

EMC in Fast Simulation

Chih-hsiang Cheng

Caltech

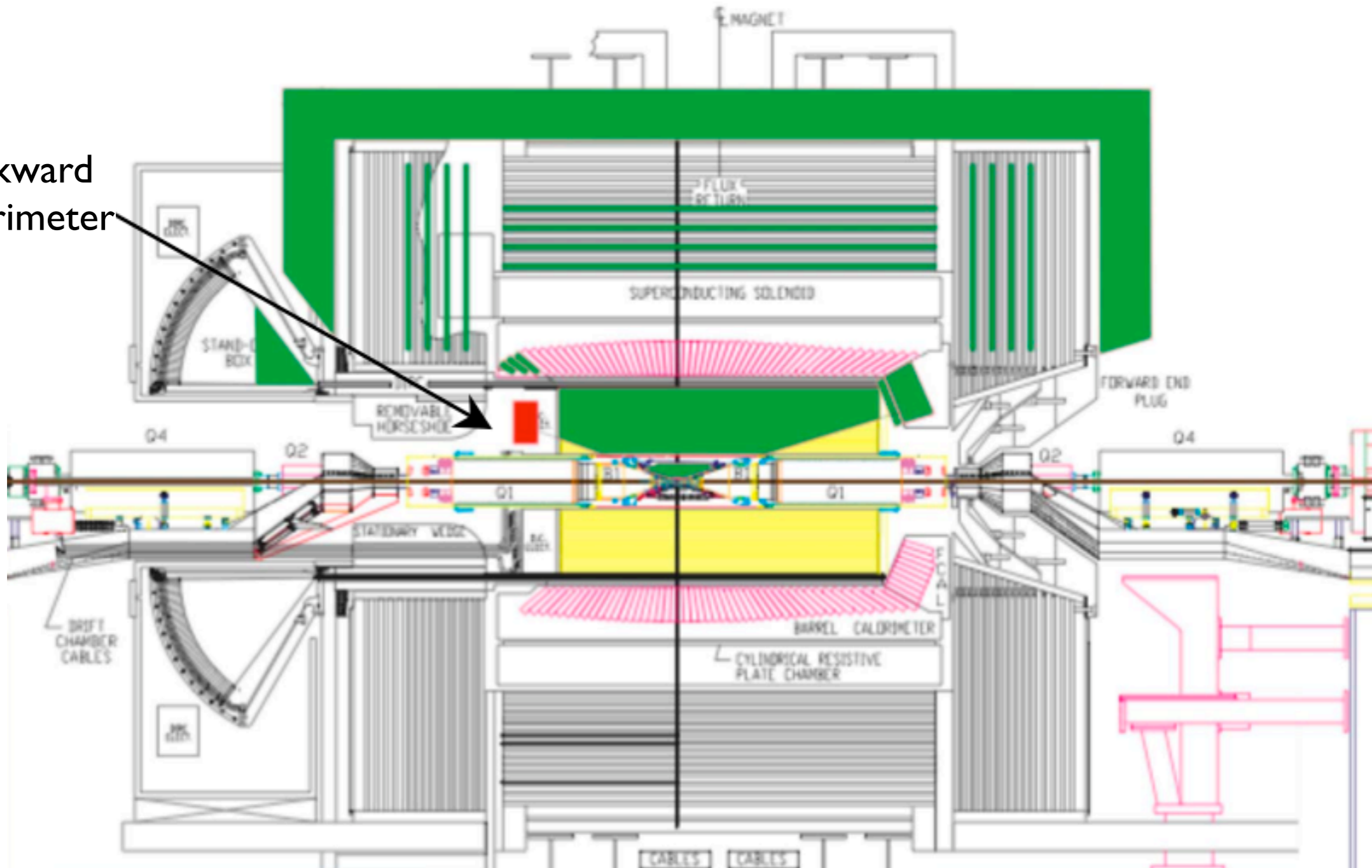
2009.02.16

SuperB Workshop, LAL Orsay, France



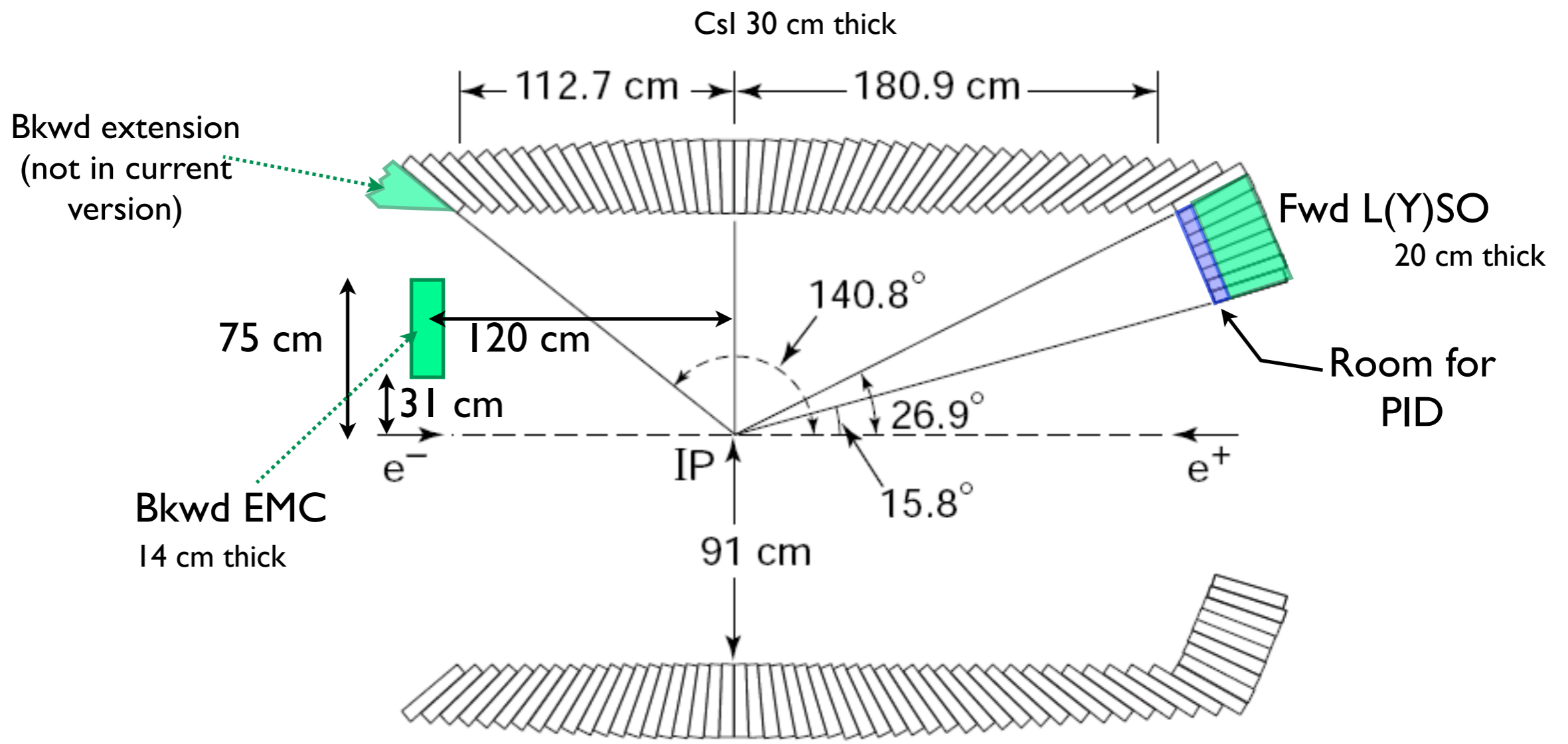
SuperB baseline design

Backward
Calorimeter

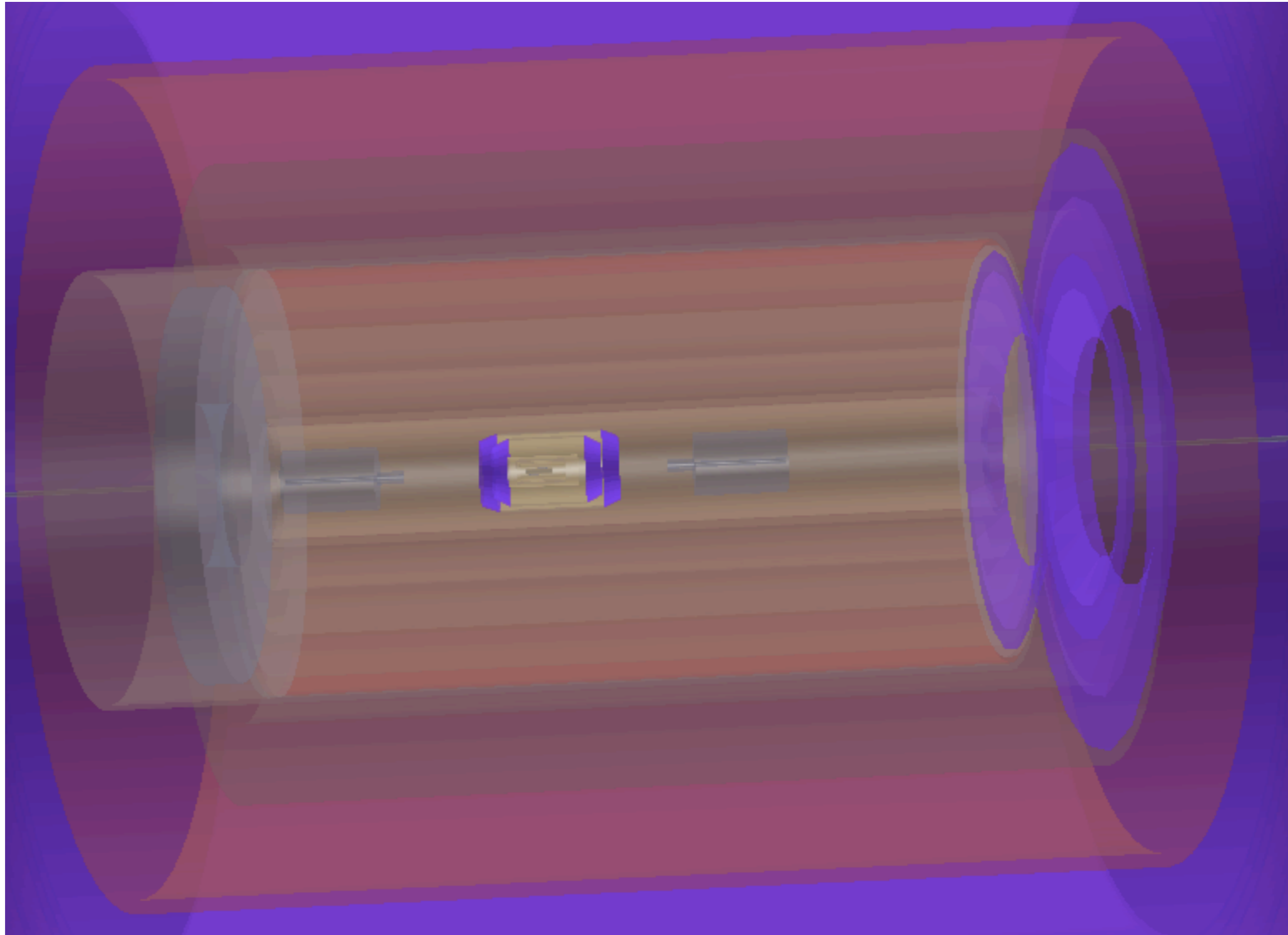


Geometry in the fastsim

- Barrel: cylinder. Fwd endcap: cone. Bkwd endcap: disk.
- 2D representation, with thickness used to calculate interaction probability and energy loss.
- Uniform $\Delta\varphi$; uniform $\Delta\theta$ in endcaps; uniform $r\Delta\theta$ in barrel.



Display



Proposed backward calorimeter

- Behind DCH I propose to place Pb-scintillator sampling calorimeter

- 2.8 mm thick Pb plates $\rightarrow 1/2 X_0$

- 3.0 mm thick scintillator tiles

- Sizes vary from 3.8 cm \times 3.8cm \rightarrow 7.8 cm \times 7.8 cm ($R_M \sim 6.0$ cm)

- cylindrical geometry, $r_i=0.31$ m, $r_o=0.75$ m

Pb: $R_M=1.5$ cm

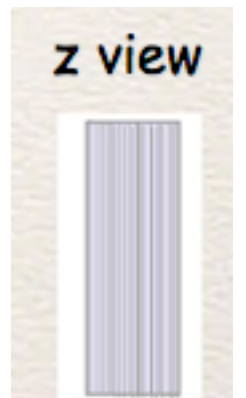
\rightarrow coverage $\sim 300\text{mr}$

- 24 planes with thickness of $12X_0$

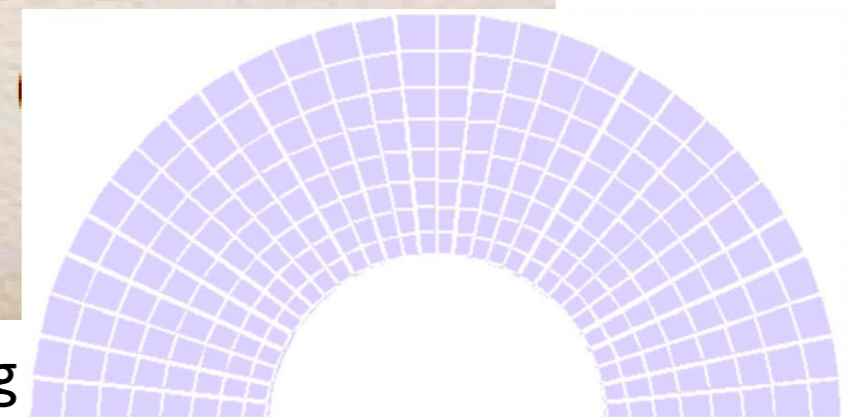
- scintillator is segmented into tiles, size increasing outwards

\rightarrow total: 11,520 channels

- Scintillator tiles are read out with WLS fibers coupled to a SIPM



Not projective



G. Eigen, SuperB meeting Elba, 31/05/2008

Gerald Eigen

8 ring, 60 tiles/ring

New materials in MaterialsList.data

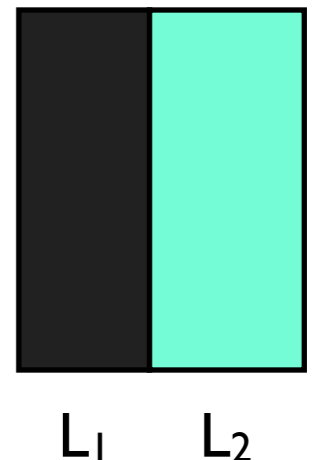
- Forward LSO: Lu_2SiO_5 ; LYSO: add ~5% Yttrium
 - $d = 7.4 \text{ g/cm}^3$; $X_0^* = 1.14 \text{ cm}$; $\lambda_l^* = 21 \text{ cm}$. ($R_M = 2.07 \text{ cm}$)
- Backward Pb-scintillator plates: (2.8mm Pb + 3.0mm scintillator tiles)x24
 - Treated as a homogeneous material, rather than sampling plates.
 - 2.8mm Pb = $0.5 X_0$; 3mm Polystyrene $\sim 0.007 X_0$.

	Z	A	X0	d	X_0^* (cm)	λ_l^* (cm)	R_M (cm)
Pb	82	207.2	6.37	11.35	0.56	17.6	1.6
Polystyrene	3.5	6.5	43.8	1.06	41.3	77.1	6.8

$$\langle d \rangle = (d_1 L_1 + d_2 L_2) / (L_1 + L_2) = 6.03$$

$$\langle A \rangle = 69.3 \quad \langle Z \rangle = 28.1$$

$$\langle X_0^* \rangle = \frac{L_1 + L_2}{L_1/X_{0;1}^* + L_2/X_{0;2}^*} = 1.14 \quad \langle X_0 \rangle = 6.87$$



New materials continued

- Moliere radius: Since the scintillator only contributes to a small amount of radiation length, only Pb contributes to creating shower particles. The spacing between Pb makes the shower profile larger, simply from geometric effect.

$$\langle R_M \rangle \simeq R_M^{\text{Pb}} \frac{L_1 + L_2}{L_1} = 3.3\text{cm}$$

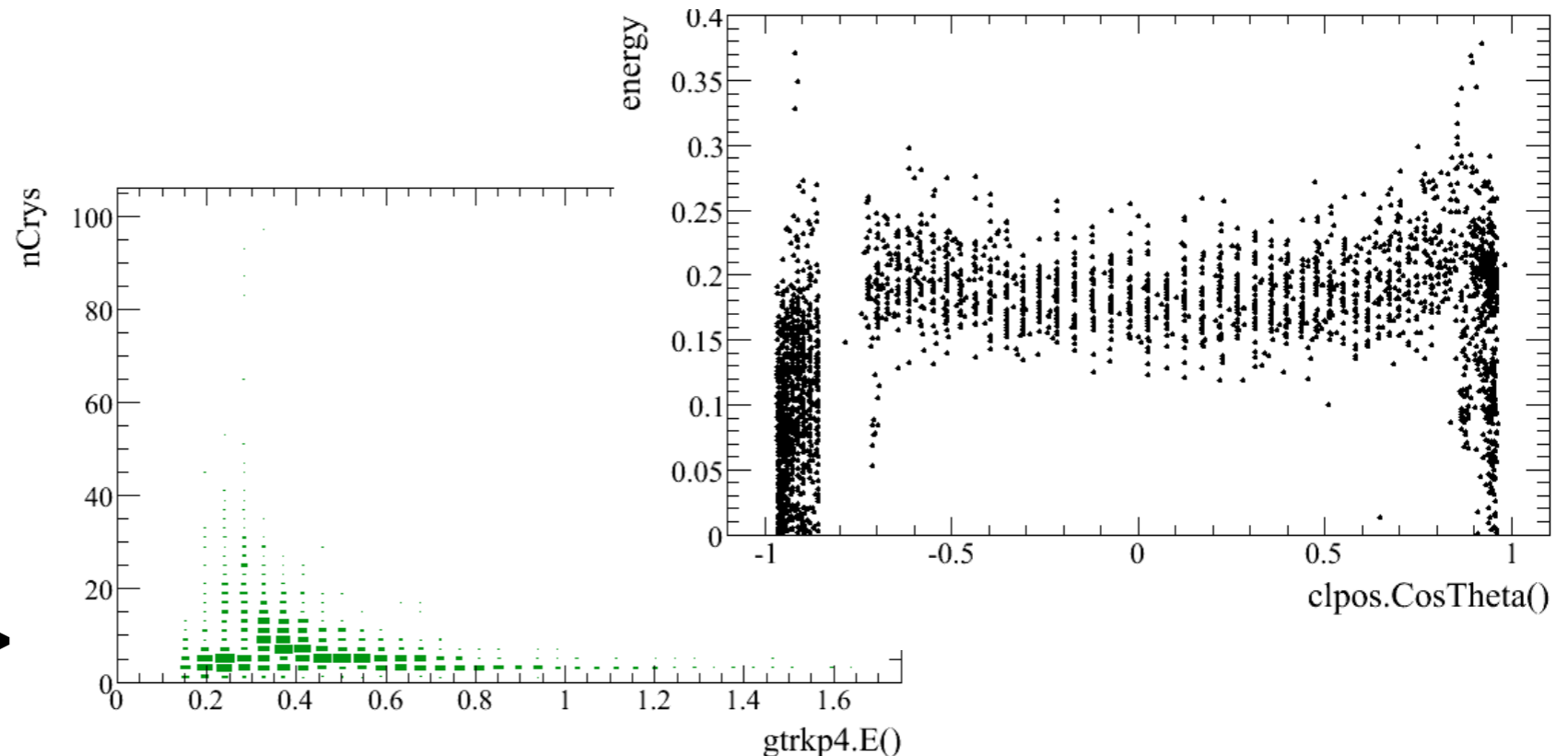
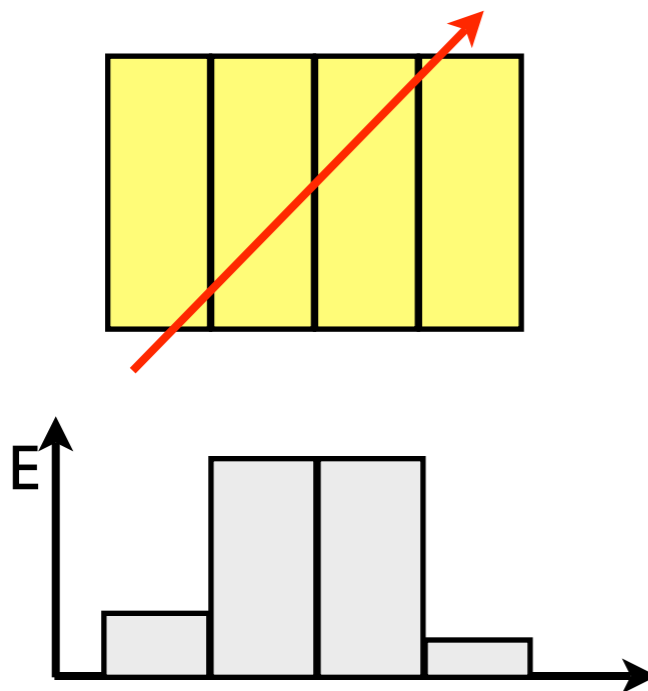
EMC clusters

- An EMC cluster is represented by the class `PacEmcCluster` (inherit from `AbsRecoCalo`), which contains a list of `PacEmcDigi`. The latter represents the energy deposition in a single crystal.
- Both classes mimic the respective classes in BaBar, but no calibration, timing, and data flow information are represented.

Ionization

- If a particle does not shower in the EMC (effects: normal, stop, interact, brems, compton, convert), we simply distribute the energy loss to the crystals it passes through. Energy is proportional to the path length in each crystal.
- Curving inside the EMC is ignored.
- Energy in each crystal is then smeared according to $\frac{\sigma_E}{E} = \frac{a}{E^d} \oplus b$

a, b, d are configuration parameters



EM shower

- The lateral shower development is assumed to be symmetric
- On average 10% of the deposited energy lies outside R_M , and about 1% outside $3.5 R_M$.
- The radial distribution can be modeled phenomenologically with

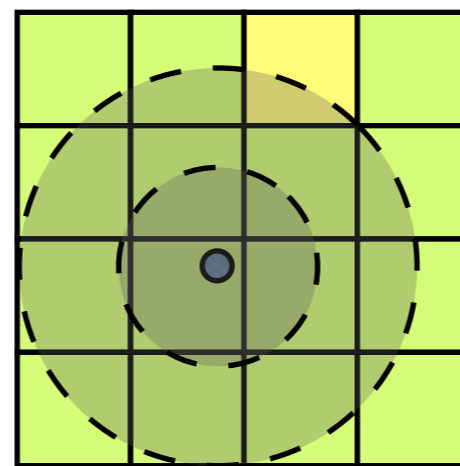
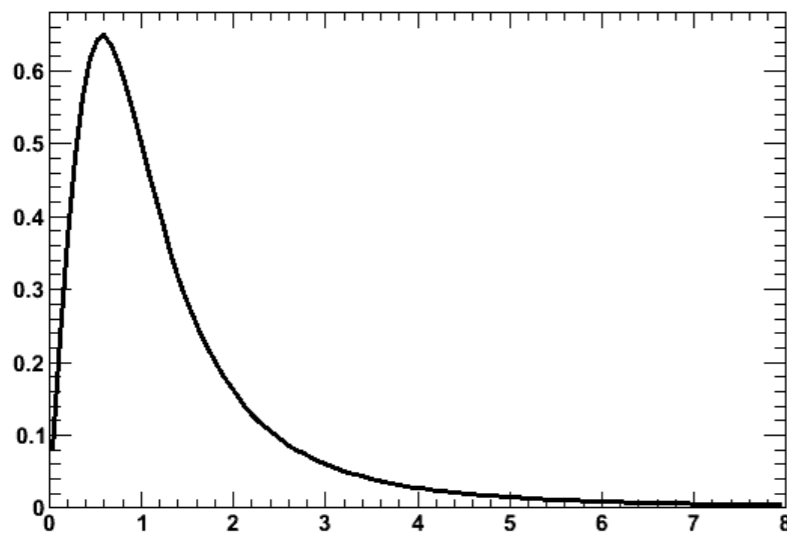
$$f(r) = \frac{2rR^2}{(r^2 + R^2)^2}$$

[PDG2008 Sec. 27.5, or NIM A290, 469]

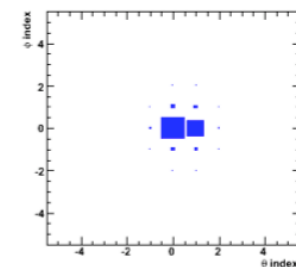
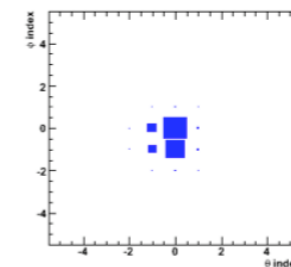
$$\int_r^\infty f(r') dr' = \frac{R^2}{r^2 + R^2}$$

$$R_M = 3R$$

$2*x/(x^2+1)/(x^2+1)$

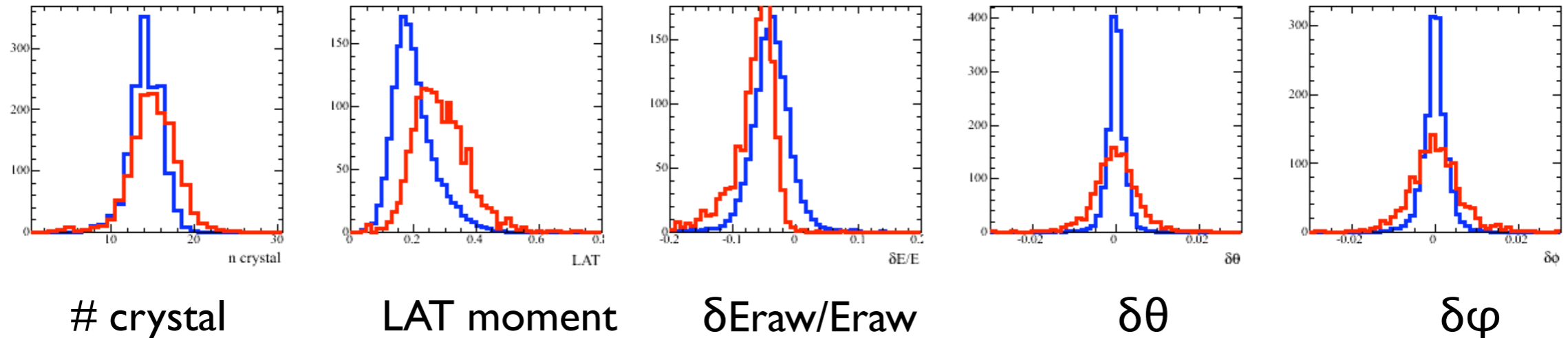


- R_M is allowed to fluctuate, so do energy in each crystal and eccentricity (axes along θ/φ , no rotation).

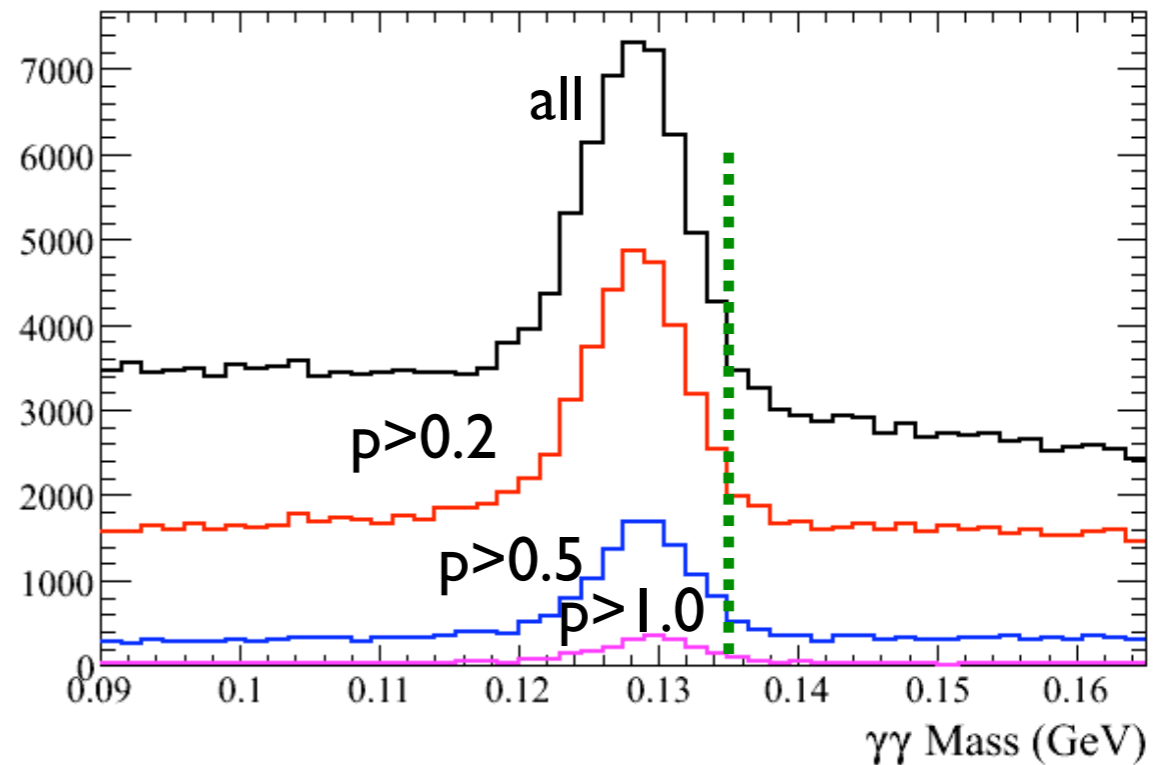
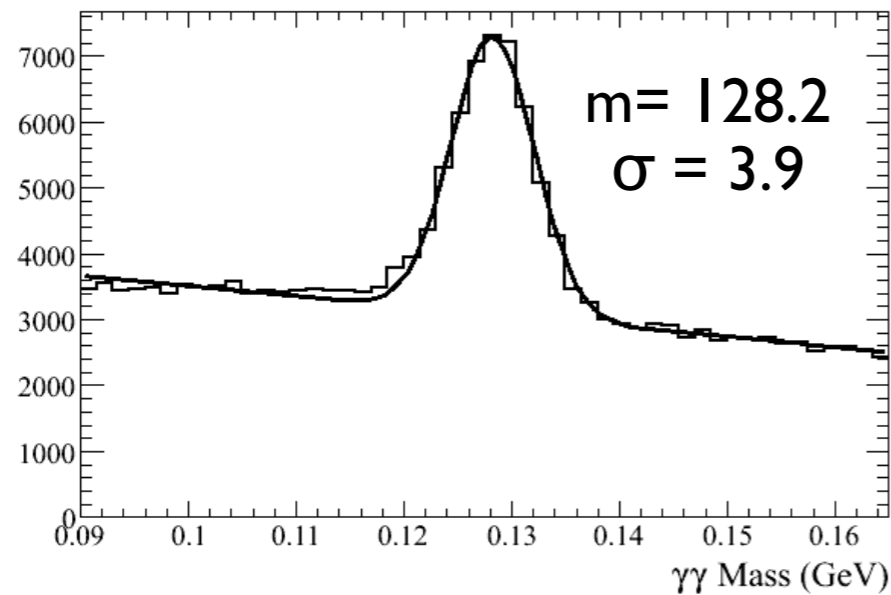


Performance

- One-GeV photons: Blue= **FastSim** (BaBar config); Red= **BaBar full Sim**



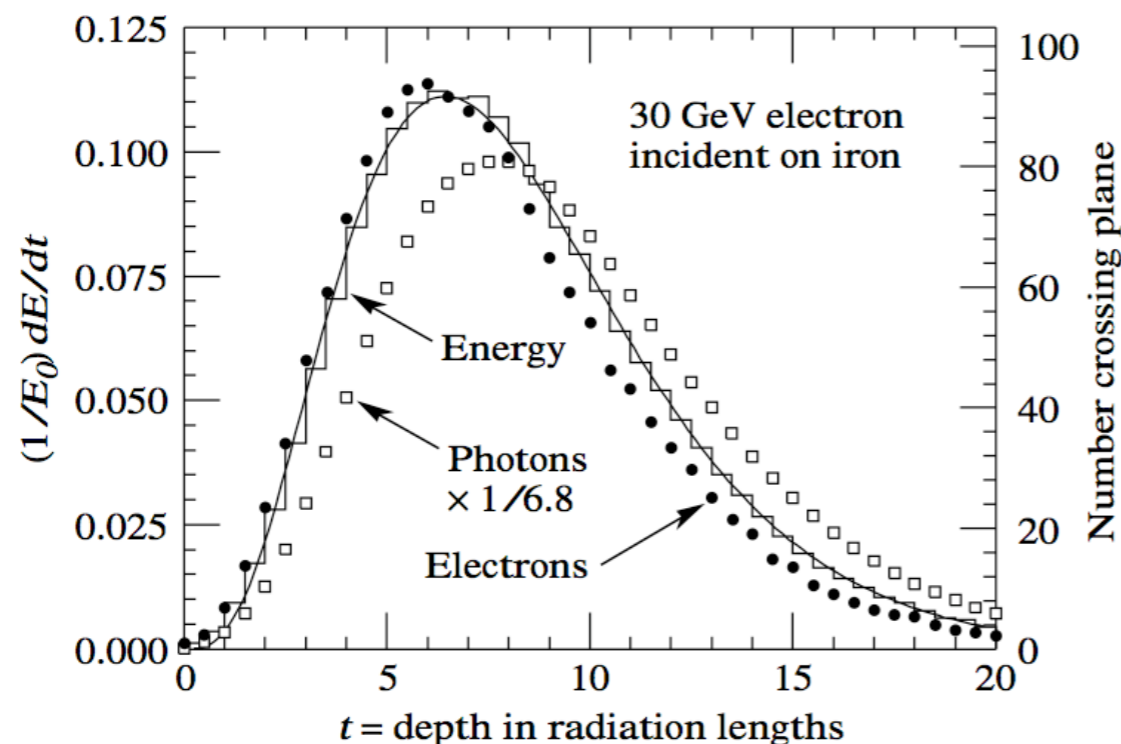
- π^0 from BB generic simulation



Need calibration. Resolution too high.

Longitudinal shower profile

- Need to know how much energy should be deposited on average, given the particle energy and the radiation lengths in and in front of the EMC, before creating clusters.
- Not in PacEmc, but calculated in PacSim (D. Brown)
- Profile depends on the material atomic number Z .
- Tricky to model when shower transits from one material to another. But important for detector study.

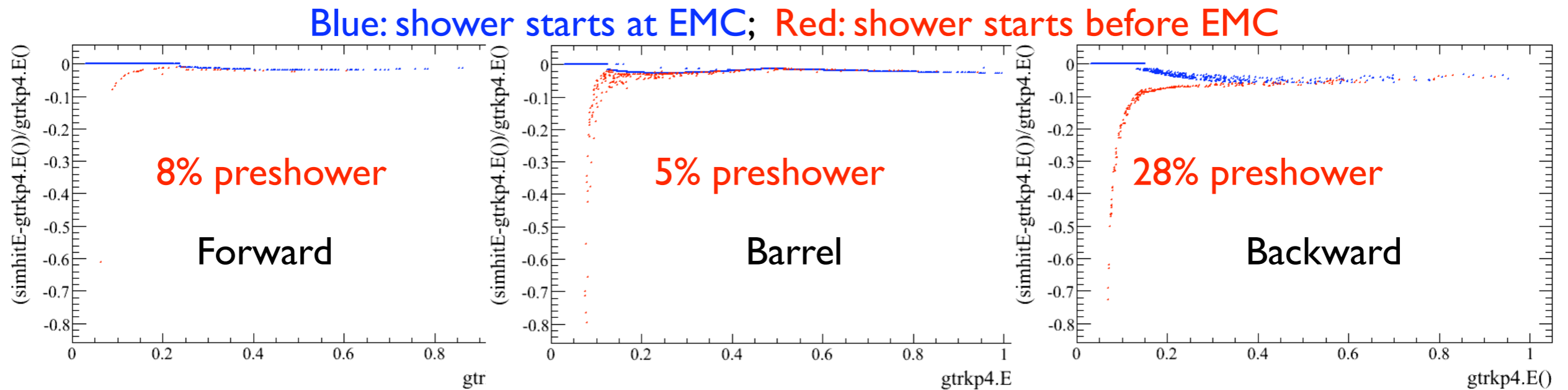


$$\frac{dE}{dt} = E_0 b \frac{(bt)^{a-1} e^{-bt}}{\Gamma(a)}$$

The integral has no closed-form solution.

We use numerical integration.

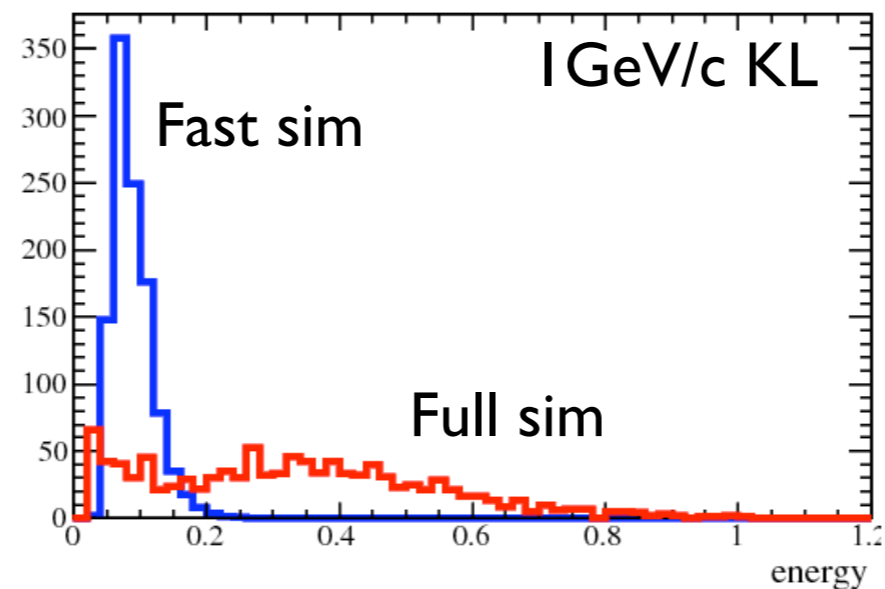
Expected energy deposition



- Except for very low energy, up to several percent energy leaks out of the back side.
- See energy loss due to preshower for low energy photons.
- Need to check against the full simulation.

Hadronic shower

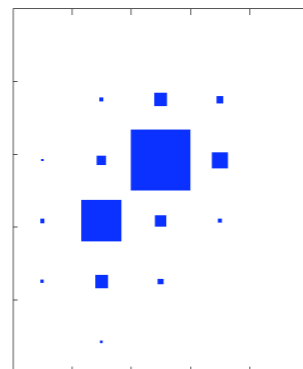
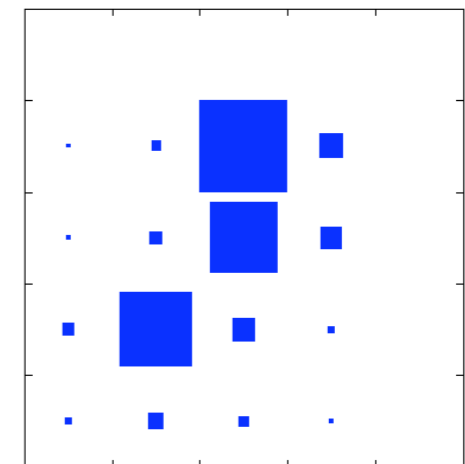
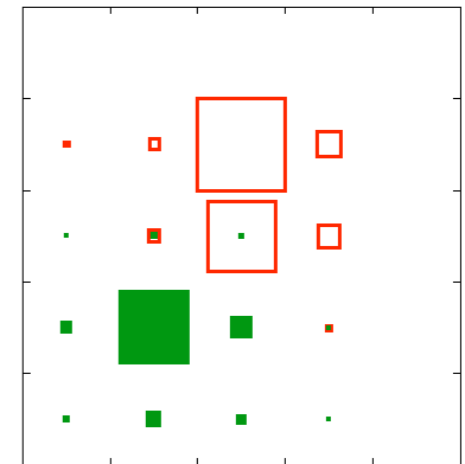
- Similar way to calculate the longitudinal integral, replacing radiation length with interaction length. However, it grossly underestimates the energy deposition.



- Hadronic showers are irregular and difficult to model with simple parametrizations.
- Use random walk to navigate through crystals and create large fluctuation to create irregular patterns in a cluster.

Cluster merging

- Merging is straight forward, simply adding energies in each crystal.
- There is no cluster reconstruction process. We know which crystals belong to the same cluster to begin with.
- We only merge clusters that are close enough to produce a single bump cluster.
 - #bumps = #local maxima
 - local maxima \equiv crystal energy higher than its neighboring eight.
 - This function hasn't been turned on in V0.0.2.



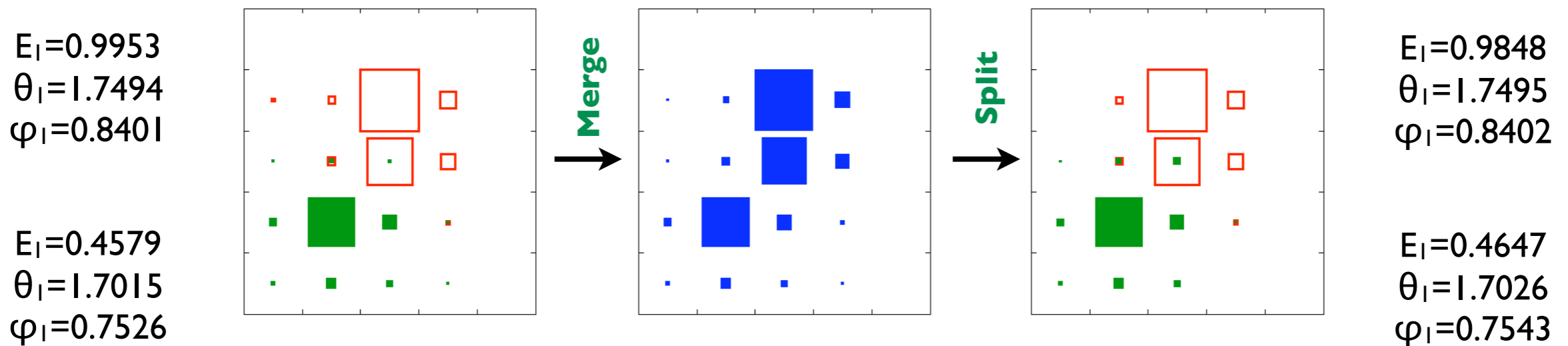
Is this one- or two-bump?

Cluster splitting

- Need to split clusters to simulate hadronic shower split-offs.
- Split clusters that have more than one local maximum.
- Assign a weight to each crystal j for each new one-bump clusters α :

$$w_j^{(\alpha)} = \frac{E^{(\alpha)} e^{-2.5|\vec{r}^{(\alpha)} - \vec{r}_j|}}{\sum_{\alpha} E^{(\alpha)} e^{-2.5|\vec{r}^{(\alpha)} - \vec{r}_j|}} \quad E^{(\alpha)} = \sum_j w_j^{(\alpha)} E_j \quad \sum_{\alpha} w_j^{(\alpha)} = 1$$

- Calculate weights and then new $E^{(\alpha)}$ and $r^{(\alpha)}$ iteratively until converge.



This function hasn't been turned on in V0.0.2.

More to do

- Ability to divide forward endcap into two regions (Csl, and LYSO).
- Track-cluster matching information.
- Truth association:
 - 1 GTrack \leftrightarrow n clusters (split); n GTracks \leftrightarrow 1 cluster (merge)
- Validation; QA; parameter tuning.
- Energy calibration.
- Hadronic shower.
- Background frame mixing.