Energy correction for PANDA EMC

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Outline



- 1 Electromagnetic calorimeter
- 2 EMC in PandaRoot
- 3 Energy normalization
- 4 Energy correction
- 5 Summary

Electromagnetic calorimeter PANDA EMC



PANDA EMC

- · Components: Target + Forward
- Goal
- Requirement
 - Coverage
 - ♦ Energy threshold (10(20) MeV)
 - Geometry coverage (99%x4π in c.m.s)
 - Energy range (up to 0.7 GeV BW,
 7.3 GeV B, 14.6 GeV FW)
 - Resolution
 - Single crystal threshold (σ_{E,noise} = 1MeV, E_{xtl} = 3MeV)
 - ♦ Energy Resolution $(1\% \oplus 2\%/\sqrt{E})$
 - ♦ Spatial Resolution (0.5° BW, 0.3° B, 0.1° FW)

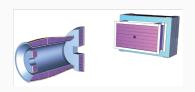


Figure: PANDA EMC

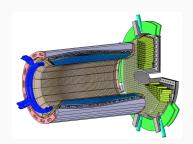


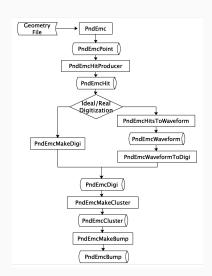
Figure: Target EMC

EMC in PandaRoot



FMC in PandaRoot

- Functionalities
 - Simulation
 - Digitization
 - Calibration
 - Reconstruction
 - Cluster finding
 - Bump splitting
 - ⋄ Correction
- Current status

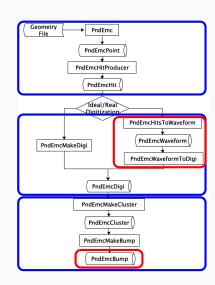


EMC in PandaRoot



EMC in PandaRoot

- Functionalities
 - Simulation
 - Digitization
 - Calibration
 - Reconstruction
 - Cluster finding
 - Bump splitting
 - Correction
- Current status
 - Functionalities are checked
 - Need Optimization



Procedure



Procedure

- Calibration
- Crystal type normalization
 - crystal type related
 - energy related
- Bump correction
 - Leakage correction
 - Preshower correction

Energy normalization



Why should we do energy normalization?

Energy

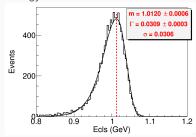
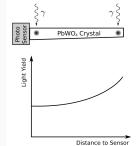


Figure: Example with 1 GeV γ . Resolution is 3.06%, worse than TDR; Events with Energy Higher than 1 GeV

Reasons

 Nonuniformity: The amount of light collected at sensor depends on the position of the energy deposition.



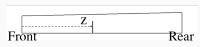
Energy normalization



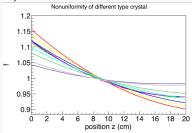
Implementation of nonuniformity in PandaRoot

light yield in a crystal

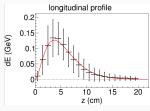
$$\mathit{Iy} = c_0 + \mathit{z}(c_1 + \mathit{z} \cdot c_2)$$



light yield of different type of crystals



- Advantage
 - More sensitive to low energy γ/e
- Disadvantage
 - Resolution Deterioration
 - Type difference of crystals
 - Light yield is not linear to incident energy (longitude depth varies with incident energy)



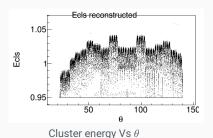
Longitude dE distribution

Energy normalization Normalization for crystal type

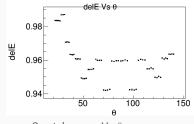


Normalization for different types of crystals

- Phenomena
 - Cluster energy varies with θ
 - One module consists of 4 crystals in θ direction
 - Different behaviours in modules



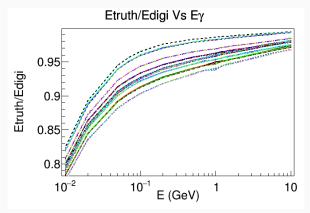
- Normalization: based on MC
 - Simulation setup: 1 GeV γ to front center of each row
 - Normalization: Scale the energy in each crystal to truth energy



Crystal energy Vs θ delE = E_{truth}/E_{digi}



Normalization for different types of crystals, energy relation

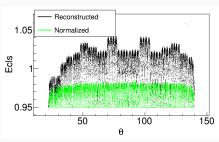


Normalization factor Vs E_{γ} , different colors for different crystal types.

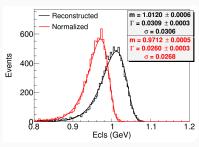
Energy normalization Normalization for crystal type



Results of the normalization



Cluster Energy Vs θ before and after normalization for crystal types.

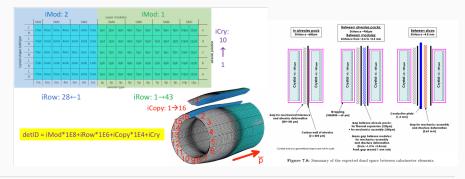


Cluster Energy of a test sample before and after normalization for crystal types..



- Gap types between crystals
 - $-\theta$ direction
 - ⋄ in a module
 - between modules
 - between supermodules

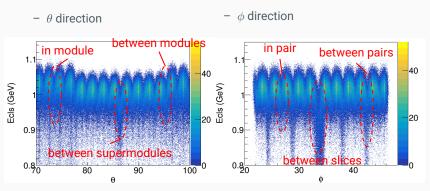
- $-\phi$ direction
 - in a pair
 - between pairs
 - between slices





Energy Leakage (including energy deposition in dead area)

Gap types between crystals





- Method description [2007 JINST 2 P04004]
- variable definition (both in θ and ϕ direction)
 - find seed crystal
 - compare energy in left and right crystals
 - ⋄ define E1 and E2

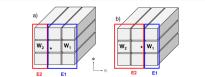


Figure 8. Definition of E1 and E2 when a) W1 < W2 and b) W1 > W2 (see text). The star represents the electron's incident position.

- Relation between E_{cluster} and In(E2/E1)
 - \diamond Relation in θ and ϕ direction

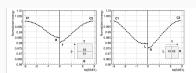
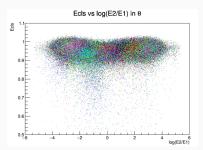
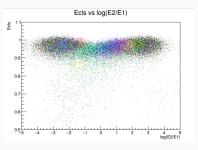


Figure 10. Normalized mean energy measured in the 3x3 array around crystal 204 versus InteZEE1) in the Φ (Heft) and γ (light) directions. The curves are 3x4 order topytomical functions fitted to the measured distributions, independently for positive and negative values of InteZEE1). The square punels represent the central crystal with various regions indicated: T (rop.) 8 (content), Left(1), R (right) and C1 and C2 (salt) off centre on each side). The labels on the distributions indicate in which region the electrons were incident.



- Apply on PANDA
- θ direction (ϕ fixed to 30.1°)
 - ♦ $E_{cluster}$ Vs In(E2/E1), where E2, E1 defined in θ direction
 - \diamond Each color corresponds to γ of a specific direction
- ϕ direction (θ fixed to 49.9°) \$\phi E_{cluster}\$ Vs In(E2/E1), where
 - \Leftrightarrow E_{cluster} VS III(EZ/ET), where E2, E1 defined in ϕ direction
 - $\diamond \;$ Each color corresponds to γ of a specific direction

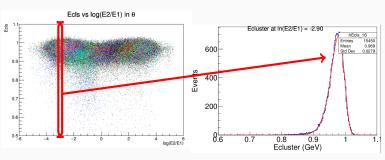






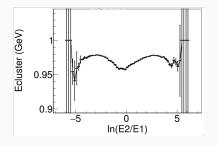
Energy Leakage (including energy deposition in dead area)

• Apply on PANDA, extract the relation split to slices \rightarrow project to $E_{cluster} \rightarrow$ fit the histogram \rightarrow Extract relation of $E_{cluster}$ Vs In(E2/E1)

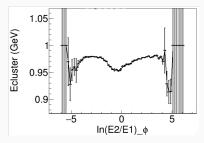




- Apply on PANDA, extract the relation
- θ direction (ϕ fixed to 30.1°)
 - ♦ $E_{cluster}$ Vs In(E2/E1), where E2, E1 defined in θ direction



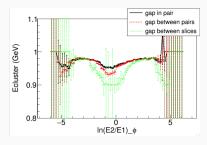
- ϕ direction (θ fixed to 49.9°)
 - \diamond $E_{cluster}$ Vs In(E2/E1), where E2, E1 defined in ϕ direction





- Apply on PANDA, extract the relation, separate different gaps
- θ direction (ϕ fixed to 30.1°)
 - \diamond $E_{cluster}$ Vs In(E2/E1), where E2, E1 defined in θ direction
 - gap between modules gap between modules gap between modules gap between modules gap between supermodules gap between supe

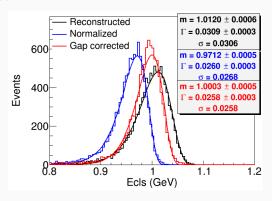
- ϕ direction (θ fixed to 49.9°)
 - \diamond $E_{cluster}$ Vs In(E2/E1), where E2, E1 defined in ϕ direction





Result

- Cluster energy corrected with crystal type has a better resolution than the one without correction. The improvement is about 10%
- Cluster energy corrected for the gap is slight better than crystal type corrected



Summary



- Functionality check
 - Functionalities of EMC part in PandaRoot has been checked from Simulation to Reconstruction
 - Digitization algorithm is updating
 - Energy and position correction are under studying
 - Calibration algorithm is in consideration
- Energy correction
 - Normalization for each crystal type has been studied, which can provide a better energy resolution
 - Energy leakage has been corrected with In(E2/E1)

Appendix Energy relation of normalization factor



