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# BeamCal electron reconstruction

Aura Rosca

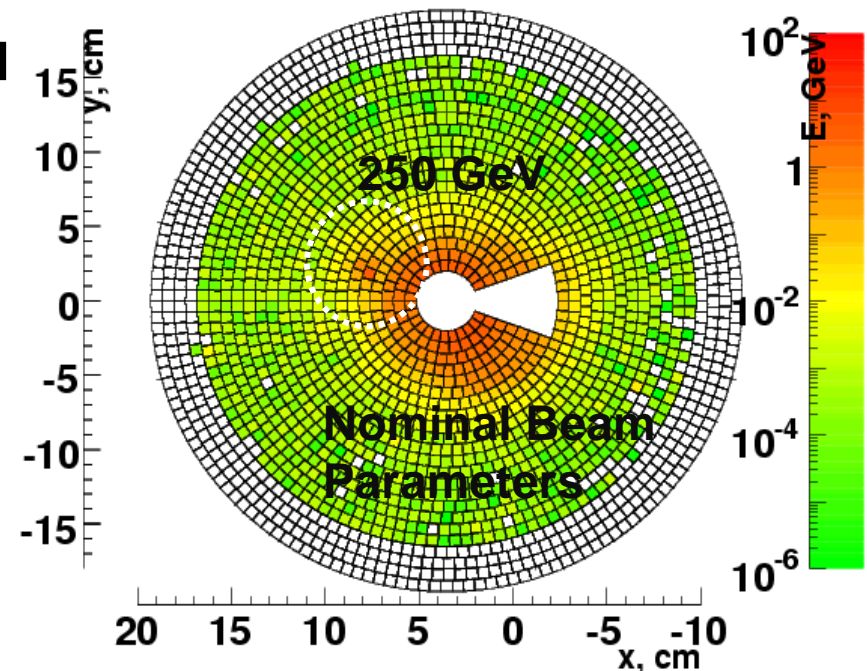
24<sup>th</sup> November 2010

# Cluster reconstruction algorithm in BeamCal

(Developed by Olga N. and Wolfgang L.)

## 1. Background subtraction procedure

- Calculate average and rms of the energy deposition of the background in each pad of the BeamCal, from 10 BX;
- Superimpose 1 BX background + 1 high energy electron;
- Subtract the value of the background average from the superposition;

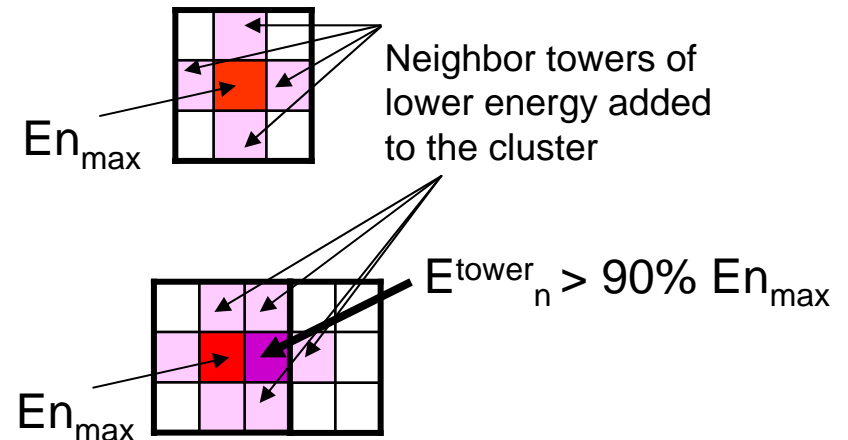
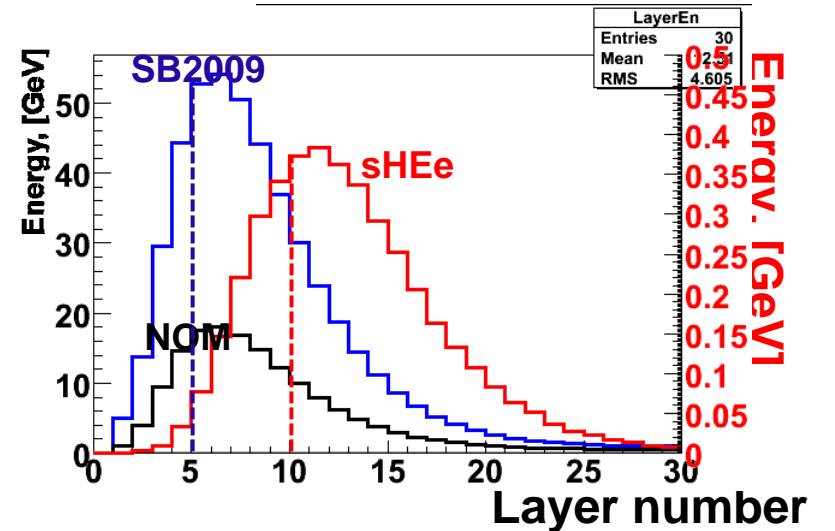


# Cluster reconstruction algorithm in BeamCal

(Developed by Olga N. and Wolfgang L.)

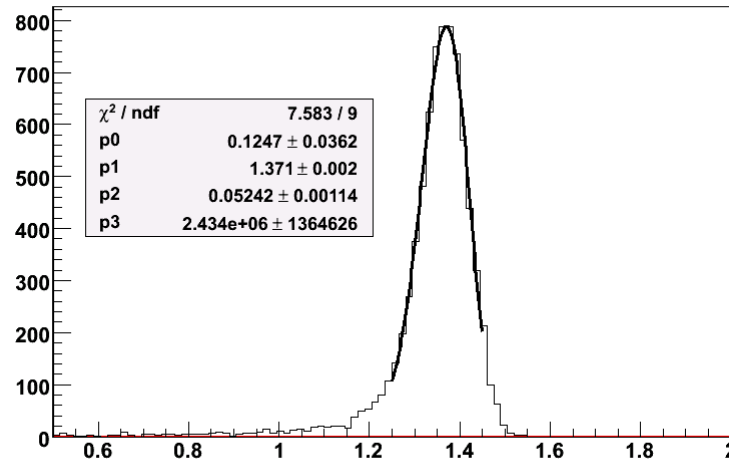
## 2. Cluster search

- Identify towers after the 5-th layer as chains of 10 consecutive fired pads;
- Search for the tower with maximum deposited energy;
- Add neighbor towers, in a 3×3 matrix around the tower with the highest energy;
- If such a neighbor tower has an energy larger than 90% of the energy of the central tower, add this tower neighbors as well;

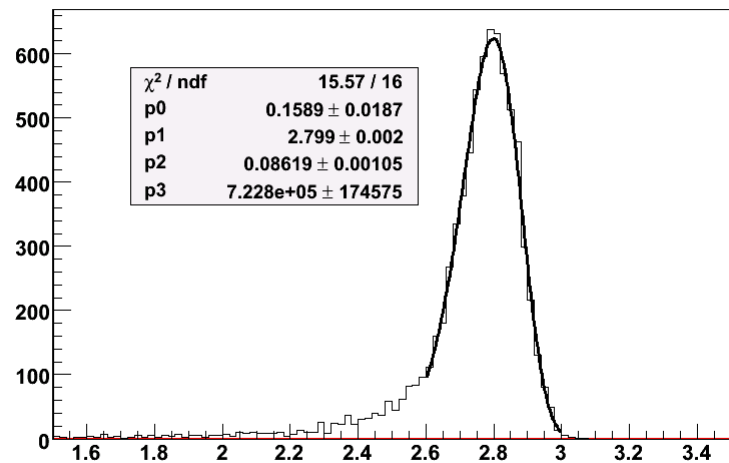


# Calibration curve

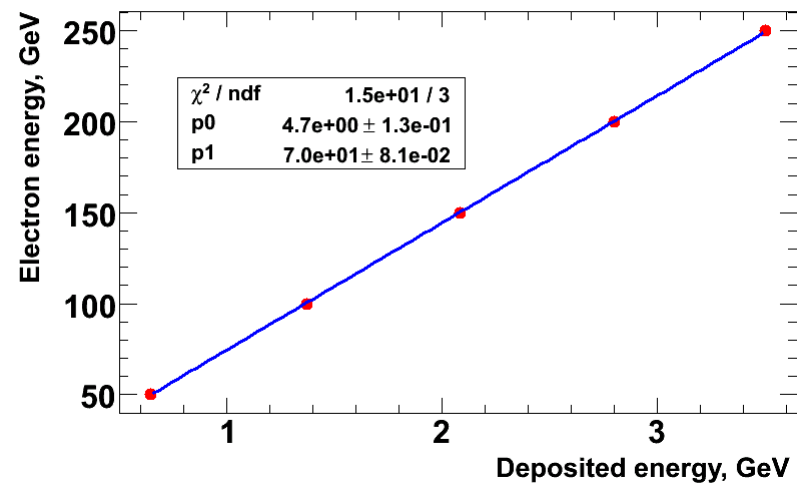
Energy deposition from 100 GeV electron showers



Energy deposition from 200 GeV electron showers



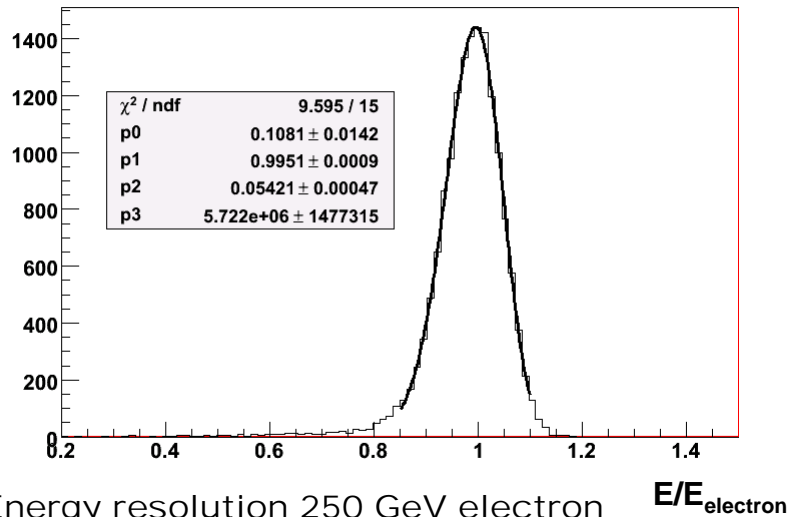
Calibration curve



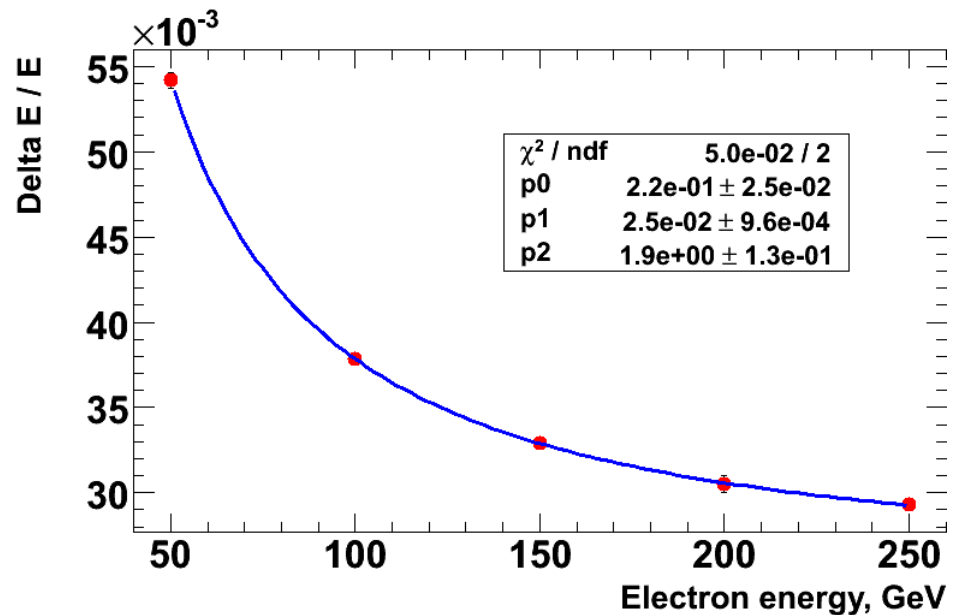
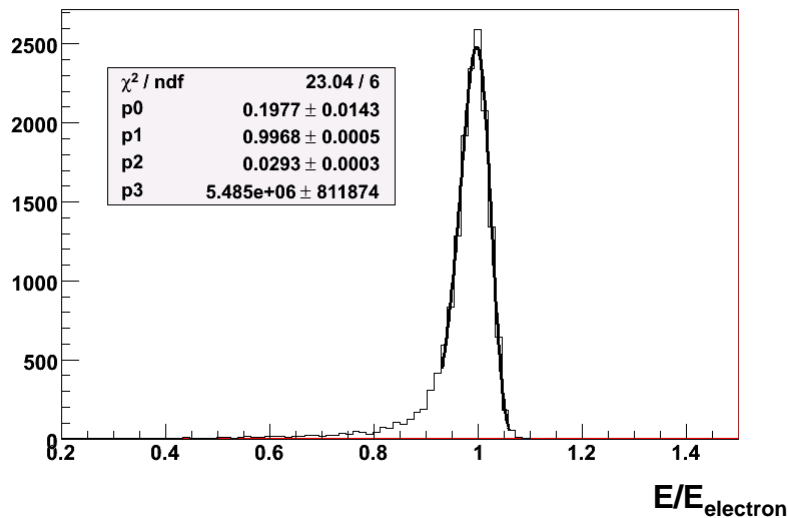
$$E = (70 \cdot E_{\text{dep}} + 4.7) \text{ GeV}$$

# Energy resolution

Energy resolution 50 GeV electron



Energy resolution 250 GeV electron



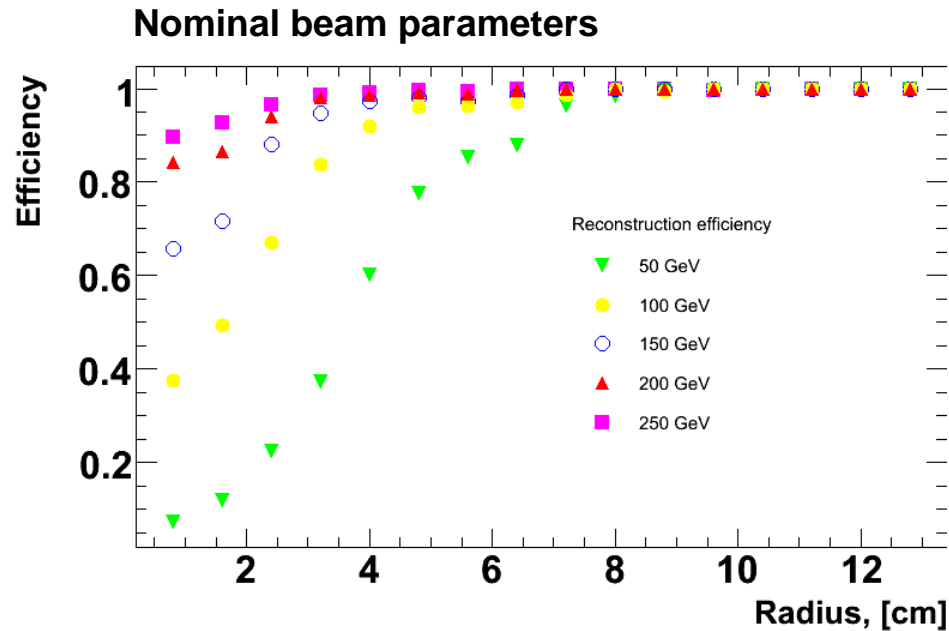
$$\sigma(E)/E = \sigma_1 / E^{1/2} \oplus \sigma_2 \oplus \sigma_3 / E$$

$$\sigma_1 = (22.0 \pm 2.50)\%$$

$$\sigma_2 = (2.5 \pm 0.1)\%$$

$$\sigma_3 = 2$$

# Reconstruction efficiency



New parameterization of the efficiency as a function of electron energy and background energy deposition, valid for nominal beam parameters:

$$p0 = f_1(\text{Energy}) = 7.534199\text{e-}02 + 7.452120\text{e-}04 * \text{Energy} - 2.001878\text{e-}06 * \text{Energy} * \text{Energy}$$

$$p1 = f_2(\text{Energy}) = 8.149966\text{e-}02 + 9.193878\text{e-}04 * \text{Energy} - 2.437620\text{e-}06 * \text{Energy} * \text{Energy}$$

$$p2 = f_3(\text{Energy}) = -1.259418\text{e-}01 + 9.974422\text{e+}03 / (\text{Energy} * \text{Energy})$$

$$\text{efficiency} = f(\text{Energy}, \text{EnDensity}_{\text{Bkgd}}) = p0 + 1 / (1 + p1 * \exp(p2 * \text{EnDensity}_{\text{Bkgd}}))$$

# Implementation in Marlin

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- Parameterization is used together with a background map of energy density deposition of the background in the BeamCal pads, in the form of a root file;
- It is of interest for physics studies that need to estimate the probability to detect a high energy electron at a position in BeamCal with a given local energy density;
- The efficiency for each particle can be accessed via the method `getGoodnessOfPID()` of the class `ReconstructedParticle` from the Marlin package;
- Parametrization should substitute the old one in the existing processor `BCalTagEfficiency`;
- Open point: who performs the implementation in the official released?

# Summary

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- Clustering algorithm in BeamCal is based mainly on finding neighbor towers of 10 consecutive pads after a background subtraction procedure is applied;
- Energy resolution in BeamCal is:  $\sigma_1 = (22.0 \pm 2.50)\%$ ;
- New parameterization exists of the electron reconstruction efficiency in BeamCal, for nominal beam parameters.