Marlin-based Algorithm for Geometry-Independent Clustering

MAGIC: v01-02



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Order of service

- Reminder of the (3-stage) clustering algorithm.
- Where to get the code and how to get started with it.
- Studies of charged/neutral shower separation at normal incidence.
- Studies of cluster reconstruction vs solid angle in full detector simulation.
- Running the algorithm on the Ecal prototype data.

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• Summary.

The algorithm and how to use it

The algorithm and how to use it...

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Clustering with MAGIC: stage 1

- Form coarse clusters by *tracking* closelyrelated hits *layer-by-layer* through the calorimeter:
 - for a candidate hit in a given layer, *l*, minimise the distance, *d*, w.r.t all (already clustered) hits in layer *l*-1;
 - if d < distMax for minimum d, assign candidate hit to same cluster as hit in layer *l*-1 which yields minimum;
 - if not, repeat with all hits in layer *l*-2, then, if necessary, layer *l*-3, *etc*., right through to layer *l*-layersToTrackBack;
 - after iterating over all hits in layer *l*, seed new clusters with those still unassigned, grouping those within proxSeedMax of hit of highest remaining density into same seed;
 - assign a direction cosine to each layer l hit:
 - if in Ecal, calculate density-weighted centre of each cluster's hits in layer *I*; assign a direction cosine to each hit along the line joining its cluster's centre in the seed layer (or (0,0,0) if it's a seed) to its cluster's centre in layer *I*;
 - if in Hcal, assign a direction cosine to each hit along the line from the hit to which each is linked (or (0,0,0) if it's a seed) to the hit itself;
 - iterate outwards through layers.



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Clustering with MAGIC: stage 2

- Try to merge backward-spiralling track-like cluster-fragments with the forward propagating clusters to which they belong:
 - for each hit in the terminating layer, *l*, of a candidate cluster fragment, calculate the distance, *p*, to each hit in nearby clusters in the same layer, and the angle, γ, between their direction cosines;
 - loop over all pairs of hits;
 - if, for any pair, both:
 - p < proxMergeMax and</pre>
 - cos γ < cosGammaMax

are satisfied, merge clusters together into one;

- iterate over clusters.



Clustering with MAGIC: stage 3

- Try to merge low multiplicity cluster "halos" (hit multiplicity < clusterSizeMin) which just fail the stage 1 cluster-continuation cuts:
 - for the hit of highest density in the seed layer, l, of a low multiplicity cluster, minimise the angle, β , w.r.t all hits in layer l-1;
 - if tan β < tanBetaMax for minimum β , merge the clusters containing the repsective hits into one;
 - if not, repeat with all hits in layer *l*-2, then, if necessary, layer *l*-3, *etc.*, right through to layer *l*-layersToTrackBack;
 - if still not, repeat above steps with the candidate hit in the seed layer of the low multiplicity cluster of next highest density, *etc.*;
 - if still not, merge the low multiplicity cluster into the nearest cluster with hits in the same layer as the low multiplicity cluster's seed layer, provided the two clusters contain hits separated by
 s < proxMergeMax;
 - iterate over clusters.



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Code organisation within LCIO/MARLIN

- Code structured as a series of 5+1 MARLIN "processors", together with a steering file: cluster.steer (read at run-time).
- Reads hits collections from LCIO file, adds LCIO clusters collections (essentially pointers back to component hits) and writes everything to new LCIO output file.
- Processors to do the reconstruction:
 - CalorimeterConfigurer
 - \rightarrow allows user to define geometrical layout of calorimeter;
 - CalorimeterHitSetter

 \rightarrow applies hit-energy threshold and adds pseudolayer and pseudostave indices to hits collection (encoded in CellIDI akin to encoding of layer and stave indices in CellIDO) as well as hit weights (= local hit density);

- CalorimeterStage1Clusterer
 - \rightarrow performs coarse cluster reconstruction;
- CalorimeterStage2Clusterer
 - \rightarrow recovers backward-spiralling track-like cluster fragments;
- CalorimeterStage3Clusterer
 - \rightarrow recovers low multiplicity cluster fragments.
- Additional processor to access MC truth (if simulation):
 - CalorimeterTrueClusterer

 \rightarrow constructs true clusters, where a true cluster is considered to comprise all hits attributable to either:

- (i) the same generator primary or any of its non-backscattered progeny, or
- (ii) the same backscattered daughter or any of its non-backscattered progeny.

User-controlled steering with MARLIN

• Detector parameters and clustering cuts set in cluster.steer (e.g. Mokka DO9 model): ProcessorType CalorimeterConfigurer

detectorType	full	<pre># "full" => barrel+endcaps; "prototype" => layers perp'r to +z</pre>
iPx	0.	<pre># x-coordinate of interaction point (in mm)</pre>
iPy	0.	# y-coordinate of interaction point (in mm)
iPz	0.	<pre># z-coordinate of interaction point (in mm)</pre>
ecalLayers	40	# number of Ecal layers
hcalLayers	40	# number of Hcal layers
barrelSymmetry	8	# degree of rotational symmetry of barrel
phi_1	90.0	<pre># phi offset of barrel stave 1 w.r.t. x-axis (in deg)</pre>
ProcessorType CalorimeterHitSet	ter	
ecalMip	0.000150	# Ecal MIP energy (in GeV)
hcalMip	0.000004	# Hcal MIP energy (in GeV)
ecalMipThreshold	0.3333333	<pre># Ecal hit-energy threshold (in MIP units)</pre>
hcalMipThreshold	0.3333333	<pre># Hcal hit-energy threshold (in MIP units)</pre>
ProcessorType CalorimeterStage1	Clusterer	
layersToTrackBack_ecal	3	<pre># number of layers to track back in Ecal</pre>
layersToTrackBack_hcal	3	<pre># number of layers to track back in Hcal</pre>
distMax_ecal	20.0	# distance cut in Ecal (in mm)
distMax_hcal	30.0	# distance cut in Hcal (in mm)
proxSeedMax_ecal	14.0	# maximum cluster-seed radius in Ecal (in mm)
proxSeedMax_hcal	50.0	# maximum cluster-seed radius in Hcal (in mm)
ProcessorType CalorimeterStage2	Clusterer?	
proxMergeMax_ecal	20.0	# Ecal proximity cut for cluster merging (in mm)
proxMergeMax_hcal	30.0	# Hcal proximity cut for cluster merging (in mm)
cosGammaMax	0.5	<pre># angular cut for cluster merging</pre>
ProcessorType CalorimeterStage3	Clusterer	
clusterSizeMin	10	# minimum cluster size to avert potential merging
layersToTrackBack_ecal	39	# number of layers to track back in Ecal for merging
layersToTrackBack_hcal	79	# number of layers to track back in Hcal for merging
tanBetaMax	6.0	<pre># angular cut for cluster merging</pre>
proxSeedMax_ecal	400.0	# Ecal proximity cut for cluster merging (in mm)
proxSeedMax_hcal	400.0	# Hcal proximity cut for cluster merging (in mm)
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Getting started with MAGIC

- Install LCIO (\geq v01-05) and MARLIN (\geq v00-07).
- Download *MAGIC* tar-ball from
 - http://www.hep.phy.cam.ac.uk/~ainsley/MAGIC/MAGIC-v01-02.tar.gz
- Two directories and a **README** file (read this first!).
- The clustering directory contains the cluster-reconstruction (and cluster-truth) code (i.e. all processors and steering file mentioned earlier).
- Takes .slcio input files containing CalorimeterHits (data) or SimCalorimeterHits (MC):
 - must be generated with hit-positions stored, *i.e.* RCHBIT_LONG=1 (data) or CHBIT_LONG=1 (MC);
 - collection names must contain the string "ecal" or "hcal" (in upper or lower case, or in some combination of these) to identify the type of hit (for energy-threshold application).
- Produces .slcio output file with cluster-related collections added:
 - CalorimeterHits \Rightarrow hits above energy threshold;
 - CalorimeterHitRelationsToSimCalorimeterHits (MC only) \Rightarrow pointers to original simulated hits;
 - CalorimeterStage1Clusters \Rightarrow clusters after stage 1 of algorithm;
 - CalorimeterStage2Clusters \Rightarrow clusters after stage 2 of algorithm;
 - CalorimeterStage3Clusters \Rightarrow clusters after stage 3 of algorithm;
 - CalorimeterTrueClusters (MC only) \Rightarrow true clusters;
 - CalorimeterTrueClusterRelationsToMCParticles (MC only) \Rightarrow pointers to original MC particles.
- The examples directory contains example analysis code which performs simple manipulations with the clusters (e.g. processors which add calibrated energies to clusters, produce the plots shown earlier, calculate the reconstruction quality... and an accompanying steering file).

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Charged/neutral shower separation

Charged/neutral shower separation...

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Charged/neutral shower separation studies

- Fire nearby charged/neutral particles into calorimeter.
- Perform standalone clustering on calorimeter hits with *MAGIC*.
- Extrapolate helix from charged track through calorimeters.
- Associate clusters/cluster fragments with charged particle if seeded within pad-size (= 1 cm) of projected helical trajectory.
- Remove corresponding calorimeter hits from further consideration; assume remainder to be the neutral shower.
- Apply energy calibration to leftover hits to reconstruct neutral particle energy.

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π^+/γ separation: D09 model (1)



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- Black cluster matched to charged track.
- Red cluster left over as neutral $\Rightarrow \gamma$ energy well reconstructed.

π^+/γ separation: D09 model (2)



- 1k single γ at 5 GeV/*c*.
- Fit Gaussian to energy distribution, calibrated according to:

$$E = \alpha [(E_{\text{Ecal; 1-30}} + 3E_{\text{Ecal; 31-40}})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}].$$

- Fix factors α , 20 by minimising χ^2 /dof.
- $\sigma/J\mu \sim 14\% JGeV$.

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- 1k γ with nearby π^+ (at 10, 5, 3, 2 cm from γ).
- Peak of photon energy spectrum well reconstructed; improves with separation.
- Tail at higher $E \rightarrow$ inefficiency in π^+ reconstruction (next page...).
- Spike at E = 0 below 3 cm \rightarrow clusters not distinguished.

π^+/γ separation: D09 model (3)



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π^+/n separation: D09 model (1)



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Black cluster matched to charged track.
Red cluster left over as neutral ⇒ n energy well reconstructed.

π^+/n separation: D09 model (2)



- 1k single n at 5 GeV/c.
- Fit Gaussian to energy distribution, calibrated according to:

$$E = \alpha [(E_{\text{Ecal; 1-30}} + 3E_{\text{Ecal; 31-40}})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}]$$

- Fix factors α , 20 by minimising χ^2 /dof.
- $\sigma/J\mu \sim 73\% JGeV$.

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- 1k **n** with nearby π^+ (at 10, 5, 3, 2 cm from n).
- Peak of neutron energy spectrum well reconstructed; improves with separation.
- Spike at E = 0 even at 10 cm \rightarrow clusters not distinguished (next page...).

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π^+/n separation: D09 model (3)



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• Nothing left over as neutral \Rightarrow n

not reconstructed (*i.e.* E = 0).



π^+/γ separation: D09Scint model

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- 1k single γ at 5 GeV/c.
- Fit Gaussian to energy distribution, calibrated General trends much as for D09 model. according to:
- $E = \alpha[(E_{\text{Ecal; 1-30}} + 3E_{\text{Ecal; 31-40}})/E_{\text{Ecal mip}} + 5E_{\text{Hcal}}/E_{\text{Hcal mip}}].$ Fix factors α , 5 by minimising χ^2/dof .
- $\sigma/J\mu \sim 14\%$ JGeV (as for D09 model).

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- 1k γ with nearby π^+ (at 10, 5, 3, 2 cm from γ).

 π^+/n separation: D09Scint model



- 1k single n at 5 GeV/c.
- Fit Gaussian to energy distribution, calibrated according to:

$$E = \alpha [(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal mip}} + 5E_{\text{Hcal}}/E_{\text{Hcal mip}}]$$

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- Fix factors α , 5 by minimising χ^2 /dof.
- $\sigma/J\mu \sim 62\% JGeV$ (cf. 73% $\overline{J}GeV$ for D09 model).

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Events 35 Events 35 Entries 1000 Entries 1000 4.701 4.523 Mean Mean 30 30 RMS 2.391 RMS 2.571 25 25 20 20 — π⁺/n at ∞ — π*/n at ∞ 15 15 - π*/n at 10 cm π*/n at 5 cm 10 10 10 15 20 / GeV (reconstruction) 10 15 GeV (reconstruction) :*/n (5 GeV/c) - D09Scint Events 25 */n (5 GeV/c) - D09Scin ۵<u>5</u> Entries 1000 Entries 1000 3.158 Mean 3.987 Mean 30 30 RMS 2.924 RMS 3.084 25 25 20 20 — π⁺/n at ∞ — π⁺/n at ∞ 15 – π*/n at 3 cm π*/n at 2 cm 10 Þ 5 E_{neutr} 10 15 20 / GeV (reconstruction) 10 15 20 / GeV (reconstruction) 5 E_____

:*/n (5 GeV/c) - D09Scint

*/n (5 GeV/c) - D09Scin

- 1k **n** with nearby π^+ (at 10, 5, 3, 2 cm from n).
- General trends much as for D09 model.

π^+ /neutral cluster separability vs separation

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5 GeV/*c* π⁺/γ



- Fraction of events with photon energy reconstructed within 1,2,3σ generally higher for D09 than for D09Scint...
- ...and absolute γ resolution similar.

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5 GeV/*c* π⁺/n



- Fraction with neutron energy reconstructed within 1,2,3σ also generally higher for D09...
- ...*but*, absolute *n* resolution is better for D09Scint.

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Clustering vs detector solid angle

Clustering vs detector

solid angle...

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- μ- fired isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha [(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}] \text{ GeV}.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) at (0,0,0).



- Default clustering cuts $\rightarrow \mu^{-}$ track fragmented at $|\cos \theta| \sim 0.78$, $|\cos \theta| \sim 0.23$ and $\phi \sim 0.20$ -0.24 (cos θ -dependent) \Rightarrow algorithm needs to know some geometry to overcome this!
- Angular-dependent clustering cuts $\rightarrow \mu^{-}$ track reconstructed with ~100% efficiency \forall (θ , ϕ).
- What detector features do these regions correspond to?

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- μ^- at cos θ = -0.75 traverses Ecal barrel, Hcal barrel and Hcal endcap.
- Track breaks on crossing from barrel to endcap \rightarrow layers of active material "missing" in the gap.



- Relax layersToTrackBack_ecal cut for $0.81 < |\cos \theta| < 0.85$ and layersToTrackBack_hcal cut for $0.72 < |\cos \theta| < 0.85$ to prevent this.
- Design the detector with as small a barrel-endcap gap as possible!

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- μ^- at cos θ = -0.24 traverses Ecal barrel module 3, Hcal barrel module 3 and Hcal barrel module 2.
- Track breaks on crossing between barrel modules at $|z| \sim 0.56$ m \rightarrow active cells "missing" near the module edges.



- Relax distMax_ecal and distMax_hcal cuts for 0.18 < $|\cos \theta| < 0.28$ to prevent this.
- Much less severe, but similar, effect at $|z| \sim 1.68$ m (0.47 < cos θ < 0.65) treated in the same way.

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• Design the detector with as small an inter-module gap as possible!

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- μ^- at $\phi = 1.58\pi$ traverses Ecal barrel stave 5, Hcal barrel stave 5 and Hcal barrel stave 6.
- Track breaks on crossing between Hcal barrel staves at ϕ -(6× π /4) = π /8 = 0.39 (curves in B-field) \rightarrow active cells "missing" near the stave edges.



• Relax distMax_hcal and layersToTrackBack_hcal cuts for 0.36 < ϕ < 0.42 if $|\cos \theta|$ < 0.82 (Hcal barrel) to prevent this.

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- No problem in Ecal \rightarrow staves overlap.
- Design the Hcal with no pointing cracks (e.g. like the Ecal)!

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- μ- fired isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha [(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}] \text{ GeV}.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) at (0,0,0).



- Default clustering cuts $\rightarrow \mu^{-}$ track fragmented at $|\cos \theta| \sim 0.78$ (barrel/endcap overlap), $|\cos \theta| \sim 0.23$ (gap between barrel modules) and $\phi \sim 0.20$ -0.24 (gap between Hcal barrel staves).
- Angular-dependent clustering cuts $\rightarrow \mu^{-}$ track reconstructed with ~100% efficiency \forall (θ , ϕ).
- Does relaxing cuts near dead zones impact on charged/neutral cluster separability though?

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- π -fired isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha [(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}] \text{ GeV}.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) at (0,0,0).



- Default clustering cuts $\rightarrow \pi^{-}$ shower fragmented at $|\cos \theta| \sim 0.83$ (barrel/endcap overlap) and $|\cos \theta| \sim 0.23$ (gap between barrel modules).
- Angular-dependent clustering cuts $\rightarrow \pi^-$ shower reconstructed with improved efficiency.
- Does relaxing cuts near dead zones impact on charged/neutral cluster separability though?

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Detector scan: γ (10 GeV)

- γ fired isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha [(E_{Ecal; 1-30} + 3E_{Ecal; 31-40})/E_{Ecal mip} + 20N_{Hcal}] GeV.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) at (0,0,0).



• Default clustering cuts $\rightarrow \gamma$ shower fragmented at $|\cos \theta| \sim 0.85$ (Ecal barrel/endcap overlap).

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- Angular-dependent clustering cuts $\rightarrow \gamma$ shower reconstructed with improved efficiency.
- · Does relaxing cuts near dead zones impact on charged-neutral cluster separability though?

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Detector scan: n (10 GeV)

- n fired isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha [(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal mip}} + 20N_{\text{Hcal}}] \text{ GeV}.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) at (0,0,0).



- Default clustering cuts \rightarrow n shower fragmented at $|\cos \theta| \sim 0.83$ (barrel/endcap overlap).
- Angular-dependent clustering cuts \rightarrow n shower reconstructed with improved efficiency.
- · Does relaxing cuts near dead zones impact on charged-neutral cluster separability though?

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Detector scan: π/γ at 5 cm (10 GeV)

- π/γ fired 5 cm apart isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: E = α[(E_{Ecal; 1-30} + 3E_{Ecal; 31-40})/E_{Ecal mip} + 20N_{Hcal}] GeV.
 Fraction of event energy in 1:1 correspondence between reconstructed and "true" clusters plotted vs $|\cos \theta|$ and $vs \phi$ (folded into first octant: $0 \le \phi < \pi/4$) on entry to Ecal.



- Default clustering cuts \rightarrow shower reconstruction/separability harder near $|\cos \theta| \sim 0.83$ (barrel/endcap overlap).
- Angular-dependent clustering cuts \rightarrow improves single-particle reconstruction, but increases potential charged/neutral confusion (cuts relaxed).
- On balance, seems beneficial \rightarrow separability barely affected.

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Detector scan: π /n at 5 cm (10 GeV)

- π -/n fired 5 cm apart isotropically into (analogue) Si/W Ecal, (digital) rpc/Fe Hcal (Mokka "D09" model).
- Cluster energies calibrated according to: $E = \alpha[(E_{\text{Ecal}; 1-30} + 3E_{\text{Ecal}; 31-40})/E_{\text{Ecal} \min} + 20N_{\text{Hcal}}] \text{ GeV}.$
- Fraction of event energy in highest-energy reconstructed cluster plotted vs $|\cos \theta|$ and vs ϕ (folded into first octant: $0 \le \phi < \pi/4$) on entry to Ecal.



- Default clustering cuts \rightarrow shower reconstruction/separability harder near $|\cos \theta| \sim 0.83$ (barrel/endcap overlap) and $|\cos \theta| \sim 0.25$ (gap between barrel modules).
- Angular-dependent clustering cuts → improves single-particle reconstruction, but increases potential charged/neutral confusion (cuts relaxed).
- On balance, may again be beneficial, but need to be careful.
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Clustering the prototype data

Clustering the prototype data...

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Steering file for the prototype

• Detector parameters and clustering cuts set in cluster.steer:

ProcessorType CalorimeterConfigurer

detectorType	prototype	<pre># "full" => barrel+endcaps; "prototype" => layers perp'r to +z</pre>
iPx	0.	<pre># x-coordinate of interaction point (in mm)</pre>
iPy	0.	<pre># y-coordinate of interaction point (in mm)</pre>
iPz	-99999.	<pre># z-coordinate of interaction point (in mm)</pre>
ecalLayers	30	# number of Ecal layers
hcalLayers	40	# number of Hcal layers
barrelSymmetry	8	<pre># degree of rotational symmetry of barrel</pre>
phi_1	90.0	<pre># phi offset of barrel stave 1 w.r.t. x-axis (in deg)</pre>
ProcessorType CalorimeterHitSet	ter	
ecalMip	0.000150	# Ecal MIP energy (in GeV)
hcalMip	0.000004	# Hcal MIP energy (in GeV)
ecalMipThreshold	0.3333333	<pre># Ecal hit-energy threshold (in MIP units)</pre>
hcalMipThreshold	0.3333333	<pre># Hcal hit-energy threshold (in MIP units)</pre>
ProcessorType CalorimeterStage1	Clusterer	
layersToTrackBack_ecal	3	# number of layers to track back in Ecal
layersToTrackBack_hcal	3	# number of layers to track back in Hcal
distMax_ecal	20.0	<pre># distance cut in Ecal (in mm)</pre>
distMax_hcal	30.0	# distance cut in Hcal (in mm)
proxSeedMax_ecal	14.0	<pre># maximum cluster-seed radius in Ecal (in mm)</pre>
proxSeedMax_hcal	50.0	<pre># maximum cluster-seed radius in Hcal (in mm)</pre>
ProcessorType CalorimeterStage2	Clusterer	
proxMergeMax_ecal	20.0	# Ecal proximity cut for cluster merging (in mm)
proxMergeMax_hcal	30.0	# Hcal proximity cut for cluster merging (in mm)
cosGammaMax	0.5	<pre># angular cut for cluster merging</pre>
ProcessorType CalorimeterStage3	Clusterer	
clusterSizeMin	10	# minimum cluster size to avert potential merging
layersToTrackBack_ecal	39	# number of layers to track back in Ecal for merging
layersToTrackBack_hcal	79	# number of layers to track back in Hcal for merging
tanBetaMax	6.0	<pre># angular cut for cluster merging</pre>
proxSeedMax_ecal	400.0	# Ecal proximity cut for cluster merging (in mm)
proxSeedMax_hcal	400.0	# Hcal proximity cut for cluster merging (in mm)
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Prototype data (Run 100121) : e⁻ (1 GeV)



- 14 layers (analogue) Si/W Ecal; > 50k 1 GeV e- events.
- Default clustering cuts \rightarrow events generally reconstruct as single clusters (no tracking info used).
- On average, **98.93** ± **0.03** % of event energy contained in highest energy reconstructed cluster (cluster energies calibrated according to: $E = \alpha(E_{\text{Ecal}; 1-10} + 2E_{\text{Ecal}; 11-14})$ GeV).

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Conclusion

Conclusion...

Summary & outlook

 Current version of Marlin-based Algorithm for Geometry-Independent Clustering available from:

http://www.hep.phy.cam.ac.uk/~ainsley/MAGIC/MAGIC-v01-02.tar.gz

- Will also put into CVS.
- Compliant with LCIO (\geq v01-05) / MARLIN (\geq v00-07) \Rightarrow input parameters (set at run-time) kept distinct from reconstruction (pre-compiled).
- Code straightforwardly applicable to any detector geometry comprising an *n*-fold rotationally symmetric barrel closed by endcaps \rightarrow just need to specify *n*, barrel orientation, and layer positions as input.
- User specifies geometry and clustering cuts (user-defined angular-dependence in next version) at run-time.

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- Algorithm can be used to compare different calorimeter designs straightforwardly (early hints of a preference for rpc over scintillator for Hcal using Mokka models).
- Please try it out!

The end

That's all folks...

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Generalising the calorimeter (1)



- Layer index changes discontinuously at barrel/endcap boundary.
- On crossing, jumps from *l* to 1 (first Ecal layer).



- Define a "*pseudolayer*" index based on projected intersections of physical layers.
- Index varies smoothly across boundary.
- Pseudolayer index = layer index, *except* in overlap region.

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Generalising the calorimeter (2)



- Layer index changes discontinuously at boundary between overlapping barrel staves.
- On crossing, jumps from *l* to 1 (first Ecal layer.



- Again, define "*pseudolayer*" index from projected intersections of physical layers.
- Again, index varies smoothly across boundary.
- Again, pseudolayer index = layer index, except in overlap region.

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Generalising the calorimeter (3)



- Define a "*pseudostave"* as a plane of parallel pseudolayers.
- "Pseudobarrel" pseudostaves meet boundaries with left- and right-hand "pseudoendcap" pseudostaves along 45° lines (if layer-spacings equal in barrel and endcaps).

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- "Pseudobarrel" pseudostaves meet boundaries with other "pseudobarrel" pseudostaves along 360°/2n lines (for an n-fold rotationally symmetric barrel).
- Calorimeter divides naturally into n+2 pseudostaves.

Generalising the calorimeter (4)

- Code recasts any layered calorimeter composed of a rotationally symmetric barrel closed by two endcaps into this standard, generalised form comprising layered shells of rotationallysymmetric *n*-polygonal prisms, coaxial with *z*-axis.
- Layers and staves from which calorimeter is built translated into pseudolayers and pseudostaves with which algorithm works.
- Only required inputs as far as algorithm is concerned are:
 - **barrelSymmetry** = rotational symmetry of barrel (*n*);
 - phi_1 = orientation of pseudobarrel pseudostave 1 w.r.t. x-axis;
 - distanceToBarrelLayers[ecalLayers+hcalLayers+2]
 - = layer positions in barrel layers ("+2" to constrain inside edge of first pseudolayer and outside edge of last pseudolayer); and
 - distanceToEndcapLayers[ecalLayers+hcalLayers+2]
 - = layer positions in endcap layers;
 - \rightarrow as geometry-independent as it's likely to get!

How the generalised detector shapes up



- Solid blue lines aligned along real, physical, sensitive layers.
- Dot-dashed magenta lines bound shell containing hits with same *pseudolayer* index, *l*.
- *Pseudostaves* automatically encoded by specifying n, ϕ_1 and R_l and Z_l ($\forall l$).

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Cluster-tracking between pseudolayers

From the pseudobarrel

From the pseudoendcap



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5 GeV π^+ event: 3 stages of clustering



 One backward-spiralling track and several halo clusters surround principal cluster. • Backward-spiralling track merged with principal cluster.

• Halo clusters merged with principal cluster.

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• Reconstruction works successfully not only for *intra*-stave, but also for *inter*-stave clusters

(e.g. black truth cluster spanning barrel staves 5+6 and the RH endcap correctly reconstructed).

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Code organisation within LCIO/MARLIN

• Layer positions set (for convenience) in CalorimeterConfigurer.cc:

// Create collections to store the barrel and endcap layer positions

```
LCCollectionVec* distanceToBarrelLayersVec = new LCCollectionVec(LCIO::LCFLOATVEC);
LCCollectionVec* distanceToEndcapLayersVec = new LCCollectionVec(LCIO::LCFLOATVEC);
// Fill the collections with their positions (in mm)
for(int l=0; l<=ecalLavers+hcalLavers+1; l++) {</pre>
 LCFloatVec* distanceToBarrelLayers = new LCFloatVec;
 LCFloatVec* distanceToEndcapLayers = new LCFloatVec;
 if(detectorType=="full") { // full detector
    if(1 \le 30) \{ // \text{ first 30 Ecal layers at a pitch of 3.9 mm (+ layer 0) \}
                                                                                              ← edit
      distanceToBarrelLayers->push back(1698.85+(3.9*1));
                                                                                                 edit
      distanceToEndcapLayers->push_back(2831.10+(3.9*1));
                                                                                                 edit
                                                                                                 edit
    else if(1>30 && 1<=ecalLayers) { // last 10 Ecal layers at a pitch of 6.7 mm
                                                                                                 edit
      distanceToBarrelLayers->push back(1815.85+(6.7*(1-30)));
                                                                                               ← edit
                                                                                              ← edit
      distanceToEndcapLayers->push back(2948.10+(6.7*(1-30)));
                                                                                              ← edit
    else { // 40 Hcal layers at a pitch of 24.5 mm (+ layer 81)
                                                                                              ← edit
      distanceToBarrelLayers->push back(1931.25+(24.5*(1-41)));
                                                                                              ← edit
      distanceToEndcapLayers->push back(3039.25+(24.5*(1-41)));
                                                                                              ← edit
                                                                                              ← edit
 else if(detectorType=="prototype") { ...some more code... } // prototype detector
 distanceToBarrelLayersVec->push back(distanceToBarrelLayers);
 distanceToEndcapLayersVec->push back(distanceToEndcapLayers);
// Save the collections
```

```
evt->addCollection(distanceToBarrelLayersVec, "distance_barrellayers");
evt->addCollection(distanceToEndcapLayersVec, "distance_endcaplayers");
```

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Getting started with MAGIC

- For new LCIO CalorimeterHits collection can:
 - getCellID0();
 - getCellID1() \Rightarrow pseudolayer/stave id encoded like layer/stave id in CellIDO;
 - getEnergy();
 - getPosition();
 - getType() \Rightarrow "0"=ecal hit; "1"=hcal hit.
- For all new LCIO Calorimeter*Clusters collections, can:
 - getCalorimeterHits();
 - getHitContributions(); and
 - getClusters()

(no energy/position/shape attributes set—user can set these in own private processors as desired).

- If simulation, can also use LCRelationNavigator to:
 - simHitRel->getRelatedToObjects(hit), and
 - mCParticleRel->getRelatedToObjects(trueCluster).