

# *Electrons in the AOD*

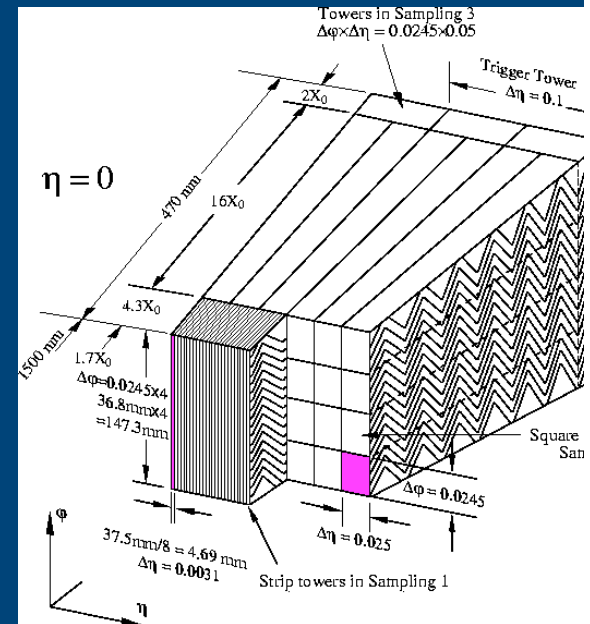
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- Outline:
    - The ATLAS EM calorimeter
      - Showers/cells/clusters & calibration  
(see my talk from yesterday)
    - Electrons in real Life
      - Identification of electrons in our detector.
      - What observables to use?
    - Electrons in the AOD
      - What variables to use for identification
      - How to access cluster/track information or other information one might need for analysis
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# Electron identification in real Life

- How to identify electrons (in very simple words):

- shower in the EM calorimeter
- energy deposited in cells/compartments of the calorimeter
- Cells/compartments form clusters
- position/energy of cluster gives information about position and energy of shower
- check if cluster is matching with a track (separate electrons and photons).
- do some cuts on calorimeter/track observables to reject jets.



- Discriminating variables from the calorimeters

- Hadronic leakage
  - Electromagnetic showers deposit a small amount of energy in the hadronic calorimeter, typically less than 2% for electrons.
- Use of the second compartment of the EM calorimeter
  - Electromagnetic showers deposit most of their energy in the second sampling of the electromagnetic calorimeter. Variables used:
    - The lateral shower shape,  $R_{\eta}(37)$ , is given by the ratio of the energy reconstructed in a  $3 \times 7$  cluster to the energy in a  $7 \times 7$  cluster. This variable shows a peak near one for electrons because of the very small lateral leakage; large tails at lower values of  $R_{\eta}(37)$  for the jets are expected.
    - The lateral width is calculated with a window of  $3 \times 5$  cells using the energy weighted sum over all cells, which depends on the particle impact point inside the cell

- Discriminating variables from the calorimeters (2)

- Use of the first compartment of the EM calorimeter

- The first compartment with its very fine granularity in rapidity can be used to detect substructures within a shower and thus isolated  $\pi_0$  discriminated against efficiently

Variables used:

- $\Delta E = E_{\max 2} - E_{\min}$

- $\Delta E_{\max 2} = E_{\max 2} / (1 + 9(5) \times 10^{-3} E_T)$ , with  $E_T$  the transverse energy of the cluster

- $\omega_{\text{tot1}} \sim \sim \sqrt{(\sum E_i \times (i - i_{\max})^2 / \sum E_i)}$ , where  $i$  is the strip number and  $i_{\max}$  the strip number of the first local maximum.

- $F_{\text{side}} = [E(\pm 3) - E(\pm 1)] / E(\pm 1)$ , where  $E(\pm n)$  is the energy in  $\pm n$  strips around the strip with highest energy.

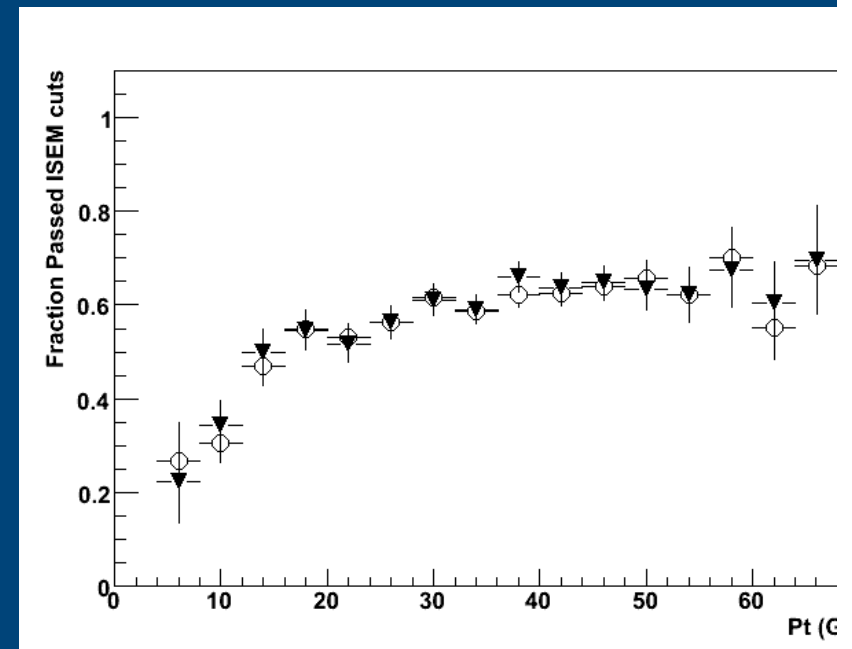
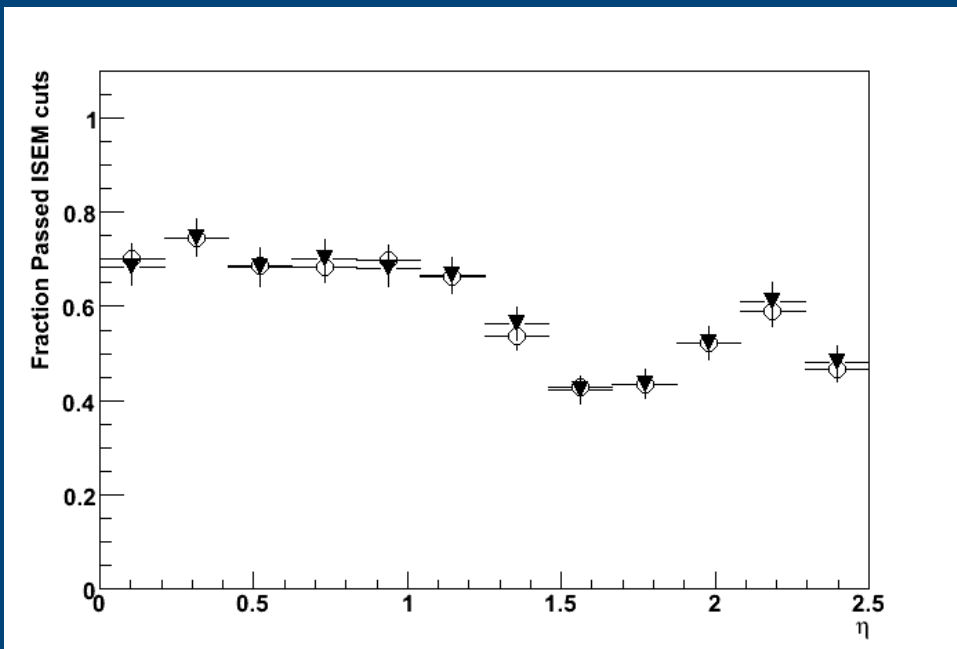
- $\omega_{3 \text{ strips}} = \sqrt{\{\sum E_i \times (i - i_{\max})^2 / \sum E_i\}}$ , where  $i$  is the number of the strip and  $i_{\max}$  the strip number of the most energetic one.

- Discriminating variables from the inner detector
    - Track quality cuts
      - at least nine precision hits (Pixel+SCT)
      - at least two hits in the pixels, one of which being in the b-layer
      - a transverse impact parameter  $A_0 < 0.1$  cm
    - Inner detector/calorimeter spatial matching information
      - $\Delta\eta = |\eta_{\text{strips}} - \eta_{\text{ID}}|$ , where  $\eta_{\text{strips}}$  is computed in the first sampling of the electromagnetic calorimeter, where the granularity is very fine, and  $\eta_{\text{ID}}$  is the pseudo-rapidity of the track extrapolated to the calorimeter.
      - $\Delta\varphi = |\varphi_{\text{middle}} - \varphi_{\text{ID}}|$ , where  $\varphi_{\text{middle}}$  is computed in the second compartment of the electromagnetic calorimeter and  $\varphi_{\text{ID}}$  is the azimuth of the track extrapolated to the calorimeter.
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- Discriminating variables from the inner detector (2)
    - Inner Detector/calorimeter energy matching information
      - The energy  $E$  measured in the electromagnetic calorimeter is compared to the momentum  $p$  measured in the Inner Detector. In the case of an electron, the momentum should match the energy.
    - Transition radiation in the TRT information
      - A further reduction of the charged hadron contamination is obtained by rejecting tracks having a low fraction of high-threshold hits. The discriminating variable used is the ratio  $N_{\text{high}}/N_{\text{all}}$  between the number of high threshold hits and the total number of TRT hits.
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# Identification efficiencies

- IsEM tight



# *Electrons in the AOD*

- The electron collection in the AOD, the keys to use are:
    - For the collection: ElectronAODCollection
    - For the details (Shower/Trackmatch): egDetailAOD
  - ElectronAODcollection
    - If a cluster has a matched track, the egamma object is added to the electron collection
  - IsEM flag
    - The candidate has to pass a series of cuts based on the shower shape properties in different compartments of the calorimeter as well as variables combining ID and Calo informations (see previous slides). If a cut is not passed, then a bit is set in the IsEM flag
    - You can use the bit mask to use only certain parts of the ID criteria
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- Predefined set of cuts:
  - Loose cuts: `electron->isem(ElectronLoose)==0`
    - Calorimeter cuts only
  - Medium cuts: `electron->isem(ElectronMedium)==0`
    - Calorimeter cuts & Track quality cuts (no b-layer)
  - Tight cuts: `electron->isem(ElectronTight)==0`
    - All cuts (previous & Track matching & TRT cuts)
  - All these cuts can be turned on or off individually using bit masks

```
enum BitDef {
  // Cluster based egamma
  ClusterEtaRange      = 0,
  ClusterHadronicLeakage = 1,
  ClusterMiddleEnergy  = 4,
  ClusterMiddleEratio37 = 5,
  ClusterMiddleEratio33 = 6,
  ClusterMiddleWidth   = 7,
  ClusterStripsEratio  = 8,
  ClusterStripsDeltaEmax2= 9,
  ClusterStripsDeltaE   = 10,
  ClusterStripsWtot    = 11,
  ClusterStripsFracm   = 12,
  ClusterStripsWeta1c  = 13,
  ClusterIsolation     = 14,
  //Track based egamma
  TrackBlayer         = 16,
  TrackPixel          = 17,
  TrackSi             = 18,
  TrackA0            = 19,
  TrackMatchEta       = 20,
  TrackMatchPhi       = 21,
  TrackMatchEoverP    = 22,
  TrackTRThits        = 24,
  TrackTRTratio       = 25,
  TrackTRTratio90     = 26
};
```



## • How to use the IsEM flag?

```
StoreGateSvc* m_storeGate;
const ElectronContainer* electronColl;
m_storeGate->retrieve(electronColl, "ElectronAODCollection");

ElectronContainer::const_iterator eleItr = electronColl->begin();
ElectronContainer::const_iterator eleEnd = electronColl->end();

for(; eleItr != eleEnd; ++eleItr) {
int isemLoose = (*eleItr)->isem(egammaPID::ElectronLoose);
int isemMedium = (*eleItr)->isem(egammaPID::ElectronMedium);
int isemTight = (*eleItr)->isem(egammaPID::ElectronTight);
if (isemLoose == 0) std::cout << "looooooose cut passed " << std::endl;
}
```

If you want to apply your own set of cuts you can use the bit masks like this:  
For example, let's say you want to use only the E/p cut and the Hadronic Leakage  
(of course you wouldn't wanna do that)

The bit definition on the previous slide tells us:

E/p cut is bit number 22

Hadronic Leakage is bit number 1

So you do:

```
(*electron)[index]-> isem(0x400002);
```

WARNING: Bit definition is different in 12.0.X

Closest to loose definition:  $(isem \& 0x7) == 0$

Closest to medium:  $(isem \& 0x3FF) == 0$

Closest to tight:  $isem = 0$

- All calorimeter observables used for identification are accessible in the AOD:
  - Shower variables
    - `egammaParameters::etcone20` : isolation energy (transverse) in a cone with half-opening angle 0.2
    - `egammaParameters::ethad1` : transverse energy in the first sampling of the hadronic calorimeters behind the cluster
    - `egammaParameters::e233` : uncalibrated energy (sum of cells) of the middle sampling in a rectangle of size  $3 \times 3$  cells
    - `egammaParameters::e237` : uncalibrated energy (sum of cells) of the middle sampling in a rectangle of size  $3 \times 7$  cells
    - `egammaParameters::e277` : uncalibrated energy (sum of cells) of the middle sampling in a rectangle of size  $7 \times 7$  cells
    - `egammaParameters::weta1`
    - `egammaParameters::weta2`
    - `egammaParameters::f1`
    - `egammaParameters::e2tsts1`
    - `egammaParameters::emins1`
    - `egammaParameters::wtots1`
    - `egammaParameters::frac1`
    - `egammaParameters::f1core`
    - `egammaParameters::f3core`
    - `egammaParameters::pos7`
    - `egammaParameters::iso`
    - `egammaParameters::widths2`
    - `egammaParameters::zvertex`
    - `egammaParameters::errz`
    - `egammaParameters::etap` : pointing eta reconstructed from the cluster (first and second sampling)
    - `egammaParameters::depth`

- EMTrackMatch variables

- egammaParameters::EtaCorrMag
- egammaParameters::EoverP : ratio of the cluster energy and the track momentum
- egammaParameters::deltaEta1 : difference between the cluster eta (first sampling) and the eta of the track extrapolated to the first sampling
- egammaParameters::deltaEta2 : difference between the cluster eta (second sampling) and the eta of the track extrapolated to the second sampling
- egammaParameters::deltaPhi2 : difference between the cluster phi (second sampling) and the phi of the track extrapolated to the second sampling

- How to access EMShower and EMTrackMatch information

```
const EMShower *shower;  
shower = (*eleItr)->detail<EMShower>("egDetailAOD")  
float etcone = shower->parameter(egammaParameters::etcone20)
```

```
const EMTrackMatch *trackmatch;  
trackmatch = (*eleItr)->detail<EMTrackMatch>("egDetailAOD")  
float etcone = shower->parameter(egammaParameters::EoverP )
```

# • Clusters and Tracks

pointer to CaloCluster :     const CaloCluster \* egamma::cluster() const  
pointer to TrackParticle :   const Rec::TrackParticle \* egamma::trackParticle() const

(\*eleItr)->trackParticle()->phi()

(\*eleItr)->trackParticle()->eta()

(\*eleItr)->trackParticle()->measuredPerigee()->parameters()[Trk::d0]

(\*eleItr)->trackParticle()->measuredPerigee()->localErrorMatrix().error(Trk::d0)

(\*eleItr)->trackParticle()->trackSummary()->get(Trk::numberOfBLayerHits)

(\*eleItr)->trackParticle()->trackSummary()->get(Trk::numberOfPixelHits)

(\*eleItr)->cluster() ->phi()

(\*eleItr)->cluster() ->eta()

(\*eleItr)->cluster() -> energy()

(\*eleItr)->cluster() -> energyBE(0)

(\*eleItr)->cluster() -> energyBE(1)

(\*eleItr)->cluster() -> energyBE(2)

(\*eleItr)->cluster() -> energyBE(3)