

PIDTOOLS FOR PARTICLE ID

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INTRODUCTION

- We have released the first version of PIDTools for ilcsoft v01– 17–07
- PIDTools module includes:
 - dE/dx calculation using TPC hit information
 - Shower profile extraction using hits in calorimeter
 - Particle Identification using all the information of tracks

• We will show usage of outputs of PIDTools modules

DE/DX

- Need to reconstruct Mokka simulation, because any information about TPC hits is lost in current public DST files.
 - After FullLDCTracking_MarlinTrk is necessary
 - Include the processor in steer file:

```
<!-- dedx calculation processor --->
cessor name="MyCompute_dEdxProcessor"/>
</processor name="MyCompute_dEdxProcessor" type="Compute_dEdxProcessor">
cessor name="EdxProcessor" type="Compute_dEdxProcessor">
cessor
```

• Just use getdEdx() in EVENT::Track collection:

like trk->getdEdx();

trk->getdEdxError(); //dEdx error, but simple 5%estimation

SHOWER PROFILE

- Need to reconstruct Mokka simulation, because any information about Cluster hits is lost in current public DST files.
 - After MarlinPandora is necessary
 - Include the processor in steer file:

```
cessor name="MyComputeShowerShapesProcessor"/>
<processor name="MyComputeShowerShapesProcessor" type="ComputeShowerShapesProcessor">
<!--Performs Shower profile extraction-->
<!--Debugging?-->
 <parameter name="Debug" type="int">0 </parameter>
<!--Name of the PFO collection-->
<parameter name="PFOCollection" type="string"> PandoraPFOs </parameter>
<!--Name of the Cluster collection-->
<parameter name="ClusterCollectionName" type="string"> PandoraClusters </parameter>
<!-- Radiation Length of Ecal-->
 <parameter name="RadiationLength Ecal" type="float">3.50 </parameter>
<!-- Radiation Length of Hcal-->
 <parameter name="RadiationLength Hcal" type="float">17.57 </parameter>
<!-- Moliere radius of Ecal-->
<parameter name="MoliereRadius Ecal" type="float">9.00 </parameter>
<!-- Moliere radius of Hcal-->
 <parameter name="MoliereRadius Hcal" type="float">17.19 </parameter>
</processor>
```

SHOWER PROFILE

- Need to reconstruct Mokka simulation, because any information about Cluster hits is lost in current public DST files.
 - Just use getShape() in EVENT::Cluster collection: like clu->getShape();
 - getShape() returns FloatVec, so there are some variables as results of shower shape fitting
 - Useful info is:

```
getShape()[0]: fitting \chi 2
```

```
getShape()[1]: maximum energy deposit(GeV)
```

```
getShape()[2]: showerMax(mm)
```

```
getShape()[3]: transverse absorption Length(mm)
```

(distance between shower axis and the point where energy deposit reduce to 1/e) getShape()[16]: xl20(mm)

(distance at the point where accumulated energy deposit is 20% of total energy deposit along shower axis(from cluster start))

```
• So far, not so sophisticate, just for PID
```

PARTICLE ID

- Need to reconstruct Mokka simulation, because any information about TPC hits and Cluster hits are lost in current public DST files.
 - Include the processor in steer file:

```
<!--processor name="MyPFOID" /-->
 cessor name="MyLikelihoodPID" />
cessor name="MyLikelihoodPID" type="LikelihoodPIDProcessor">
<!--Performs particle identification-->
<!--Debugging?-->
<parameter name="Debug" type="int">0 </parameter>
<!--Boundaries for energy binning-->
<parameter name="EnergyBoundaries" type="FloatVec">0 1.0e+07 </parameter>
<!--Name of files containing pdfs for charged particles-->
<parameter name="FilePDFName" type="StringVec">
  LikelihoodPID_Histogram_v01.root
</parameter>
<!--Name of the PFO collection-->
<parameter name="RecoParticleCollection" type="string">PandoraPFOs</parameter>
</processor>
```

 PDF file(Likelihood_Histogram_v01.root) is necessary to calculate likelihood

PARTICLE ID

 Particle ID outputs are stored to ParticleID class, so we can use standard method:

Direct access(I think it is easy…)

<pre>lcio::ReconstructedParticle *pfo;</pre>	//get PFO particles
pfo->getParticleIDs()[0]->getPDG();	<pre>//get absolute PDG value(no charge)</pre>
<pre>pfo->getParticleIDs()[0]->getLikelihood();</pre>	//get posterior prob.
<pre>pfo->getParticleIDs()[0]->getParameters()[0];</pre>	//electron hypothesis
<pre>pfo->getParticleIDs()[0]->getParameters()[1];</pre>	//muon hypothesis
<pre>pfo->getParticleIDs()[0]->getParameters()[2];</pre>	//pion hypothesis
<pre>pfo->getParticleIDs()[0]->getParameters()[3];</pre>	//kaon hypothesis
<pre>pfo->getParticleIDs()[0]->getParameters()[4];</pre>	//pion hypothesis

- So far Particle ID method is only one, so getParticleIDs()[0] is enough
- But, in next version, it is not guaranteed

PARTICLE ID

 Particle ID outputs are stored to ParticleID class, so we can use standard method:

• Using PIDHandler(easy?)

PIDHandler *pidh= new PIDHandler(pfocol); //register PFO collection to PIDHandler lcio::ReconstructedParticle *pfo; //get PFO particles pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getPDG(); //get absolute PDG value(no charge) pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getLikelihood(); //get posterior prob. pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[0]; //electron hypothesis pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[1]; //muon hypothesis pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[2]; //pion hypothesis pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[2]; //pion hypothesis pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[3]; //kaon hypothesis pidh->getParticleID(pfo, pidh->getAlgorithmID("LikelihoodPID")).getParameters()[4]; //proton hypothesis

Seems safe for extension of PID

PLOTS



Shower profile:
 ex) Absorption length





PLOTS

o @500GeV

• Momentum dep. of PID efficiency



SUMMARY & TODO

- We have released the first version of PIDTools
 - We can use dE/dx information
 - We can use shower profile information(but still messy)
 - We can use Particle ID using all the track information

o Todo:

- Forming several types of PID, using just part of track information e.g.) use dE/dx only, use shower profile only, etc.
- Some variables are not yet included, so we need to use them(backups)
 - o μ / π separation will become better
- Many minor updates…



SHOWER PROFILE -STRUCTURE IN THE CLUSTER



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INTRODUCE VARIABLES Variables are almost same as cut based: 0

- E/P, EM/(EM+HAD), |d0|, |Z0|, cone energy •
- Using these variables as pdf

Introduce new variables: 0

 $Q \times \Delta x$, Δz - distance between the cluster position and expected position when tracks are • extrapolated to the radius of the cluster position(Q is charge)

