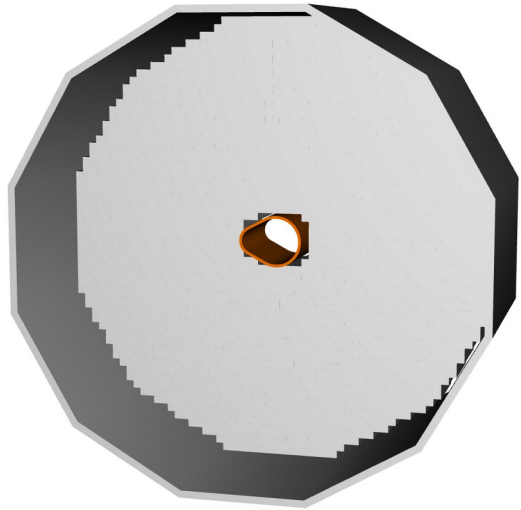


# Test Beam Prototype Simulations

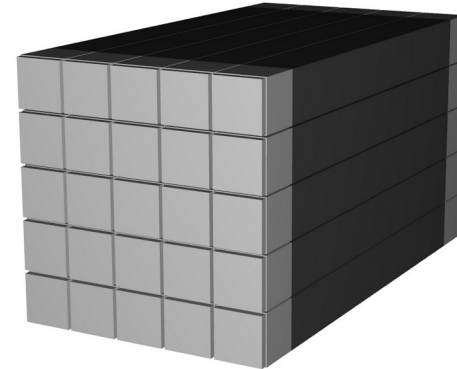
**Artur Hoghmrtsyan**

# Simulation Setup

[https://github.com/eic/epic/tree/main/build/epic\\_eemcal\\_only.xml](https://github.com/eic/epic/tree/main/build/epic_eemcal_only.xml)



## Prototype Simulations



### Particle Gun

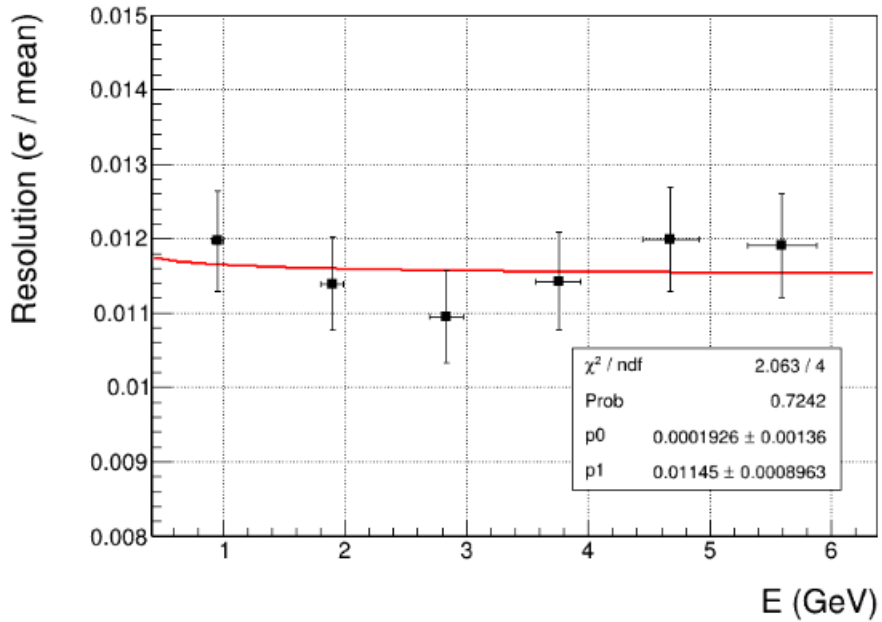
Position (0,0,-40mm)

Particle -  $e^-$

Energies - 1,2,3,4,6 (Gev)

Energy Spread - 158 (MeV)

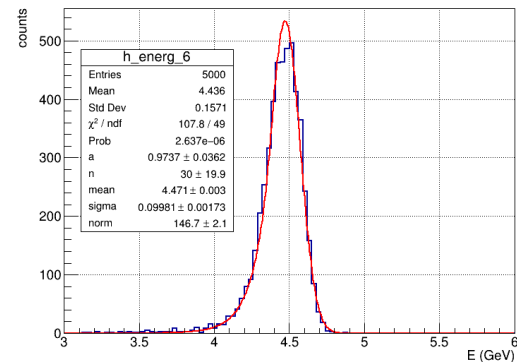
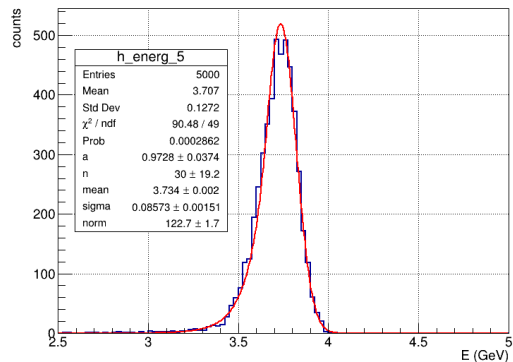
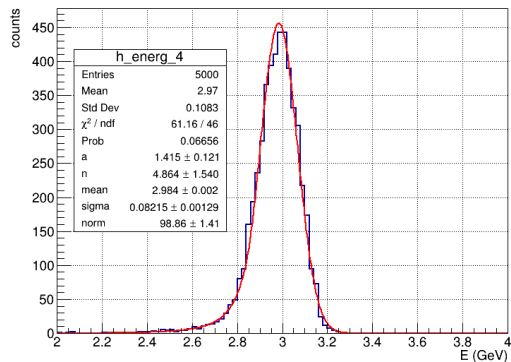
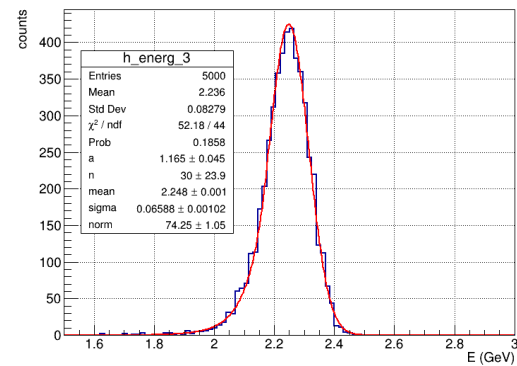
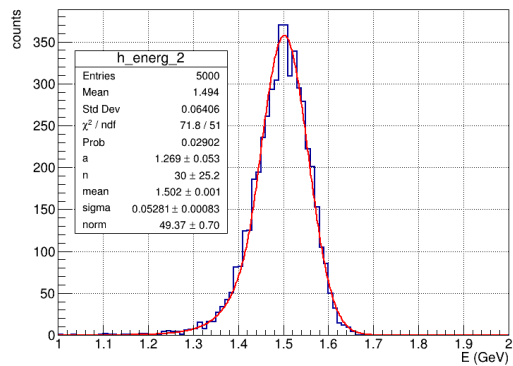
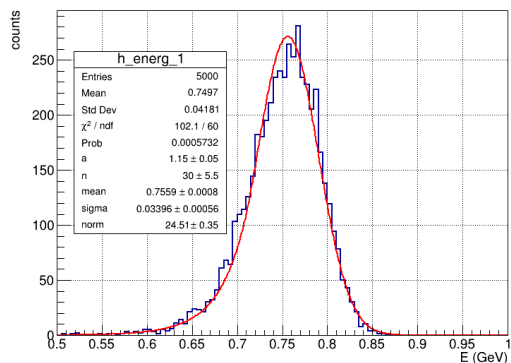
## Problem from the last update



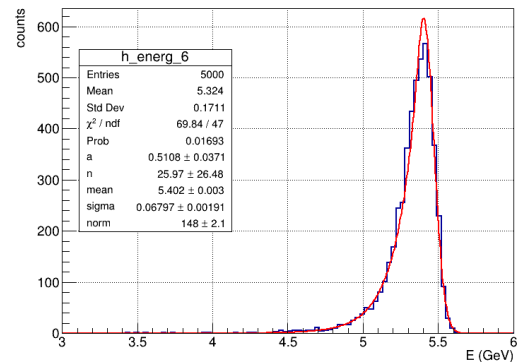
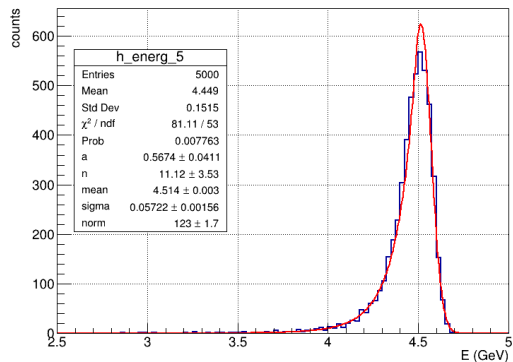
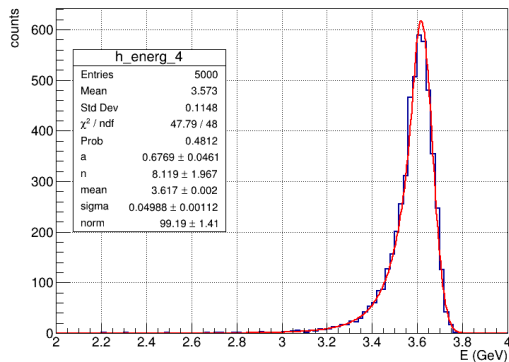
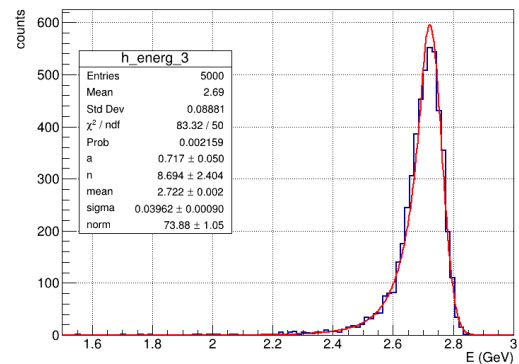
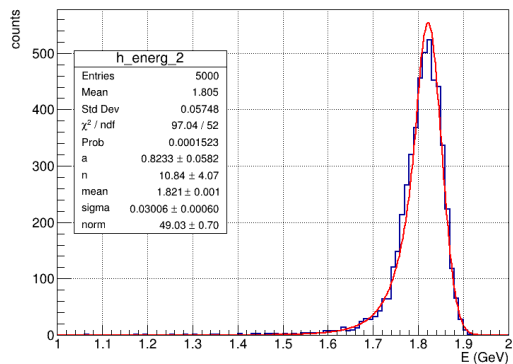
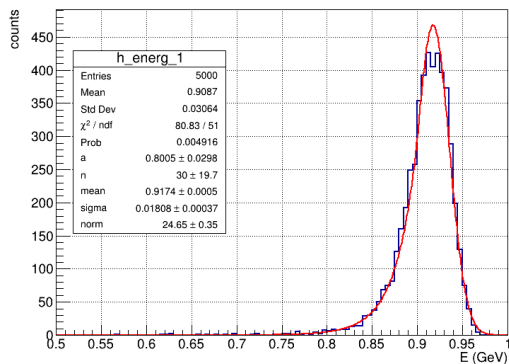
**Fixed:**

Distance between crystals was more than 1mm

# Deposited energy in Mono energetic case 1x1

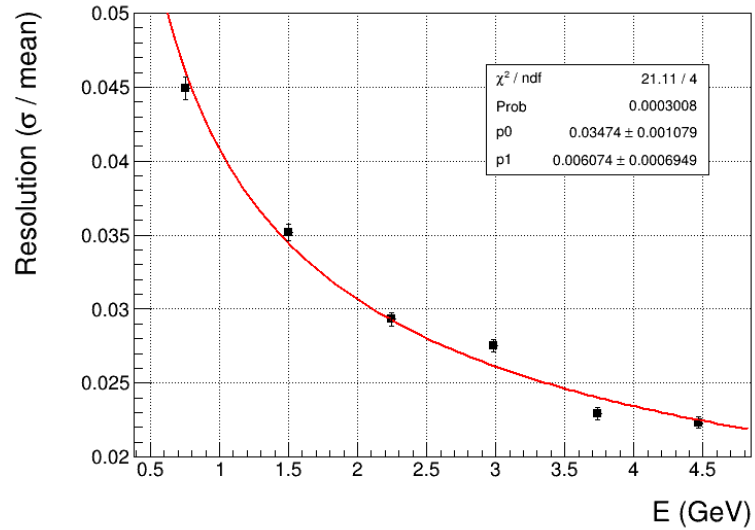


# Deposited energy in Mono energetic case 3x3 configuration

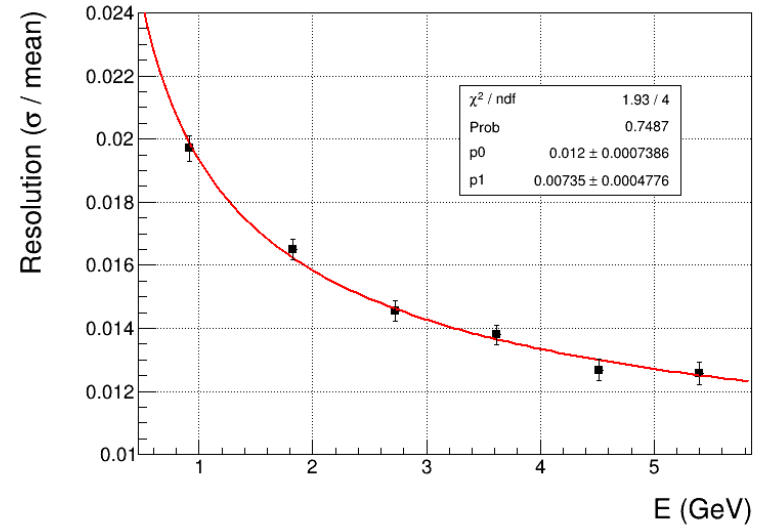


# Energy Resolution for Monoenergetic 1x1 vs 3x3 cases

## 1x1

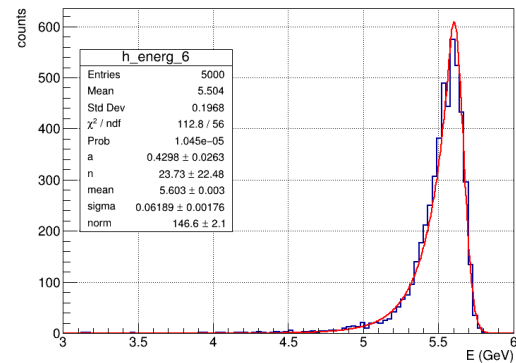
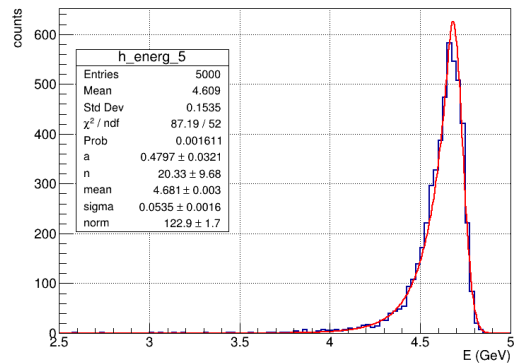
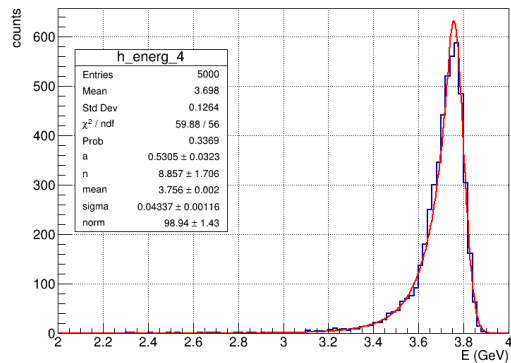
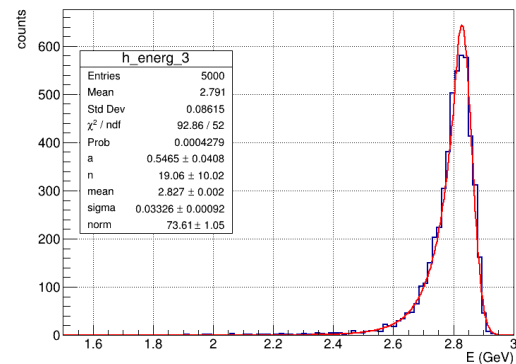
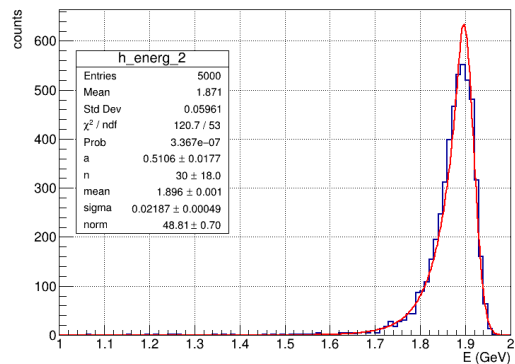
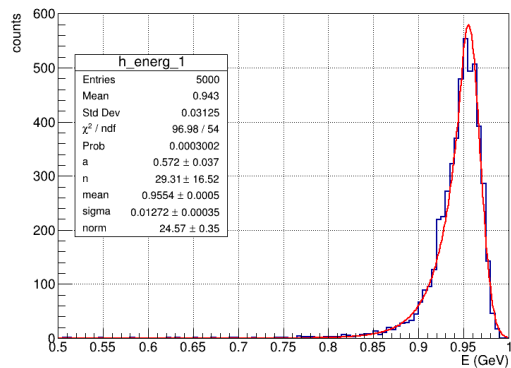


## 3x3



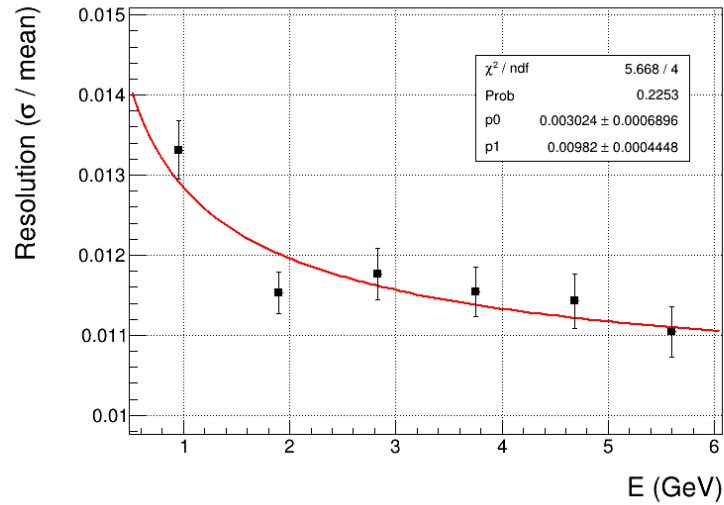
Everything looks normal

# Deposited energy in Monoenergetic case 5x5

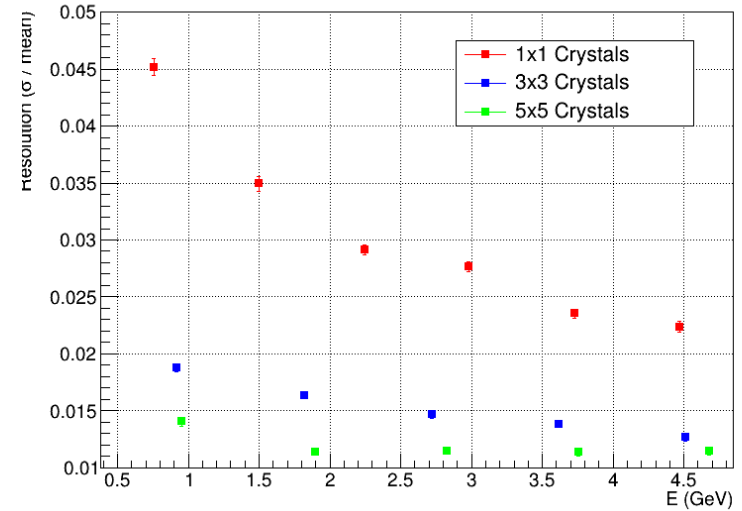


# Energy Resolution Plots

## Monoenergetic 5x5

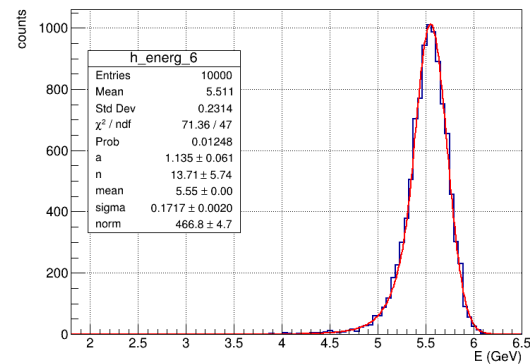
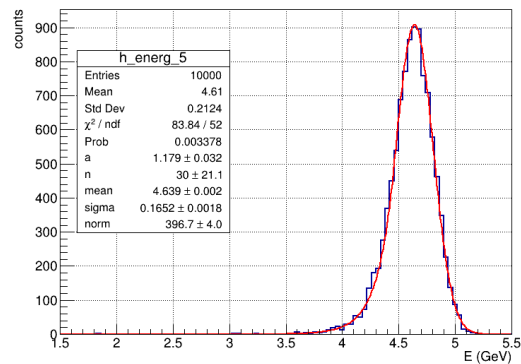
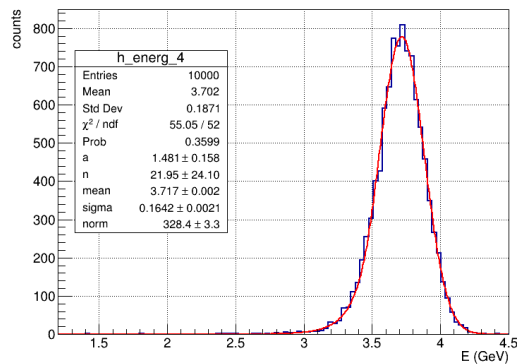
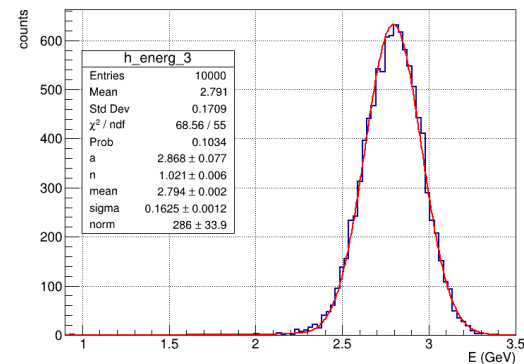
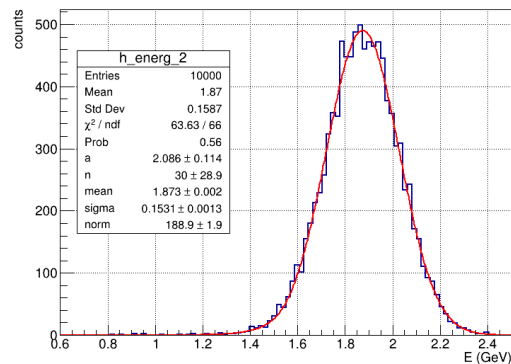
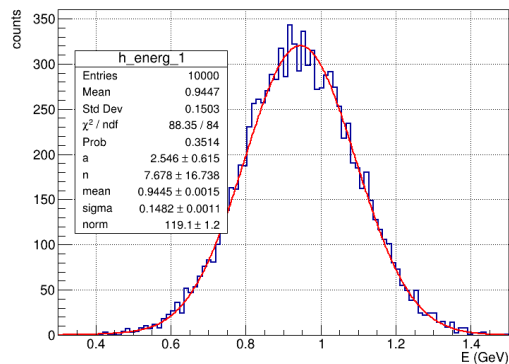


## All configurations

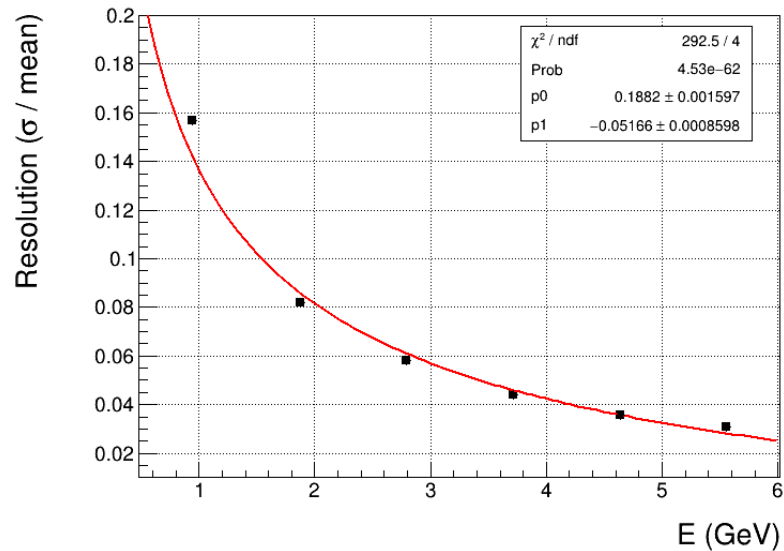
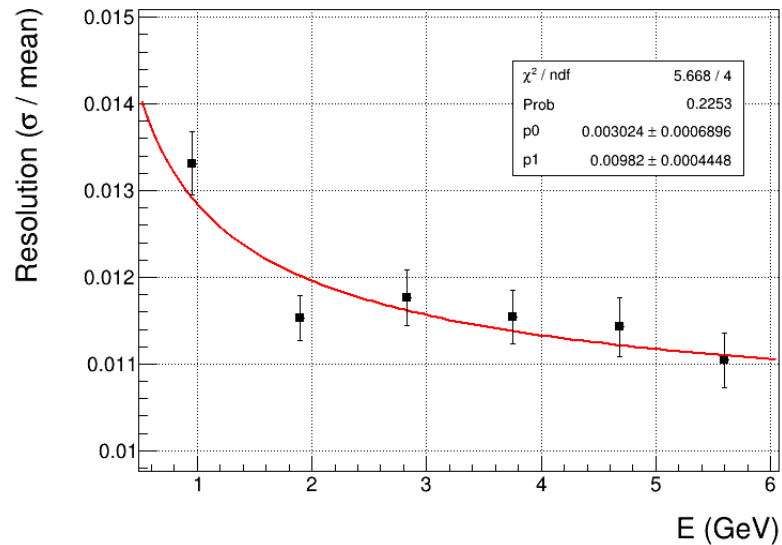




# Deposited energy in DESY energetic case 5x5

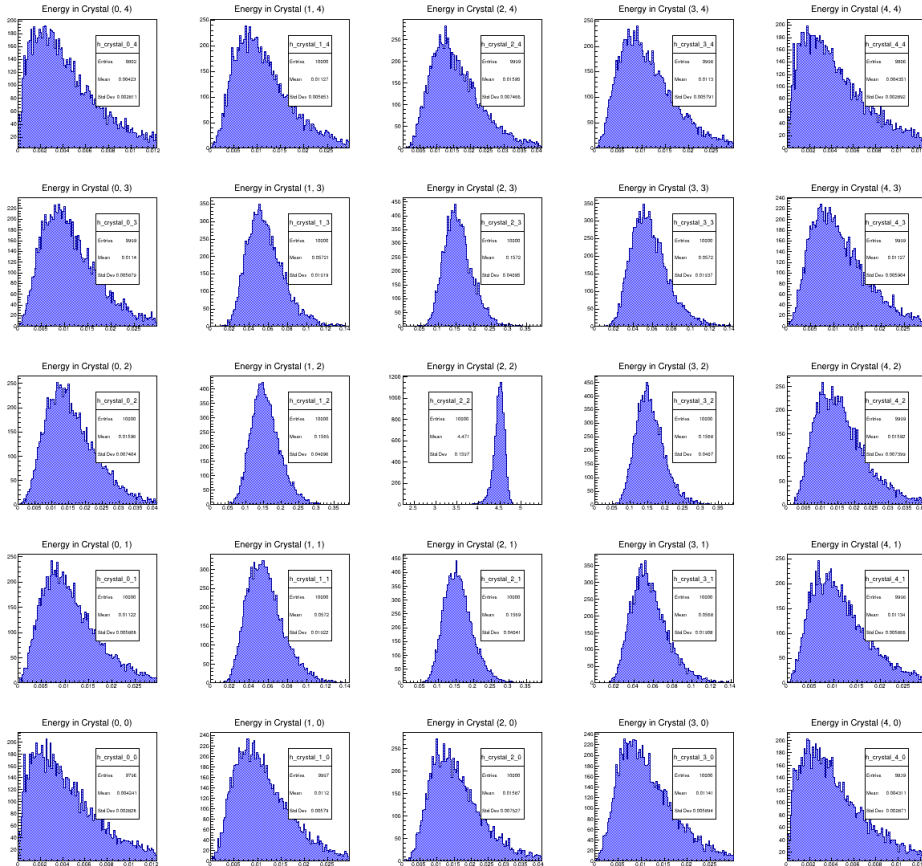


# Energy Resolution for Mono vs DESY 5x5 matrices



Big effect for low energies

# Smearing by number of Photoelectron



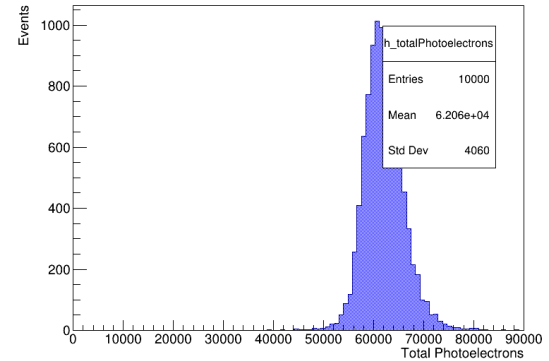
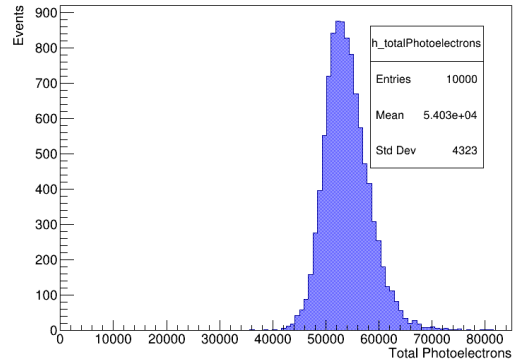
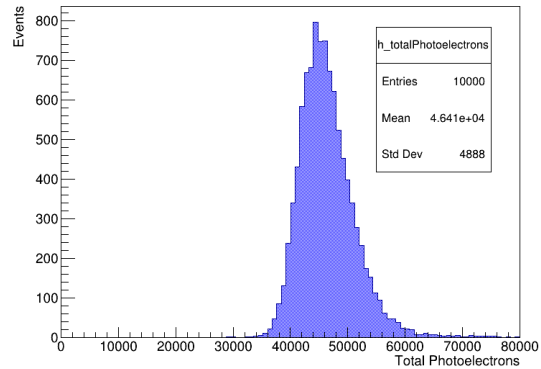
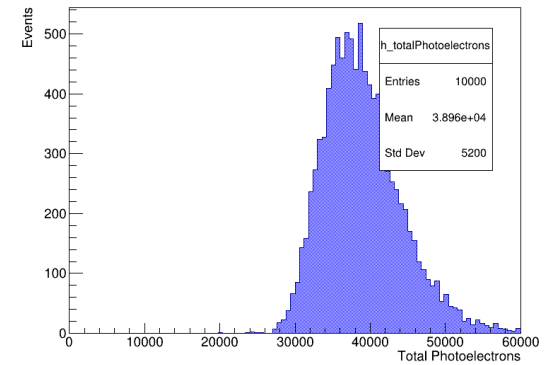
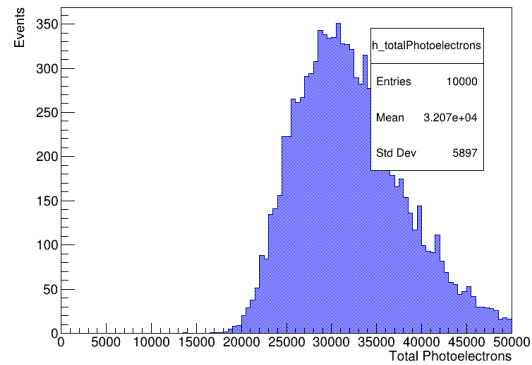
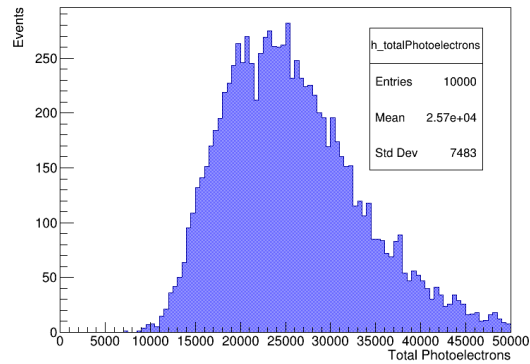
Gaussian Smearing Calculation: For each deposited energy  $E_i$  (in GeV)

$$\sigma = \frac{1}{\sqrt{10000 \cdot E_i}}$$

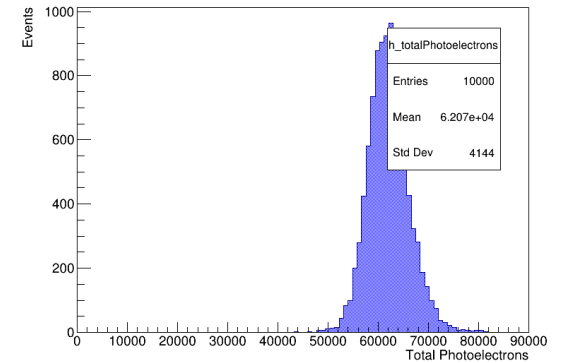
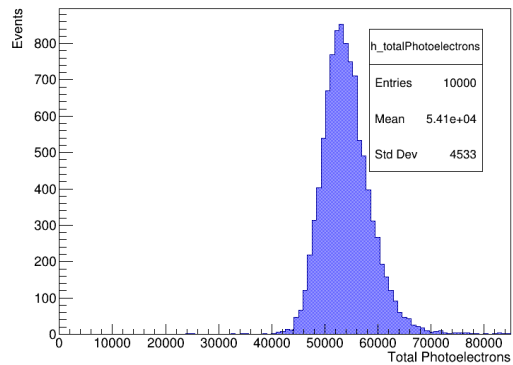
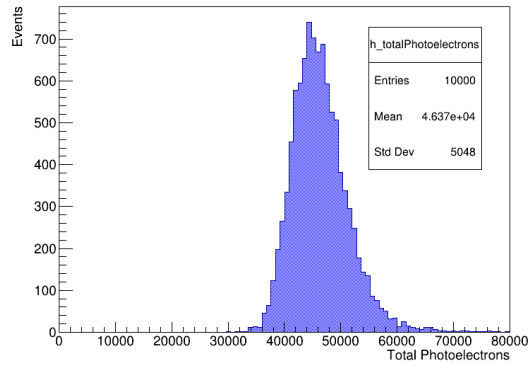
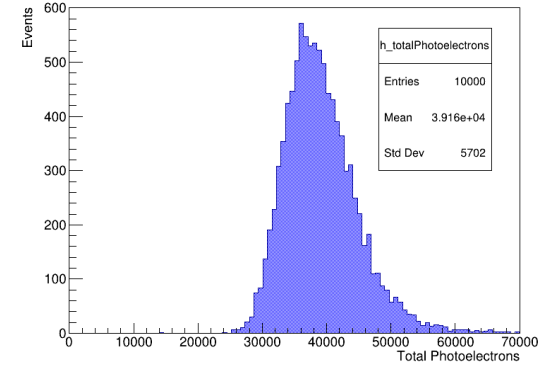
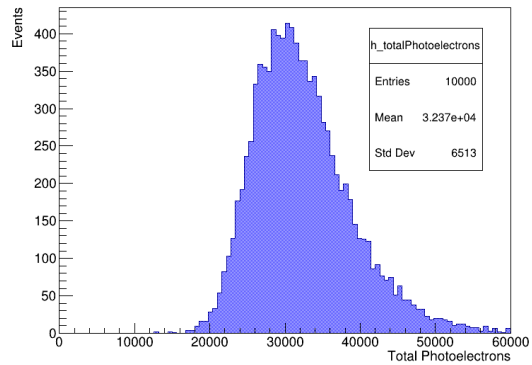
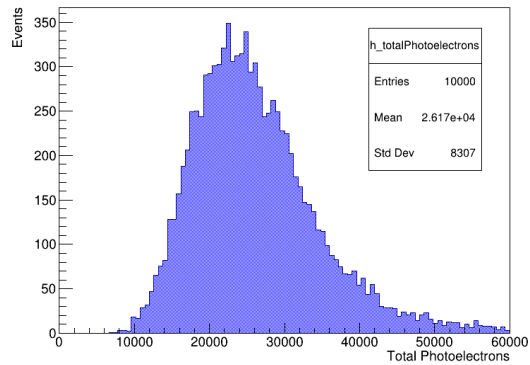
PWO - 30 photo-electrons/MeV

SiPM – 30% efficiency

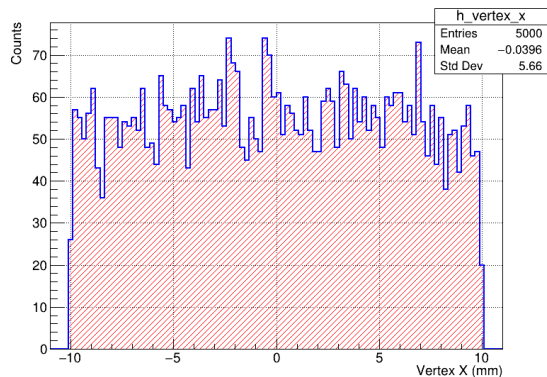
# Sum of Photoelectrons for Monoenergetic case



# Sum of Photoelectrons for DESY case



# Position Resolution analysis



Uniformly distributed  
x coordinate over the  
width of central crystal

Logarithmic weights

$$x = \frac{\sum_i w_i x_i}{\sum_i w_i} \quad w_i = \max \left\{ 0, \left[ W_0 + \ln \left( \frac{E_i}{E} \right) \right] \right\}$$

Rec. vs Gen. X positions

