Setup



<u>Objective</u>: Use clusters to distinguish between neutron/pion shower reconstruction.

```
□ (1 n + 1 \pi) / event. ---- <u>Standalone ddsim</u>

□ φ = 45°

■ θ<sub>n</sub> = 155° (η = -1.51) ----- <u>fixed</u>

■ θ<sub>π</sub> = 155° (η = -1.51), 158° (η = -1.64),

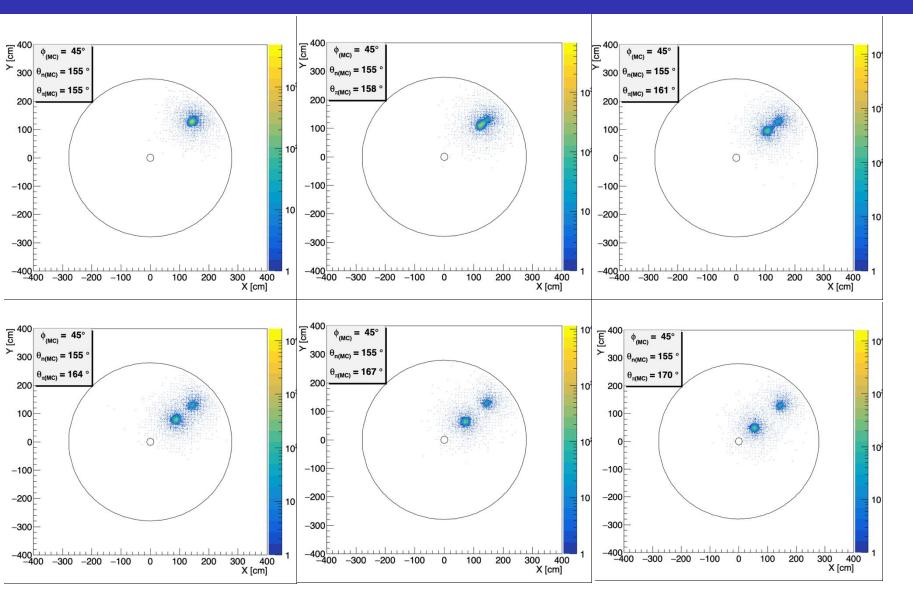
161° (η = -1.79), 164° (η = -1.96),

167° (η = -2.17), 170° (η = -2.44)
```

- Only Backward HCal was taken into account [not the whole ePIC geometry – scattering effects neglected]
- $-4.14 < \eta < -1.18$
- Alternating Steel and Scintilator slices
- 10 cm. x 10 cm. Polystyrene tiles

Cluster Positions (xy coordinates)





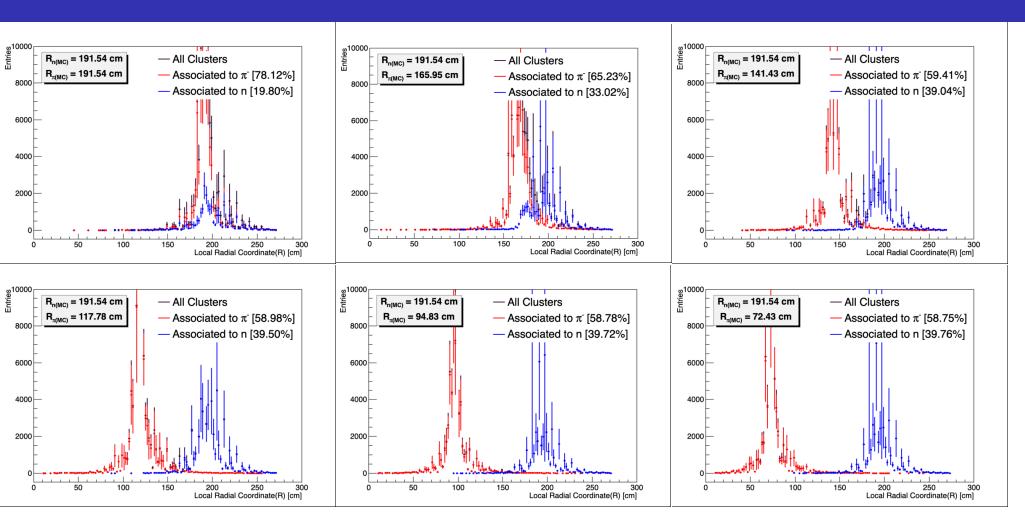
Cluster (x,y) are shown along with simulated angular coordinates

p = 1 GeV/c

[neutron showers in outer region; pion showers in inner region]

Distributions are becoming more smeared more distinguishable as $(\theta_{\pi} - \theta_{\rm n})$ increases...

Cluster Radial Coordinates





p = 1 GeV/c

~ 80% of the clusters associated with pions

~ 20% of the clusters associated with neutrons

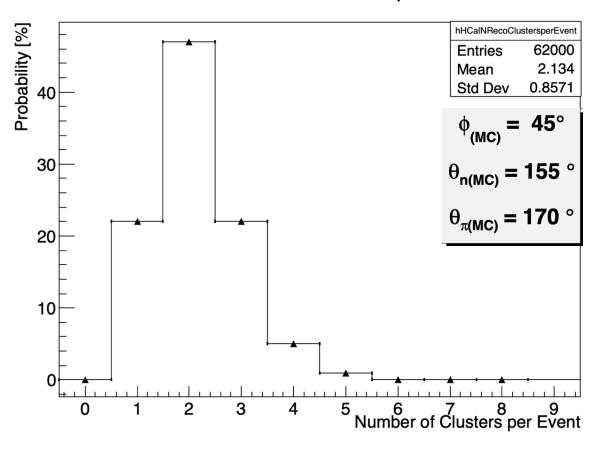
Neutron Clusters start to shift inwards as $(R_n - R_{\pi})$ increases

Percentages (fraction of clusters identified as π -/n clusters) are based on ClusterMCParticle associations [better performance as the π -n distance increases]

Clusters are in the expected positions



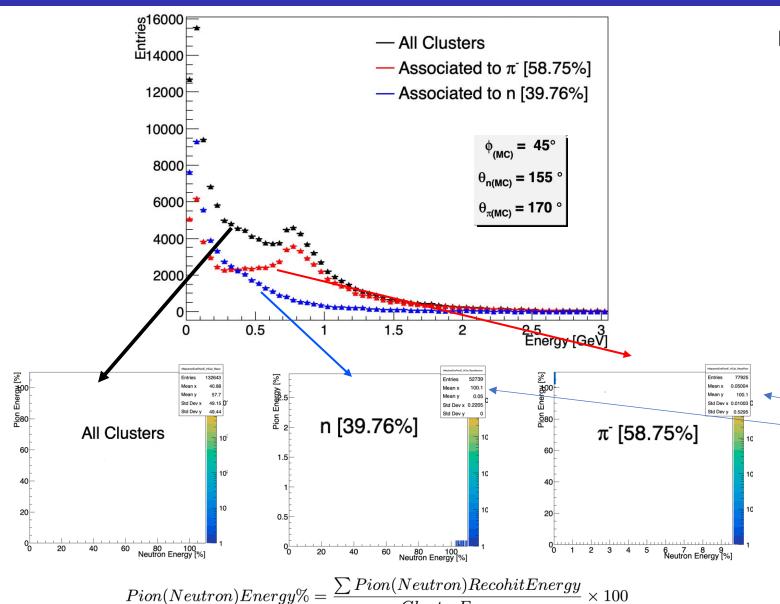
Number of Reco Clusters per Event



Expectation: 2 clusters per event

Observation: > 90%
events have 1
reconstructed cluster
Mean = 2.134

Clusters are being merged...?
Better Performance.



ClusterEnergy

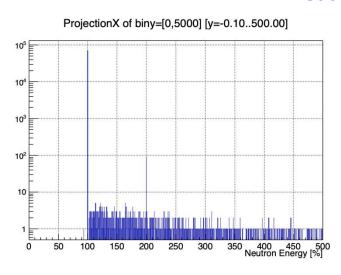
Delivation of the mergedRecoHits

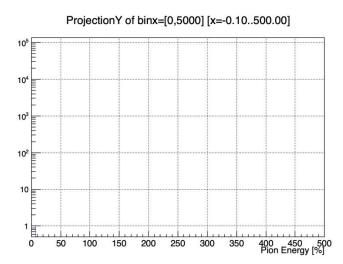
Cluster energies have been traced back to the constituent recoHits

mergedRecoHits [can be accessed via cluster.getHits()] which were tagged as pion/neutron hits based on the most energetic hit contribution of the mapped simHit [can be accessed by comparing cellIDs] of the mergedRecoHits.

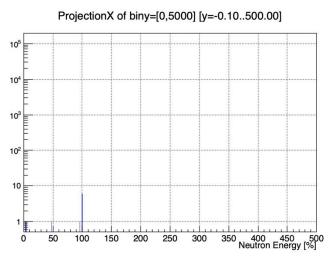
Cluster-MCparticle Association works well.

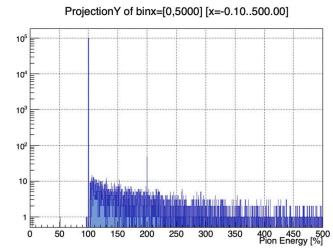
Neutron Clusters



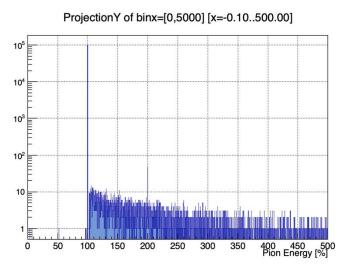


Pion Clusters

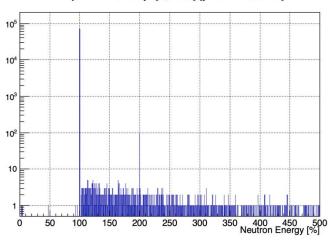




All Clusters



ProjectionX of biny=[0,5000] [y=-0.10..500.00]



^{*}discussion on next slides with an example



```
Processing event 84959/85000... # Event containing 2 clusters
```

```
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (0, 0, 16, 59, 3)
Cluster: energy = 0.721966. nHits = 6
Rec Hit: energy = 0.0263415, time = 17.3, pos.z() = -4107.5
Sim Hit: energy = 0.000220031
Contrib: energy = 0.000220031, time = 17.303
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (0, 1, 24, 89, 3)
Cluster: energy = 0.721966, nHits = 6
Rec Hit: energy = 0.132992, time = 14.29, pos.z() = -4107.5
Sim Hit: energy = 0.00127383
Contrib: energy = 0.00121976, time = 14.2872
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (duster, reco hit, sim hit, contrib, particle) ID = (0, 2, 23, 84, 3)
Cluster: energy = 0.721966, nHits = 6
Rec Hit: energy = 0.462582, time = 13.83, pos.z() = -4107.5
Sim Hit: energy = 0.00295112
Contrib: energy = 0.000721806. time = 14.039
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (0, 3, 30, 102, 3)
Cluster: energy = 0.721966, nHits = 6
Rec Hit: energy = 0.0424034, time = 14.62, pos.z() = -4107.5
Sim Hit: energy = 0.000395054
Contrib: energy = 3.9479e-06, time = 68.6456
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (duster, reco hit, sim hit, contrib, particle) ID = (0, 4, 35, 132, 3)
Cluster: energy = 0.721966, nHits = 6
Rec Hit: energy = 0.0841643, time = 14.22, pos.z() = -4107.5
Sim Hit: energy = 0.000807706
Contrib: energy = 0.000807706, time = 14.2187
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (0, 5, 37, 134, 3)
Cluster: energy = 0.721966, nHits = 6
Rec Hit: energy = 0.090589, time = 14.8, pos.z() = -4107.5
Sim Hit: energy = 0.000865865
Contrib: energy = 0.00016734, time = 14.7966
Particle: energy = 1.00969, pdg = -211, gen status = 1
```

Cluster 1:

cluster.getEnergy(): 0.721966 hcalreco neutronE: 0 hcalreco pionE: 0.839073

contribution

Pion energy > cluster energy

```
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (1, 0, 16, 59, 3)
Cluster: energy = 0.117107, nHits = 6
Rec Hit: energy = 0.0263415, time = 17.3, pos.z() = -4107.5
Sim Hit: energy = 0.000220031
Contrib: energy = 0.000220031, time = 17.303
Particle: energy = 1.00969, pdg = -211, gen status = 1
[Contrib #0] (cluster, reco hit, sim hit, contrib, particle) ID = (1, 1, 24, 89, 3)
Cluster: energy = 0.117107, nHits = 6
Rec Hit: energy = 0.132992, time = 14.29, pos.z() = -4107.5
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Sim Hit: energy = 0.000865865
Contrib: energy = 0.00016734, time = 14.7966
Particle: energy = 1.00969, pdg = -211, gen status = 1
```

Cluster 2:

cluster.getEnergy(): 0.117107 hcalreco neutronE: 0 hcalreco pionE: 0.839073

Constituent RecHit energies can be > Cluster energies *continued discussion on next slide

> *Rec hits are mergedRechits [look at the pos.z()]



Cluster energy is determined after doing a weighted sum of the mergedhits. If the weight is too small, cluster energy can be < a constituent mergedhit energy.

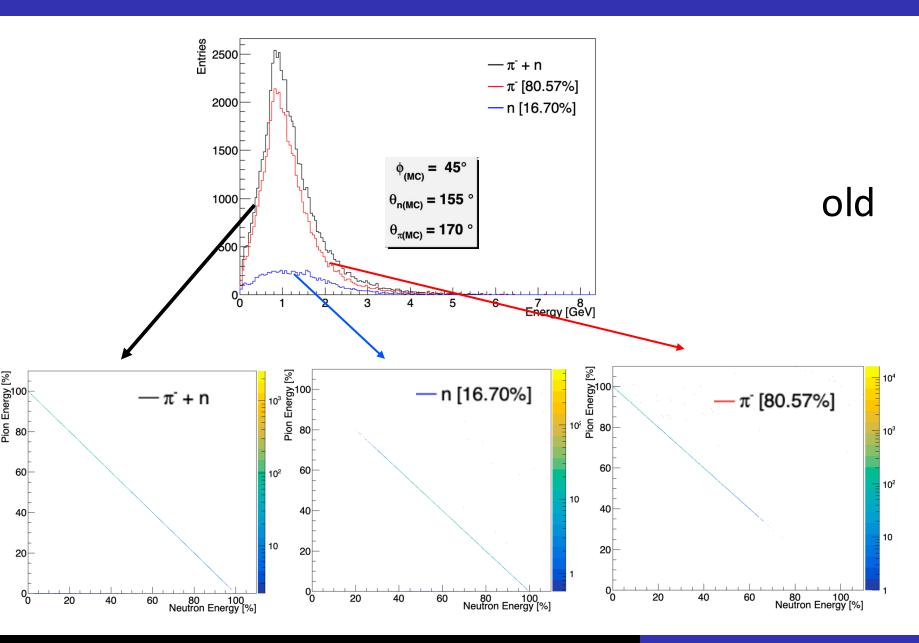
```
for (unsigned i = 0; i < pcl.getHits().size(); ++i) {
const auto& hit = pcl.getHits()[i];
const auto weight = pcl.getWeights()[i];
debug("hit energy = {} hit weight: {}", hit.getEnergy(), weight);
auto energy = hit.getEnergy() * weight;
totalE += energy;
time += (hit.getTime() - time) * energy / totalE;
cl.addToHits(hit);
cl.addToHitContributions(energy);
const float eta = edm4hep::utils::eta(hit.getPosition());
if (eta < minHitEta) {</pre>
minHitEta = eta;
if (eta > maxHitEta) {
maxHitEta = eta:
cl.setEnergy(totalE / m_cfg.sampFrac);
cl.setEnergyError(0.);
cl.setTime(time);
cl.setTimeError(timeError);
```

If a mergedRecoHit is far away from a local maxima; it will have a less weight to the cluster corresponding to that local maxima.

```
for (std::size t idx : group) {
size ti = 0;
// calculate weights for local maxima
for (std::size t cidx: maxima) {
double energy = hits[cidx].getEnergy();
double dist = edm4hep::utils::magnitude(transverseEnergyProfileMetric(hits[cidx], hits[idx]));
weights[j] = std::exp(-dist * transverseEnergyProfileScaleUnits / m cfg.transverseEnergyProfileScale) * energy;
i += 1;
// normalize weights
vec normalize(weights);
// ignore small weights
for (auto& w : weights) {
if (w < 0.02) {
w = 0:
vec normalize(weights);
// split energy between local maxima
for (size t = 0; k < maxima.size(); ++k) {
double weight = weights[k];
if (weight <= 1e-6) {
continue;
pcls[k].addToHits(hits[idx]);
pcls[k].addToWeights(weight);
```

*no of local maxima = no ofclusters





Thank You



