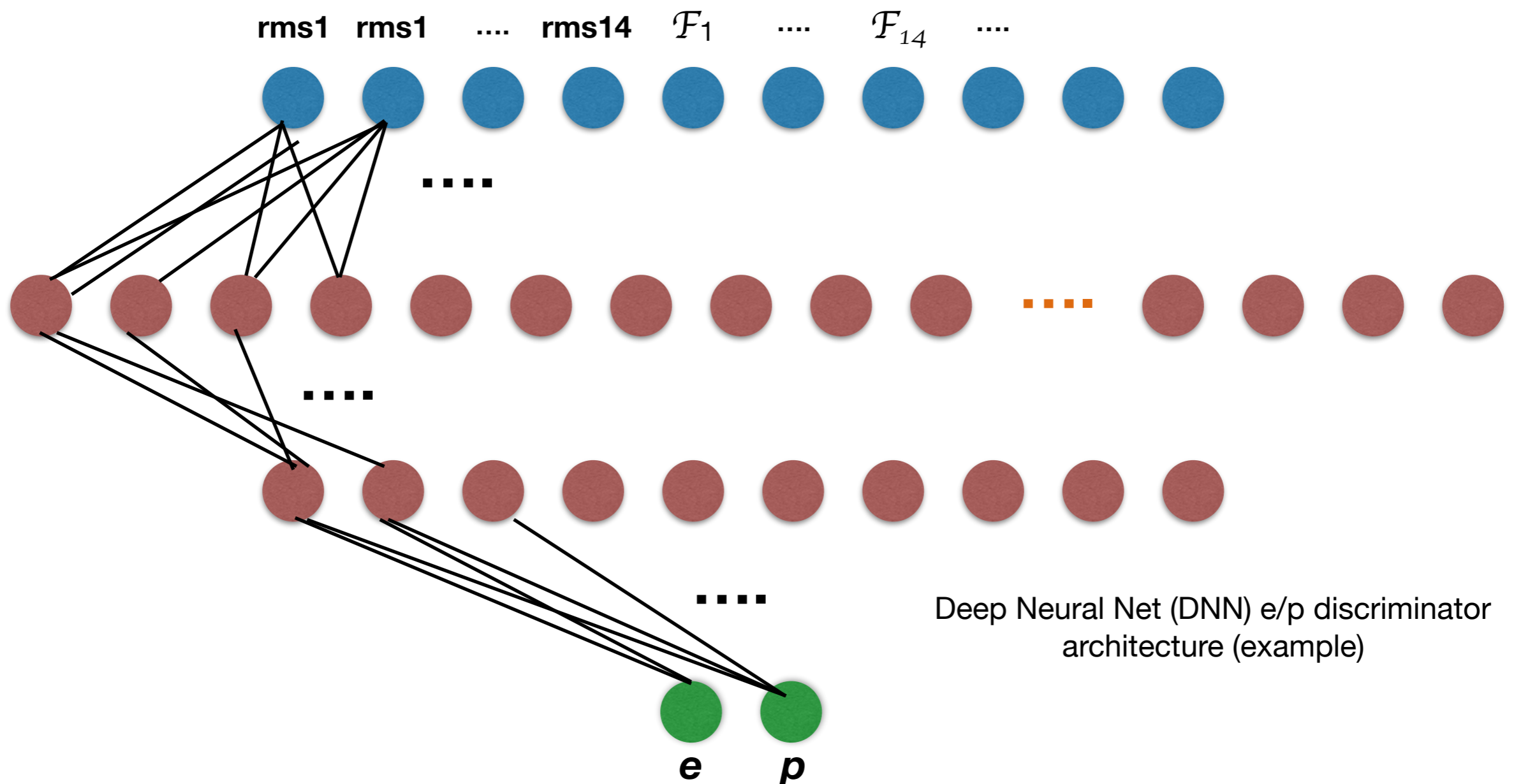
Discrimination of e/p & Simulation

- Very good data-MC agreement of e/p classifier at beam-test energies.
- Good agreement for flight data, 8% uncertainty at 4 TeV (at ~80% electron selection efficiency)



Discrimination of e/p & Simulation

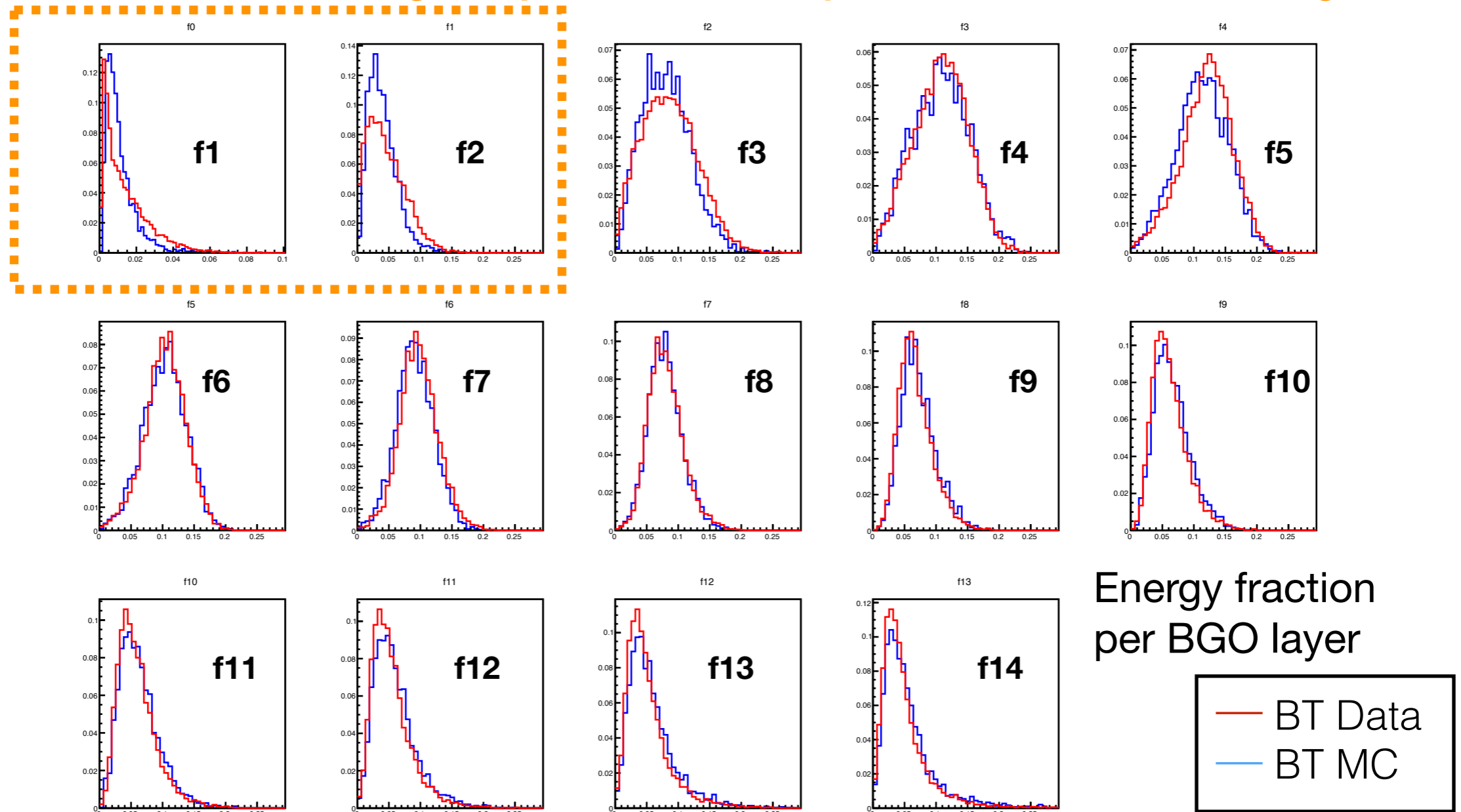
- At >5 TeV energies Deep Learning or similar approaches appear promising for e/p discrimination \rightarrow represents challenge for the simulation.



Discrimination of e/p & Simulation

- e/p classifier optimised with the simulation → relies significantly on the MC precision
- Works well with MC — does it work with the real data? How to estimate background?

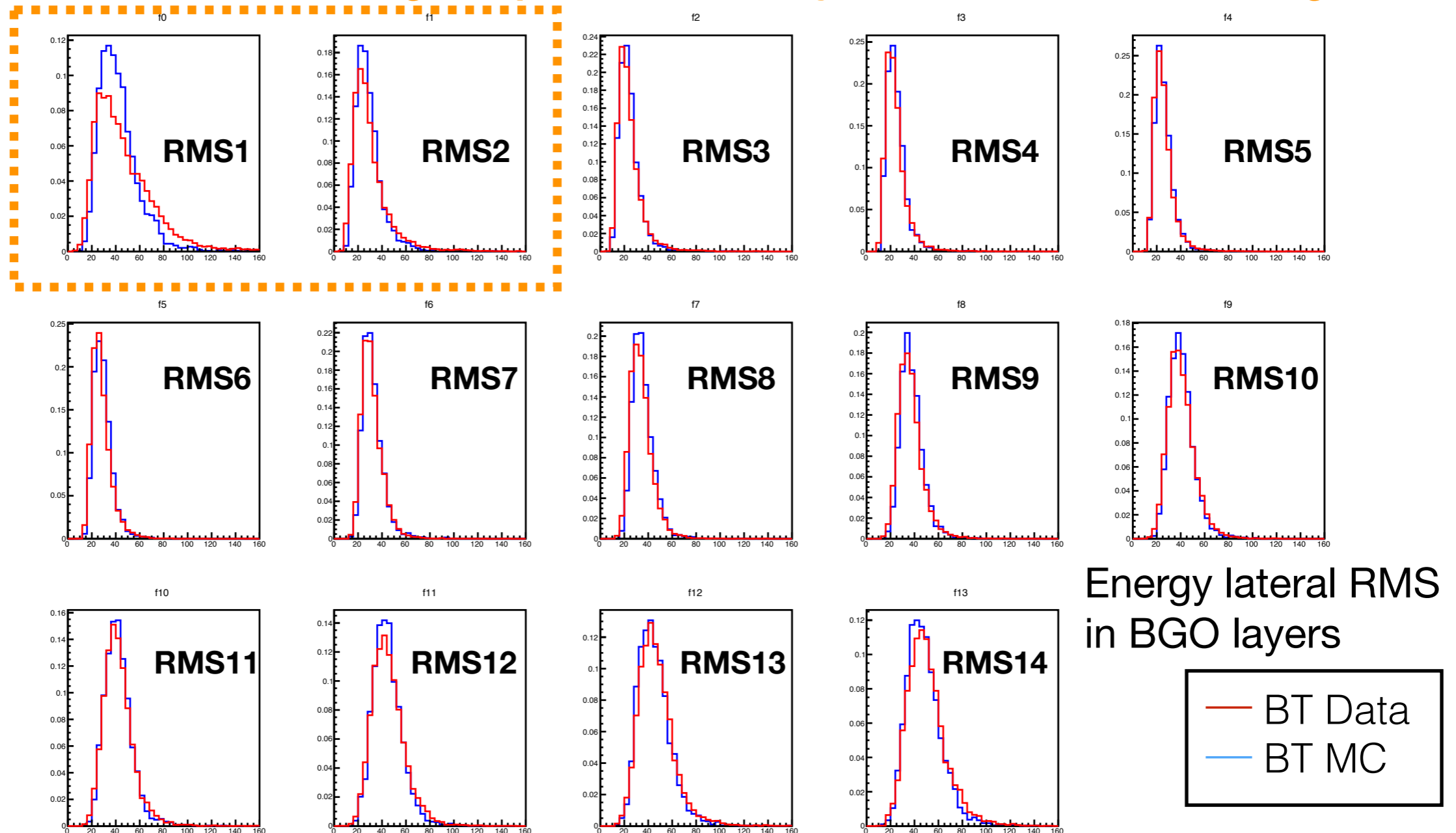
Mismodelling of input variables may throw off the data—MC agreement



Discrimination of e/p & Simulation

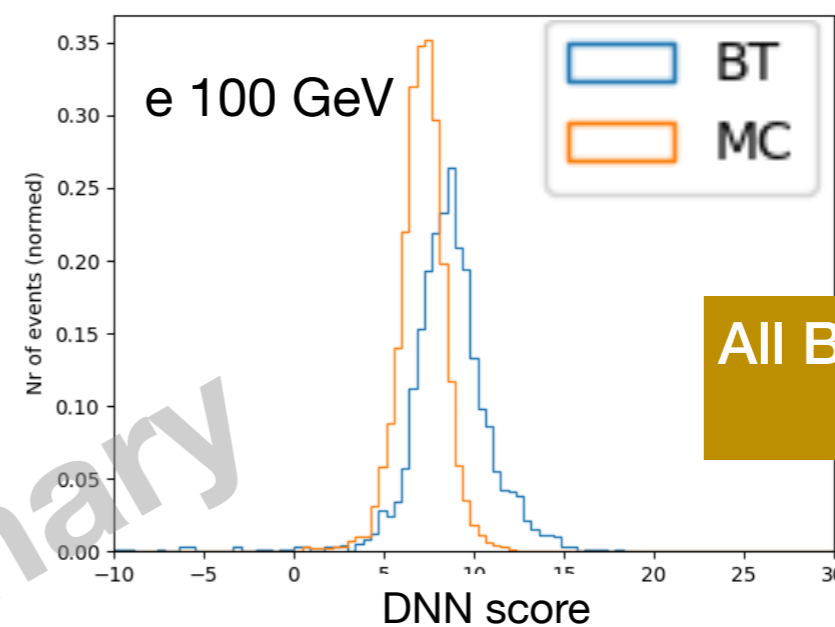
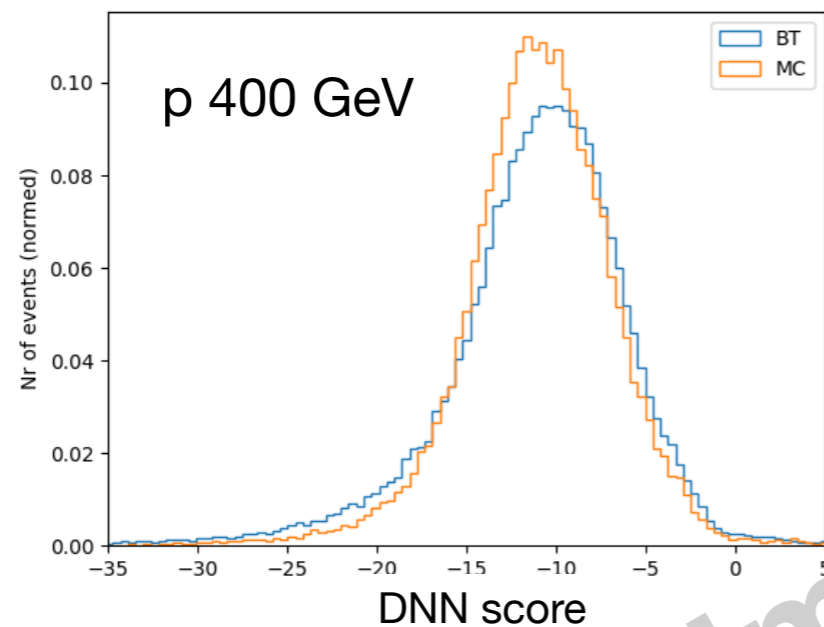
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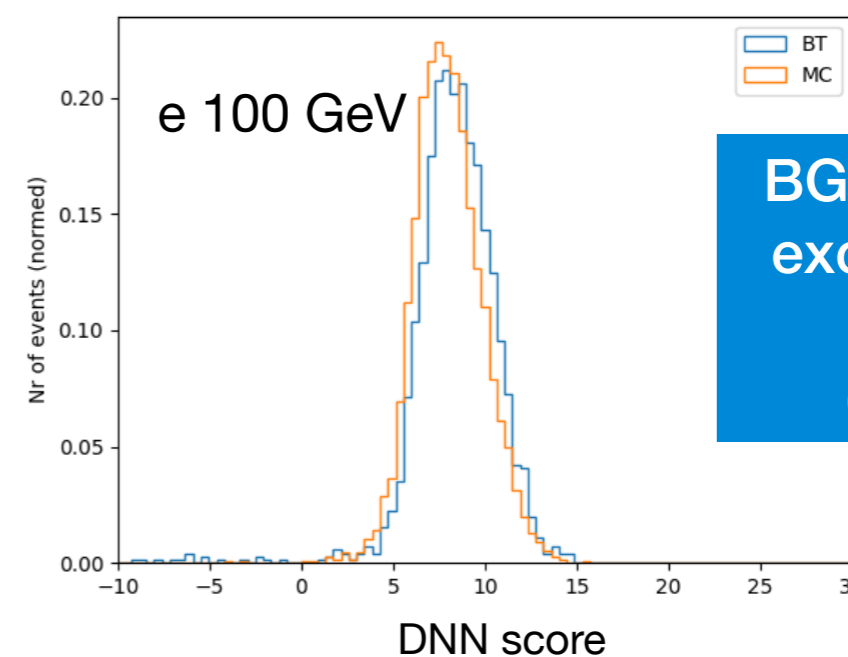
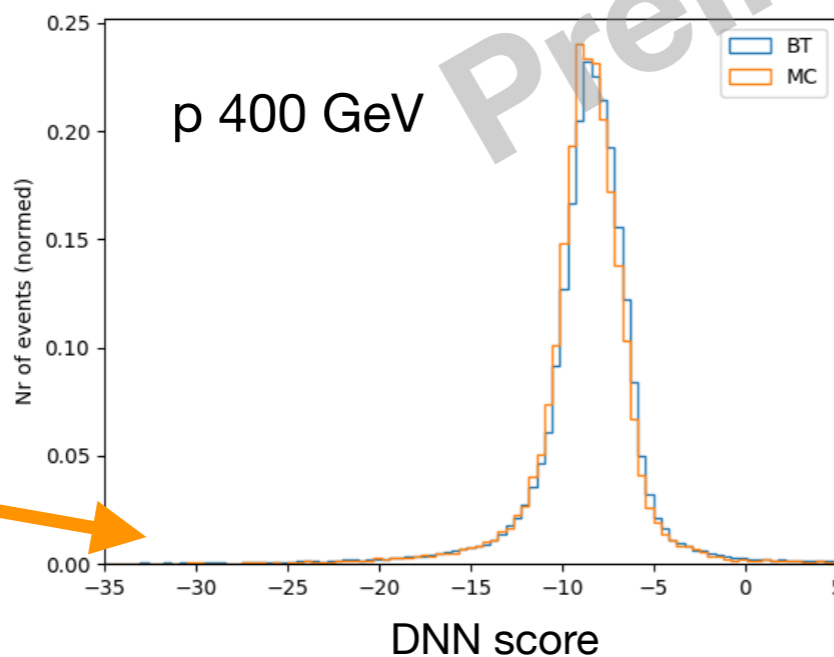
Discrimination of e/p & Simulation

- Excluding some non well-modelled variables from DNN allowed to maintain good data-MC agreement of the classifier score, while keeping (almost) the same DNN performance.



All BGO variables used

Good data – MC agreement of classifier score



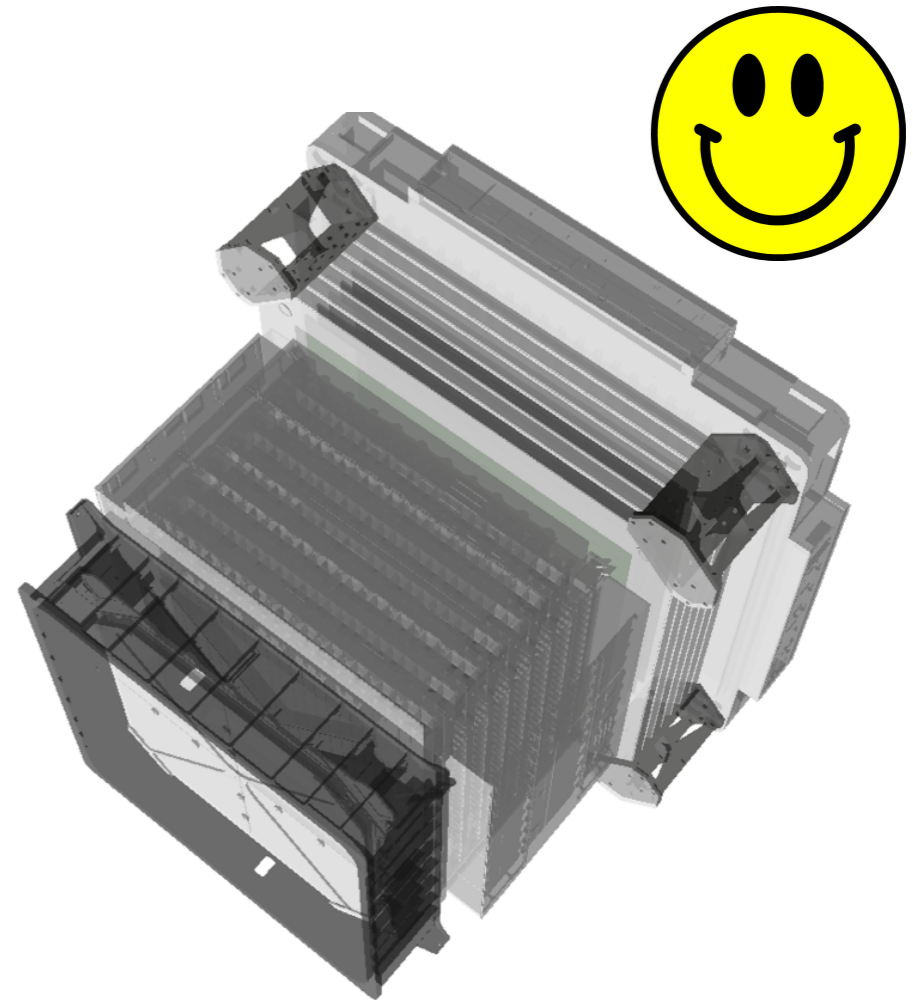
BGO layers 1,2 excluded from the DNN classifier

Conclusions

- Geant4 DAMPE simulation proven to work well: good agreement with beam-test and orbit data.
- Solutions for performing simulations above 100 TeV:
 - FLUKA
 - GEANT4 + CRMC
- Simulation in CR proton/ion analysis:
 - Energy unfolding
 - Trigger efficiency
 - Charge simulation
 - Hadronic cross-sections
- Good quality of simulation is important for CR electron analysis and e/p discrimination optimisation.

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Thank you!