

analysis tools in e14lib developing version

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Csl data analysis

data from FADC (in KEKCC:/hsm/had/koto/)

↓ convert to rootfile

pulse shape information

Feb. data (in run40/conv_data/crate[0-10])

run????_conv.root

Jun. data (in run43/fiber_out/node[0-10])

run????_conv_crate[0-11]_node[0-9].root

conv
data

↓ convert (pulse shape info are lost)

ID, energy(integratedADC) , time

Feb. data (in run40/sumup_data/)

Sum????_root

Jun. data (in run43/fiber_merged)

Sum????_fiber.root

sumup
data

**MC simulation
of KL decays**

ID, energy, time

cluster finding

π^0 or K_L reconstruction etc..

Csl data analysis

data from FADC (in kekcc:/hsm/had/koto/)

↓ convert to rootfile

pulse shape information of each FADC crate

Feb. data (in run40/conv_data/crate[0-10])

run????_conv.root

Jun. data (in run43/fiber_out/node[0-10])

run????_conv_crate[0-11]_node[0-9].root

↓ convert (pulse shape info are lost)

ID, energy(integratedADC) , time

Feb. data (in run40/sumup_data/)

Sum????_root

Jun. data (in run43/fiber_merged)

Sum????_fiber.root

el4lib_dev

**MC simulation
of KL decays**

ID, energy, time

gsim4
(geant4 based MC)

cluster finding

π^0 or K_L reconstruction etc..

**today's
topic**

How to install e14lib dev. ver.

- **install to your own PC (for bash)**

0, Prepare geant4, ROOT and clhep.

1, In the directory where you want to install the library,

`svn co https://justice.hep.sci.osaka-u.ac.jp/svn/e14lib/trunk/e14`

(you'll be asked username and password. user: koto, passwd:[usual one])

2, add follows to your .bashrc.

`export E14_TOP_DIR=foo/bar/e14 (where you installed e14lib)`

`source $E14_TOP_DIR/scripts/e14.sh` `>& /dev/null`

`export PATH=$E14_TOP_DIR/bin:$PATH` Comment out here at first
time to see error message.
If any messages, follow it.

3, compile the library.

`source ~/.bashrc`

`cd $E14_TOP_DIR`

`make install-headers` (←Please ignore error messages in this stage.)

`make`

for KEKCC users

- **for KEKCC user (bash)**

The e14lib for common use is already prepared in

/sw/koto/e14lib/dev/installed20120803/e14/.

To use it, add follows to ~/Env/bashrc_post

geant4, root, clhep setting

source /sw/koto/e14lib/dev/installed20120803/requiredEnvironment

e14lib dev. version setting

source /sw/koto/e14lib/dev/installed20120803/setup.sh >& /dev/null

(In the file “**requiredEnvironment**”, proper environments of g4,root and clhep are set. Don't change those environments after this.)

* Also, if you want to prepare the private e14lib in KEKCC for some reasons, you can use the installation procedure in the previous page.

eI4 analysis tools

- We already have a lot of know-hows for various analysis which were accumulated through e391a and MC study .
⇒ some of them are available in eI4lib developing version.

source code: `$EI4_TOP_DIR/sources/ana/rec/`
provide ...

- cluster finding → class ClusterFinder
- gamma identification → class GammaFinder
- pi0 or KL reconstruction → class Rec2g , RecKlong
- energy correction → class EI4GNAnaFunction
- position correction → class EI4GNAnaFunction
- shower shape cut → class ShapeChi2, ShapeANN

examples

example program : \$E14_TOP_DIR/examples/el4gnana/

- el4clustering -- clustering from MC

bin/el4clustering [input file(MC data)] [output file]

- el4g2ana -- pi0 reconstruction for KLpi0nunu event

bin/el4g2ana [input file(clustering result)] [output file]

- el4g4ana -- KL reconstruction for KLpi0pi0 event

bin/el4g4ana [input file(clustering result)] [output file]

- el4g6ana -- KL reconstruction for KL3pi0 event

bin/el4g4ana [input file(clustering result)] [output file]

→ **see backup slides for how to use these programs**

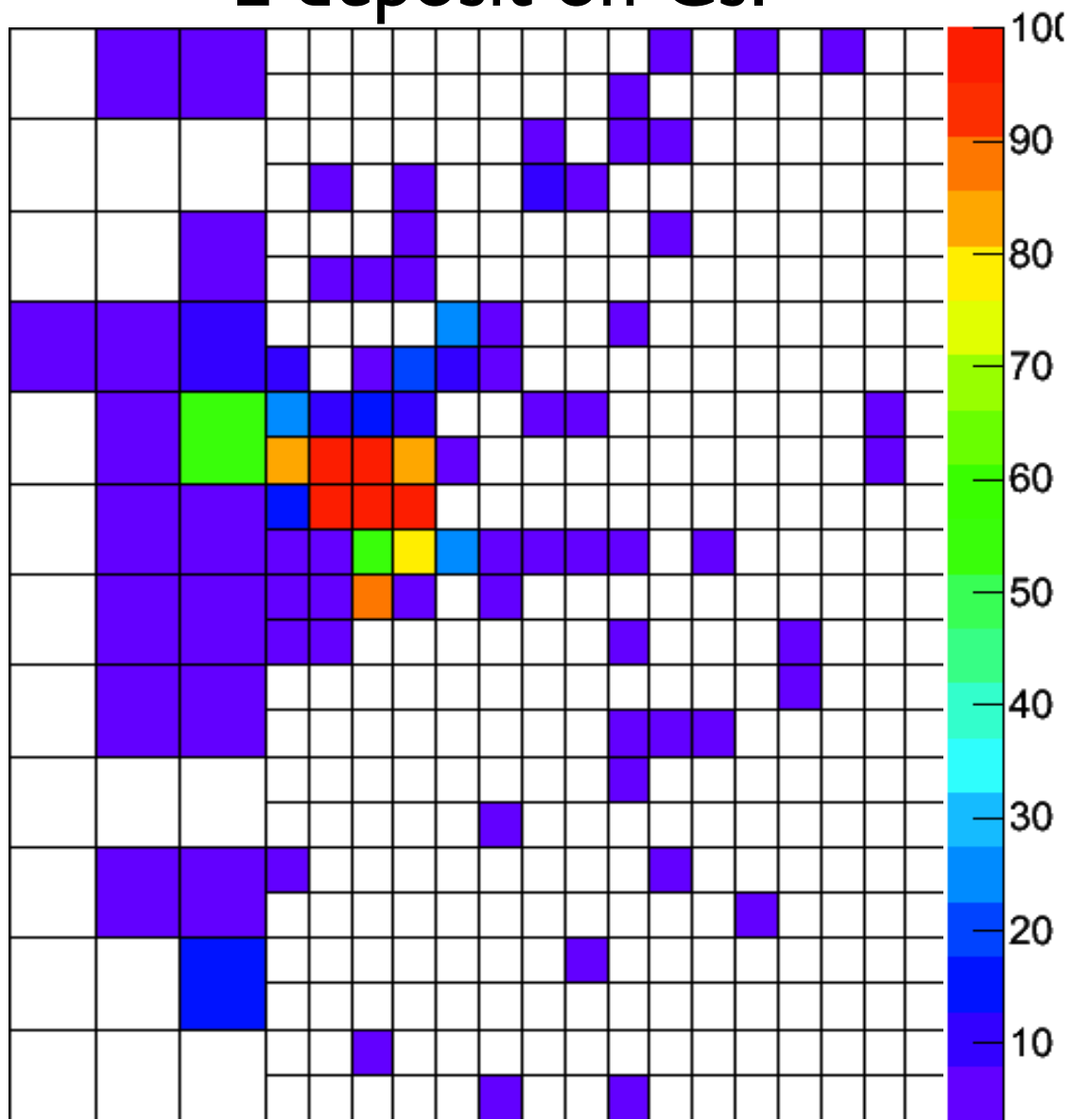
flow of analysis

I.cluster finding

- To measure gamma energy using Csl calorimeter, we should make “cluster”.
 - cluster = a group of Csl crystals with energy deposit exceeding a threshold.
 - class “ClusterFinder” provides the function for clustering
- ```
std::list<Cluster> findCluster(int nCsi,int *ID,double *E,double *time)
```

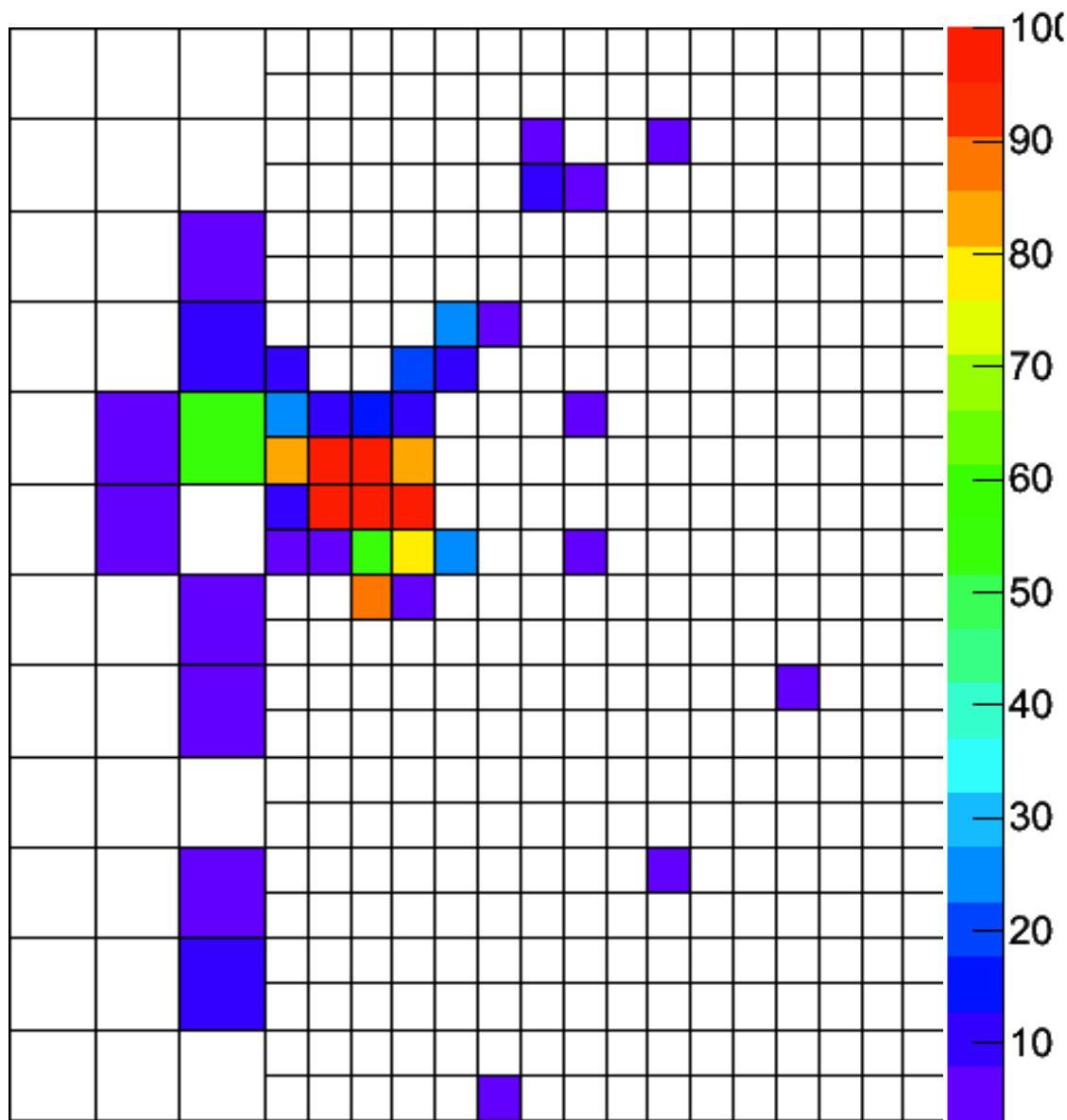
# finding clusters

E deposit on Csl



consider the case like  
left figure.

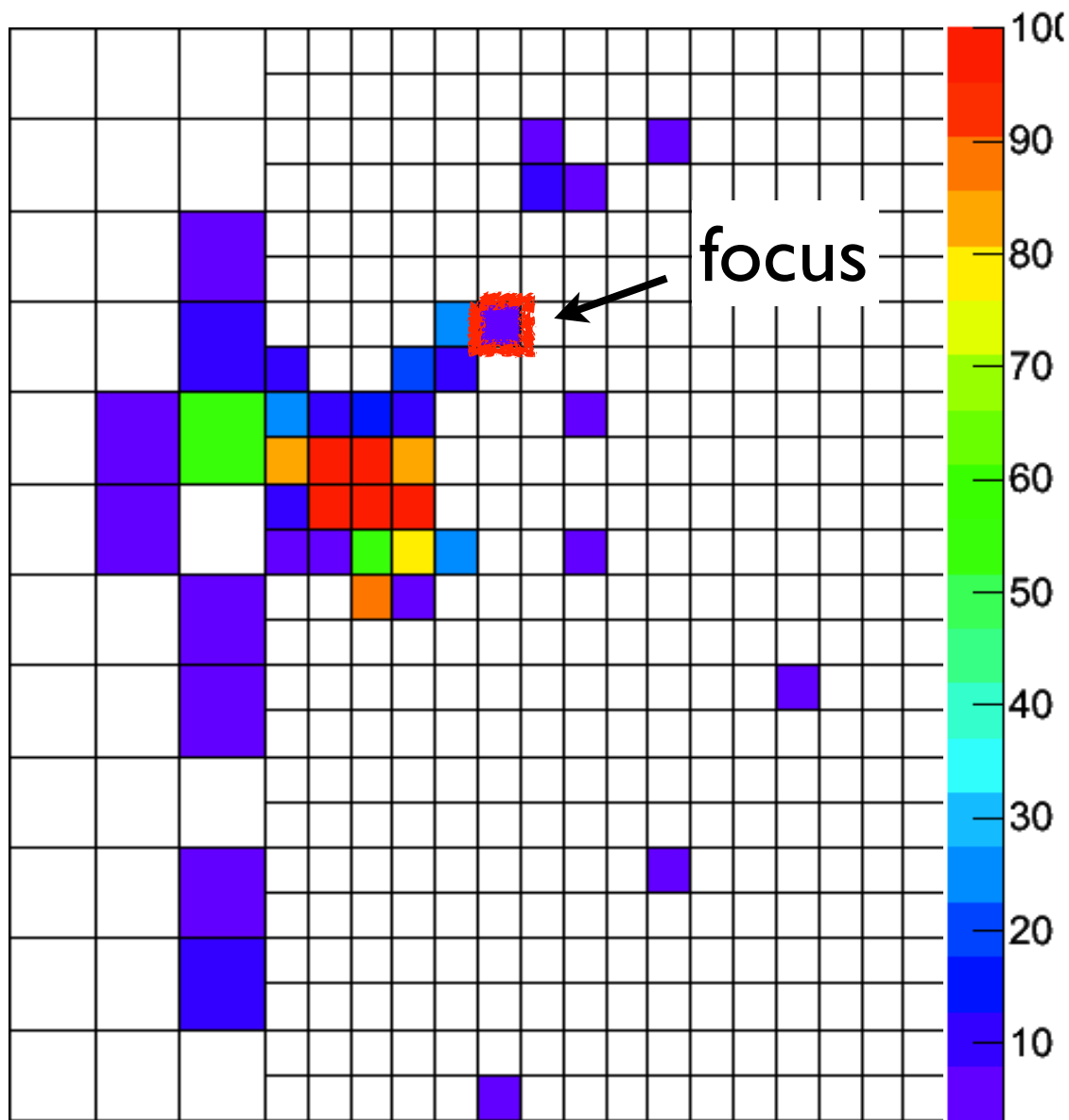
# finding clusters



! ,Set threshold  
(default = 3MeV)

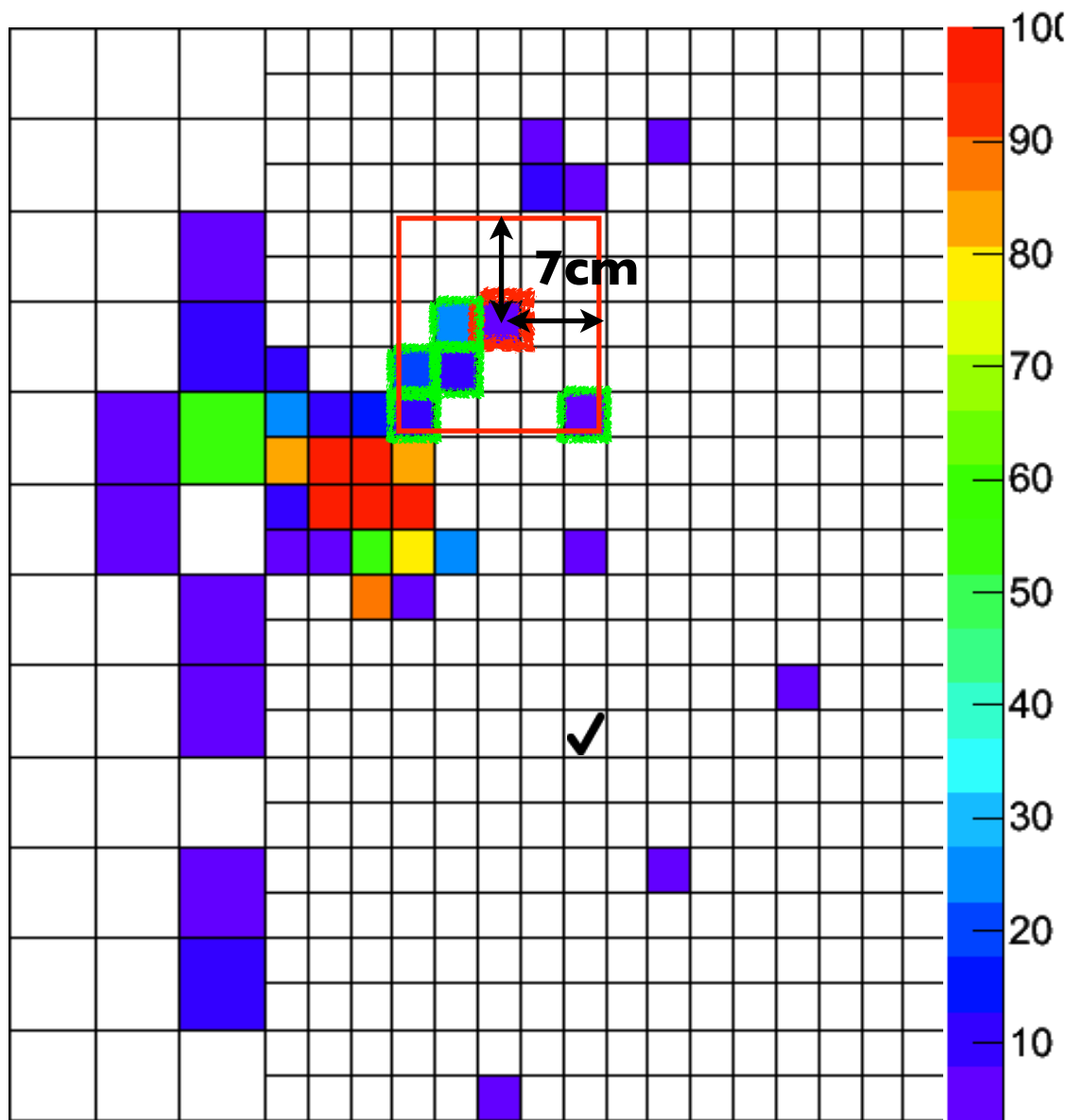
Ignore crystals whose  
energy deposit is less than  
threshold.

# finding clusters



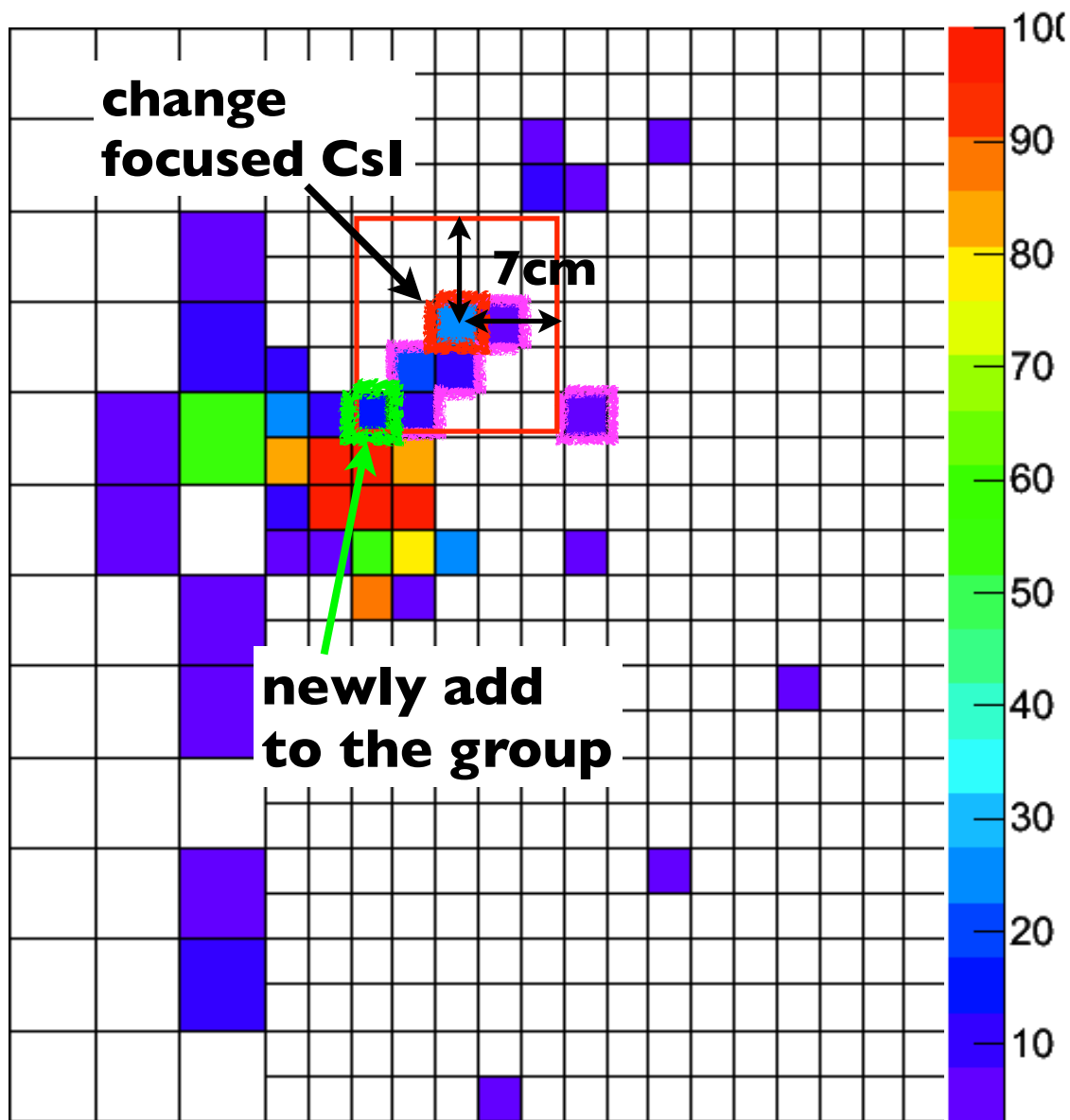
2, focus on one **crystal**.

# finding clusters



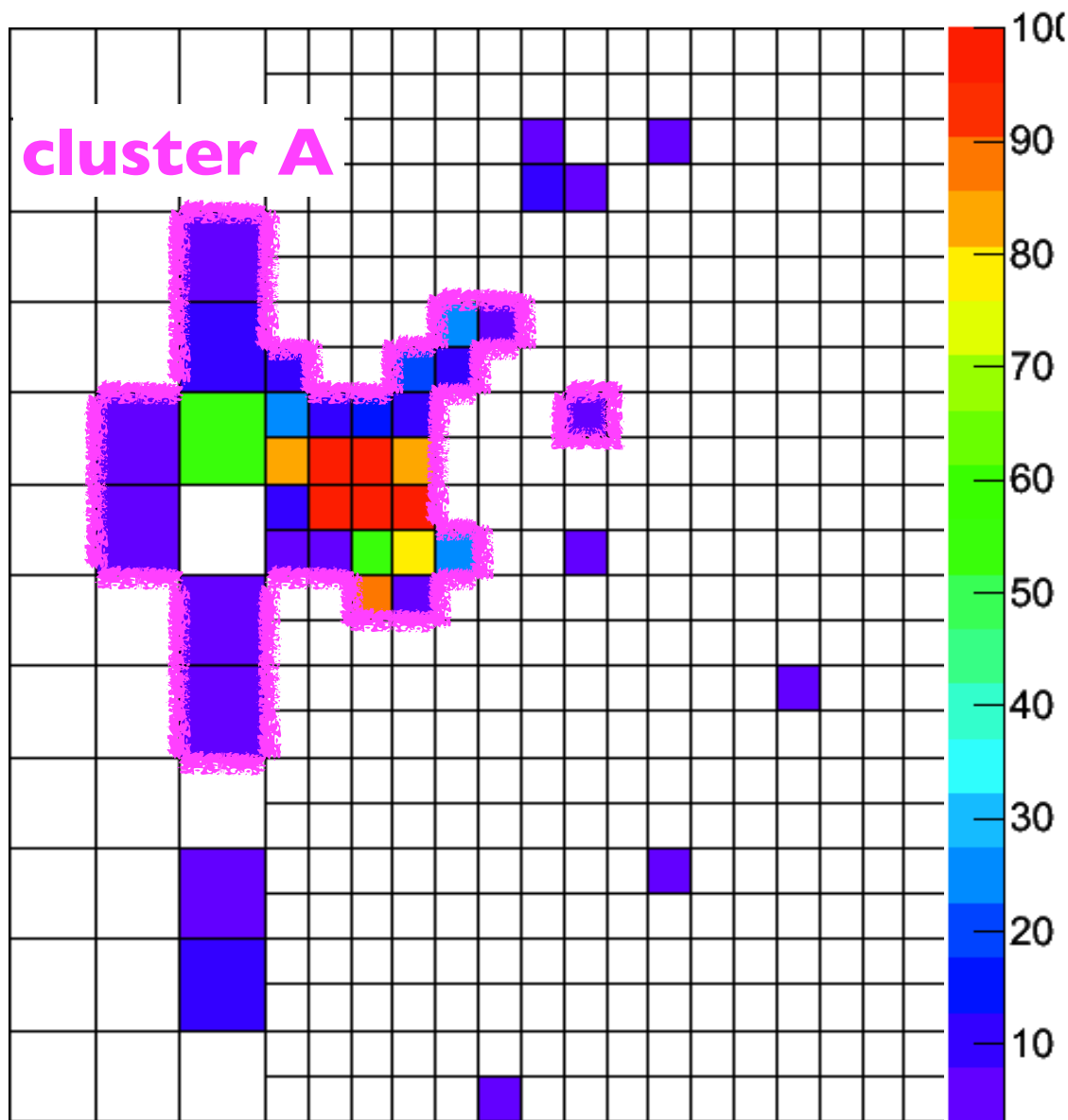
3, group **crystals** of which center positions are less than 7cm apart from the center of **focused crystal**

# finding clusters



4, focus on another **crystal** in the same group and repeat process 3.

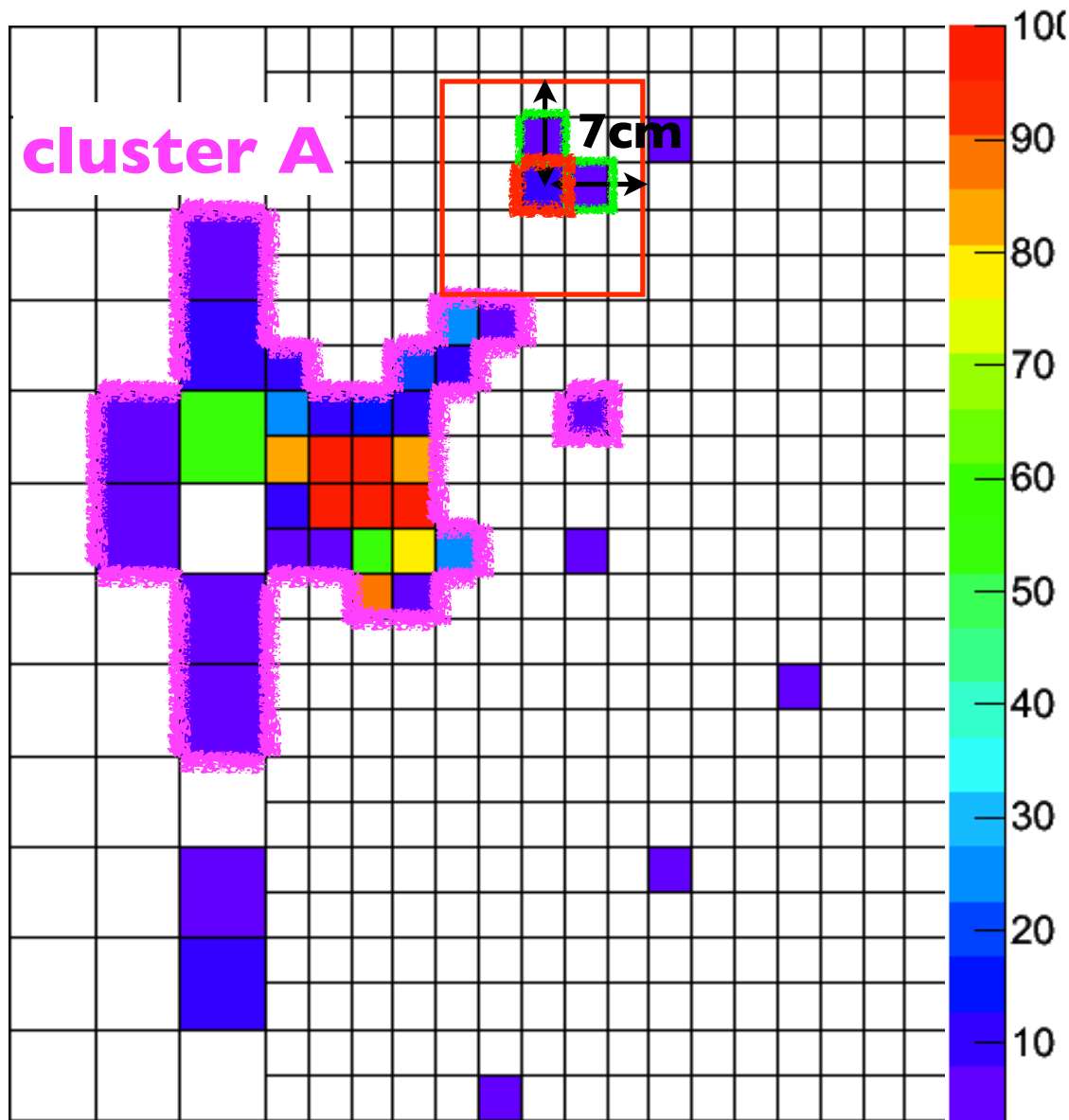
# finding clusters



4, focus on another **crystal** in the same group and repeat process 3.

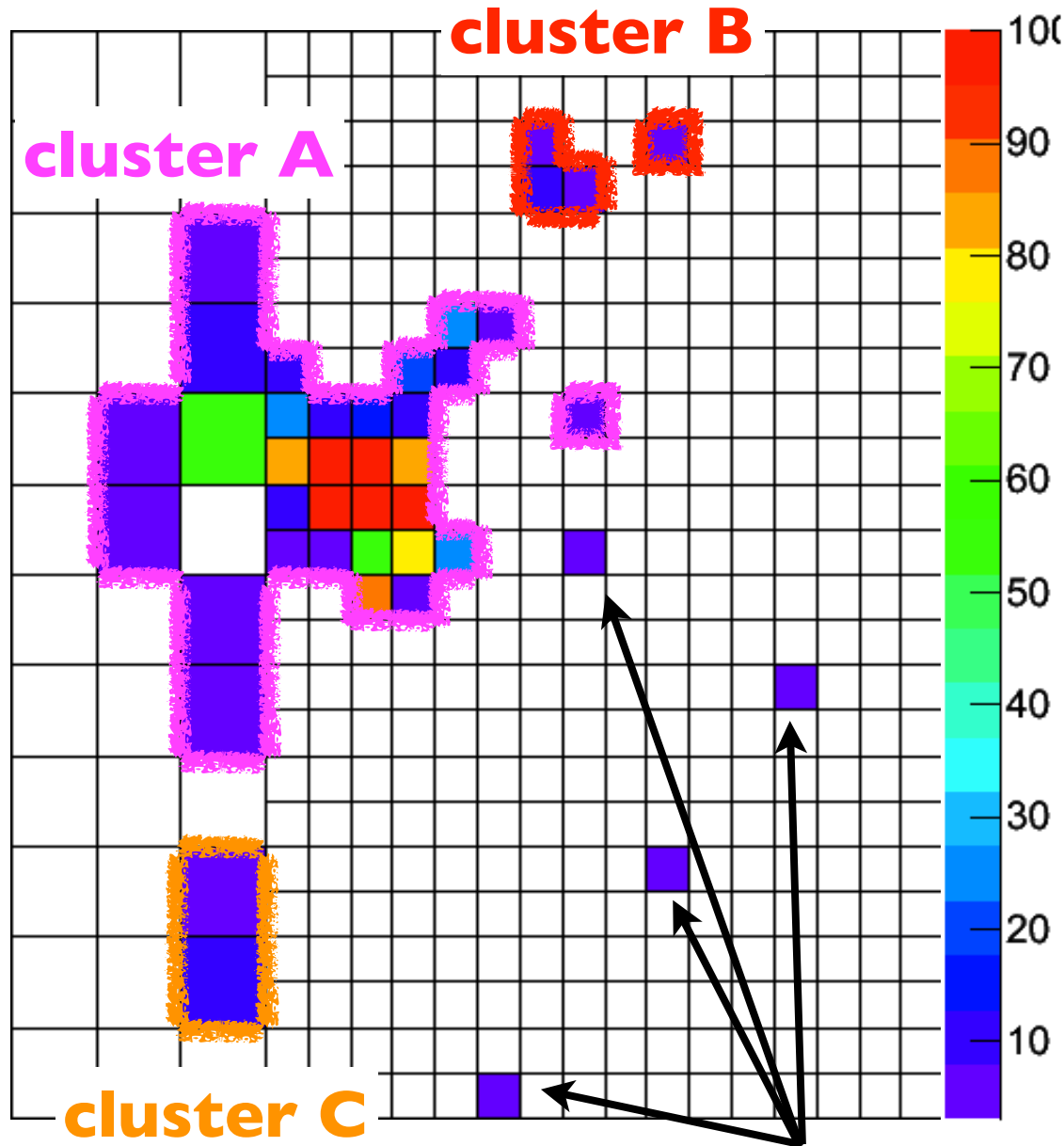
→ run over until there is no crystal to add to the group.

# finding clusters



5, focus on a remaining crystal and repeat the procedure above until all crystals are used.

# finding clusters



finally, in this case,  
3 clusters were found.  
( notice that single hits are  
not “cluster” )

single isolated crystals

# energy, position, time of cluster

following cluster informations are calculated

- sum of energy deposit
- position of center-of-energy

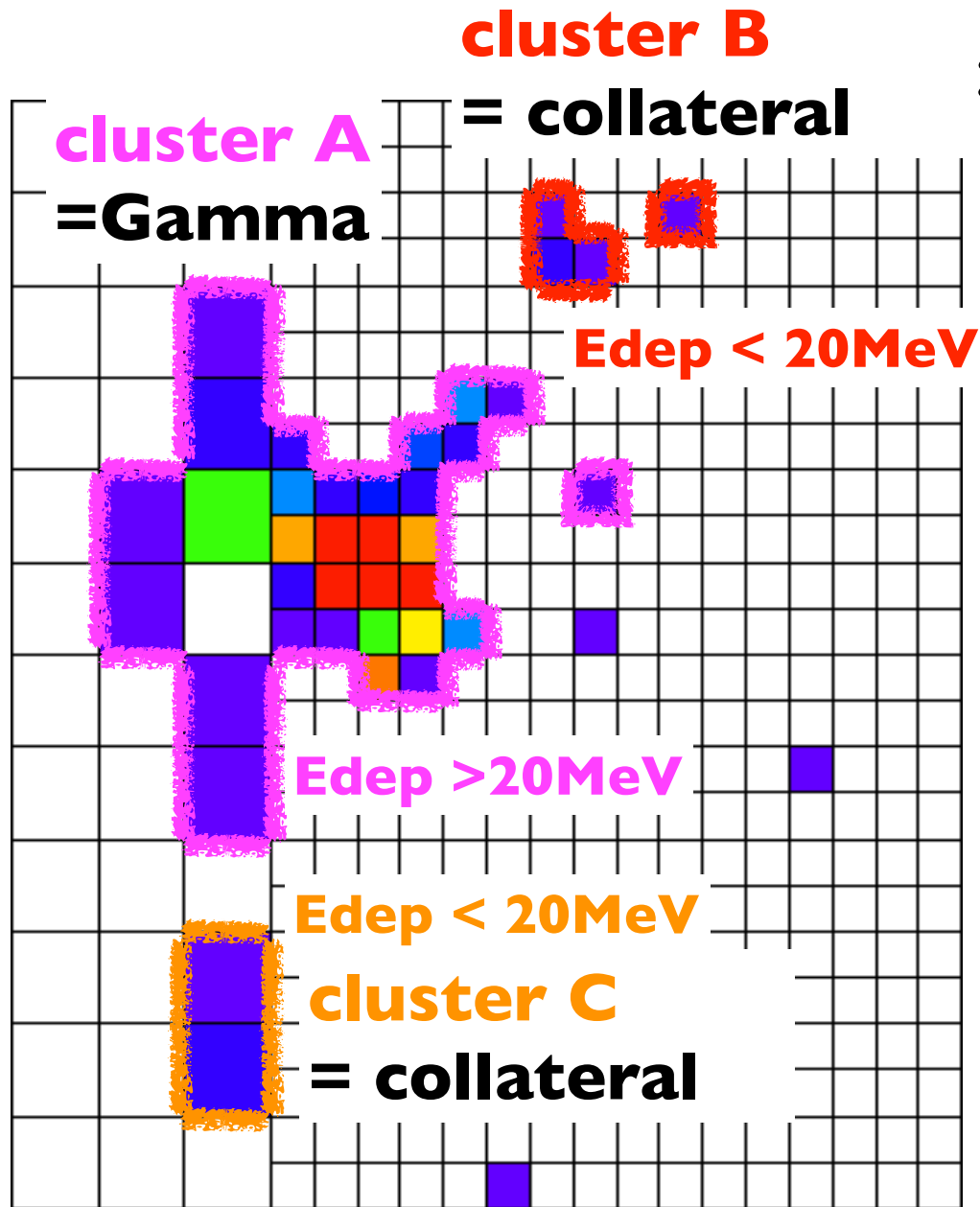
$$x_c = \sum_i^n \frac{(E_i \times x_i)}{\sum_i E_i} \quad \begin{array}{l} x_i = \text{position of each Csl} \\ E_i = \text{energy deposit each Csl} \end{array}$$

$$y_c = \sum_i^n \frac{(E_i \times y_i)}{\sum_i E_i}$$

- timing  Cluster timing is not studied sufficiently. Now I temporarily defined cluster time as earliest time of Csl hits.

# flow of analysis

## \*2. gamma finding



- gamma finding = to identify clusters as gamma cluster or not

- require energy deposit > 20 MeV as a gamma cluster

note: “gamma cluster” means “not a collateral cluster but a cluster made by particle”.

I don't mean that “gamma cluster” is made by gamma.

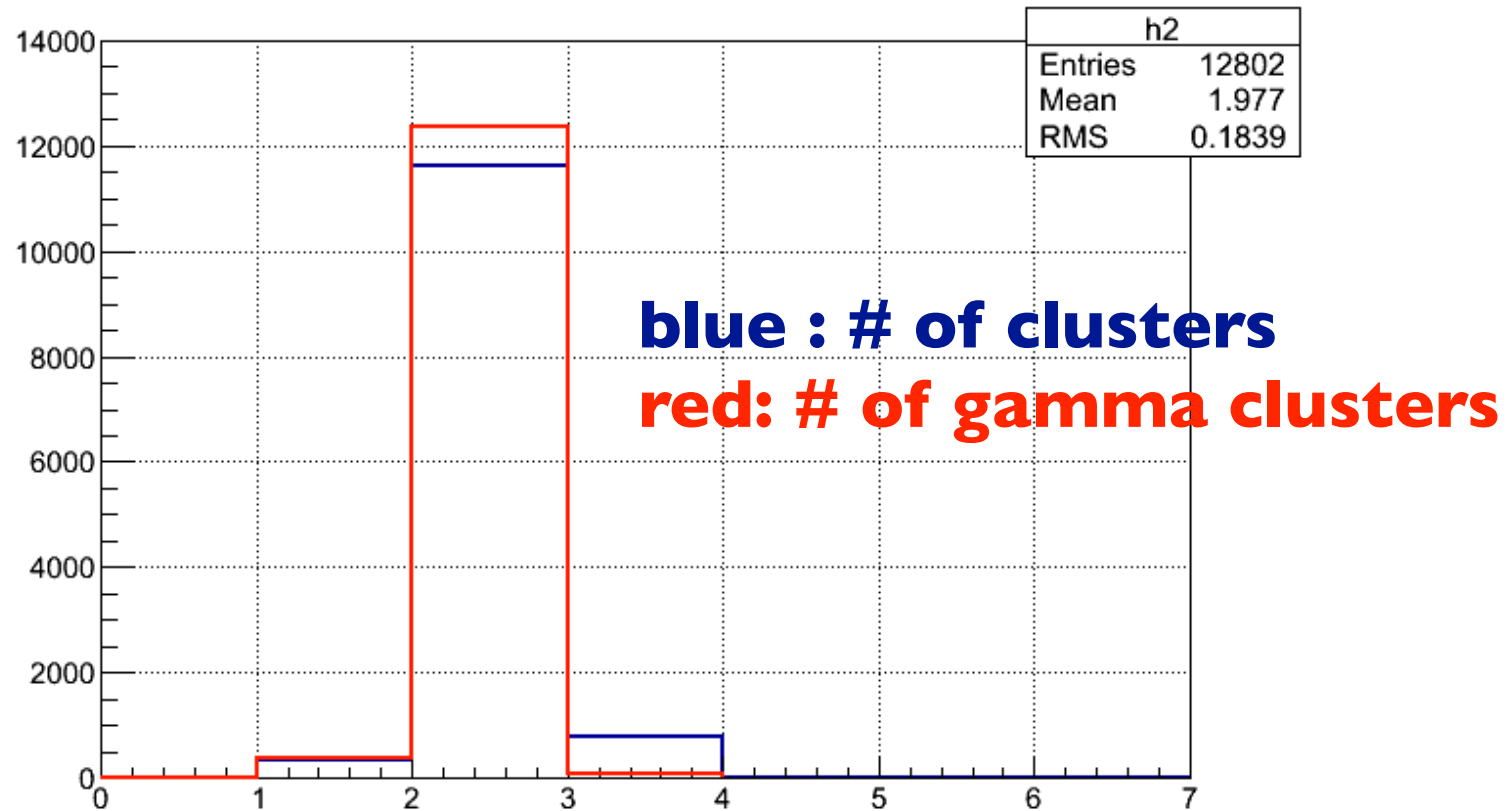
class “GammaFinder”

provide the Function for this.

```
void findGamma(std::list<Cluster> const &clist, std::list<Gamma> glist)
```

# KLpi0nunu MC

(events where 2 gammas hit on Csl fiducial)



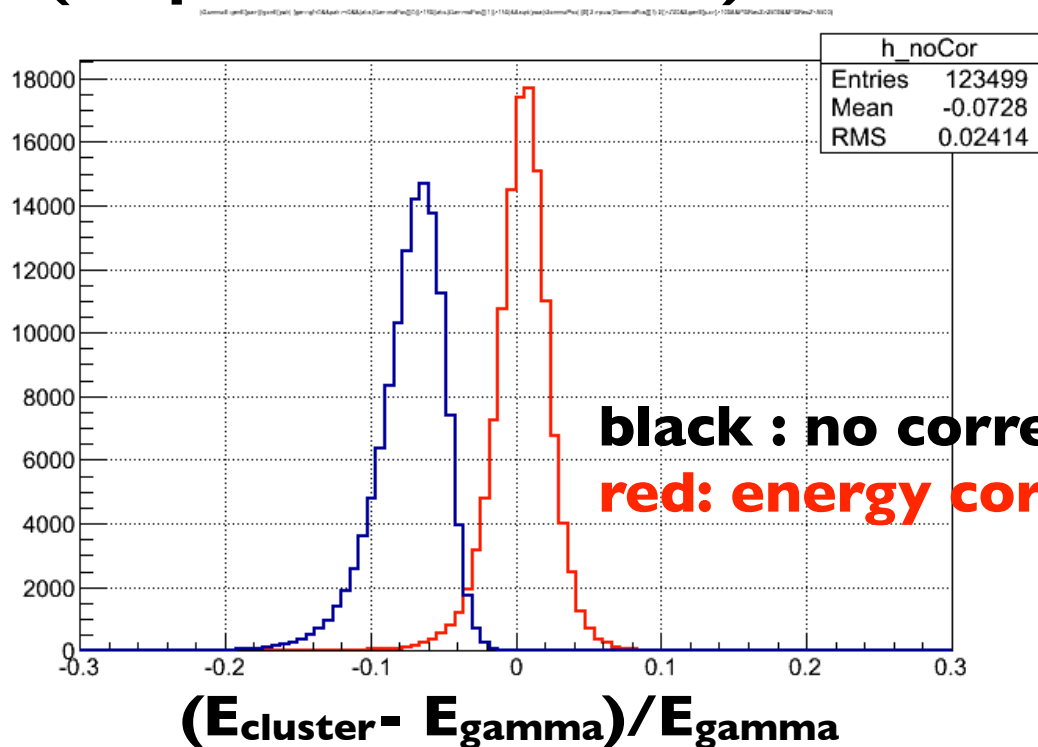
☆ 90.9% accepted → 96.6%

Now I just require very simple condition ( $E \text{ deposit} > 20 \text{ MeV}$ ).  
One who analyzes the data should study and develop this procedure.

# flow of analysis

## 3. energy correction (1st)

**energy difference of  
gamma( $E_{\text{gamma}}$ ) & cluster( $E_{\text{cluster}}$ )  
( KLpi0nunu event MC )**



- correct cluster energy  
( some energies of incident  
gamma leaked out of cluster )

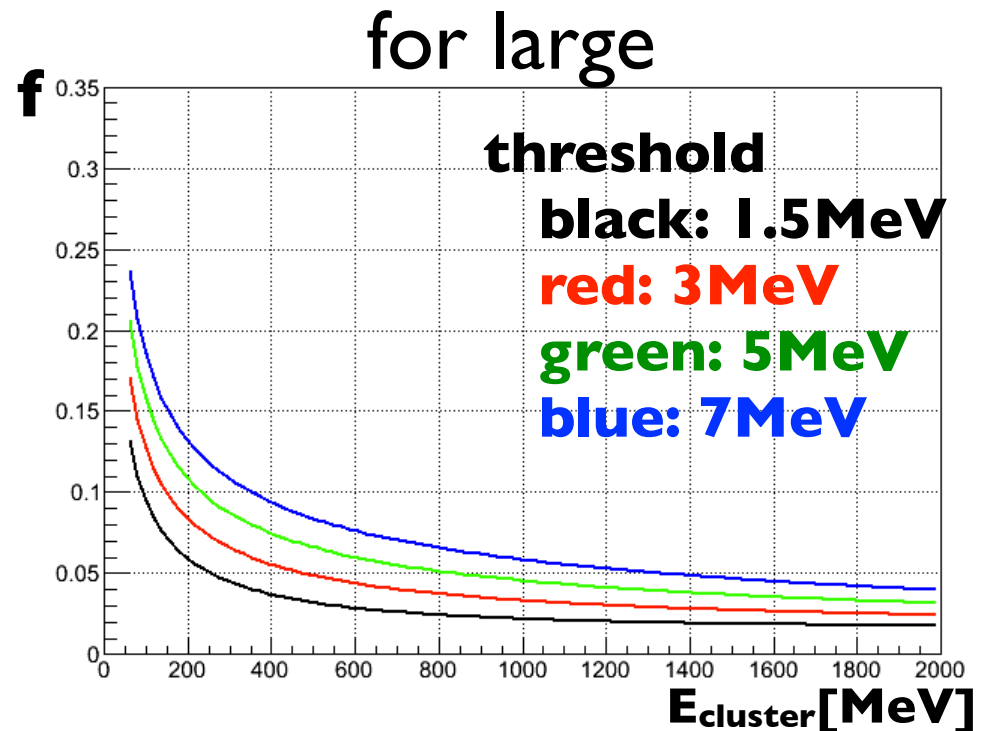
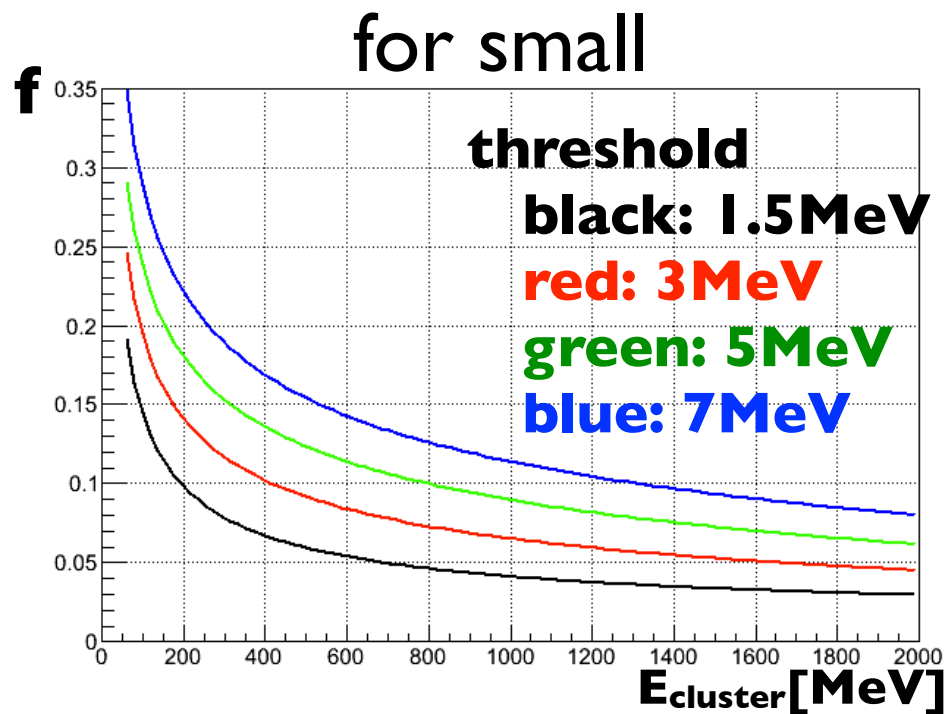
$$E_{\text{gamma}} = (1 + f(E_{\text{cluster}})) \times E_{\text{cluster}}$$

this process is also done in `GammaFinder::findGamma()`

```
void findGamma(std::list<Cluster> const &clist, std::list<Gamma> glist)
```

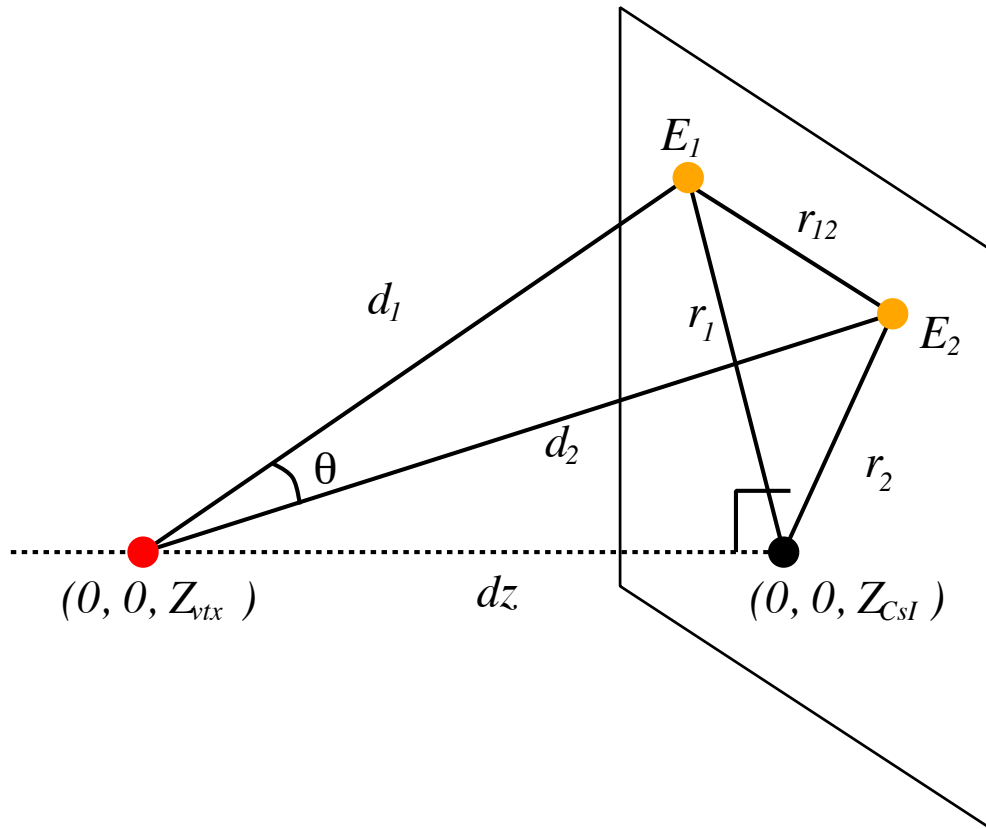
# correction function

- correction function **f** was derived from KLpi0nunu MC.
- fraction of energy leakage strongly depends on energy and threshold.  
→  **$f = f(E_{\text{cluster}}, \text{threshold})$**
- **f** are prepared for small and large crystals respectively.  
Energy weighted mean of 2 functions are used for clusters which include both small and large crystals
- **$f = 0.014 + 0.052/\sqrt{E[\text{GeV}]} - 0.0074 \log(E[\text{GeV}])$**  [Small, thre=3MeV]



# flow of analysis

## 4-1. reconstruction pi0



- for the events where 2 gamma clusters on CsI calorimeter.

- assume that 2 gamma on CsI were generated from pi0 then,

$$\cos\theta = 1 - \frac{M_{\pi^0}^2}{2E_1E_2} .$$

- assume the vertex is on beam axis

→ can calculate  $Z_{vtx}$  as left figure

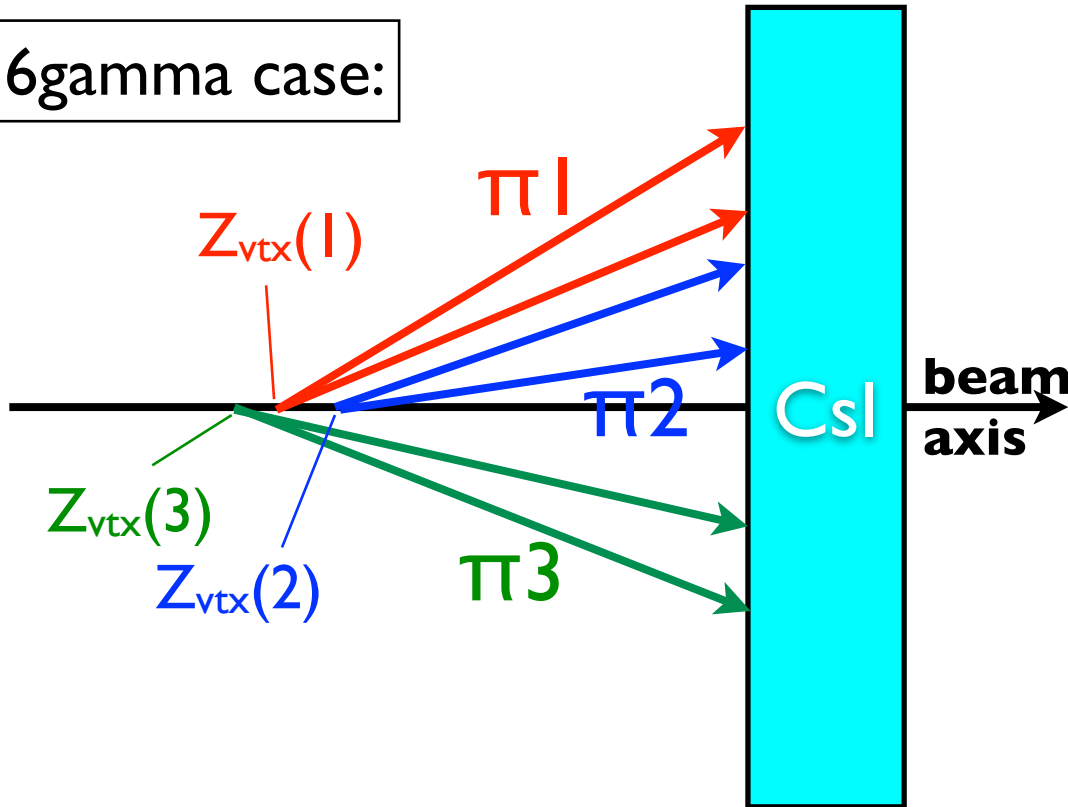
class Rec2g provide the function for pi0 reconstruction

```
std::list<Pi0> recPi0withConstM(std::list<Gamma> glist, const double M=Mpi0)
```

# flow of analysis

## 4-2. reconstruction KL

6gamma case:



- for the events where 4 or 6 gamma clusters on CsI.
- pions are reconstructed using a method in previous page.  
→  $Z_{KL} = \text{mean of } Z_{\pi}$

$$Z_{KL} = \frac{\sum_{i=1}^3 [Z_{\text{vtx}}(i) / \sigma_{\text{vtx}}(i)^2]}{\sum_{i=1}^3 [1 / \sigma_{\text{vtx}}(i)^2]}$$

- KL energy = sum of gamma energies

$$E_{K_L} = \sum_{i=1}^6 E_{\gamma}(i)$$

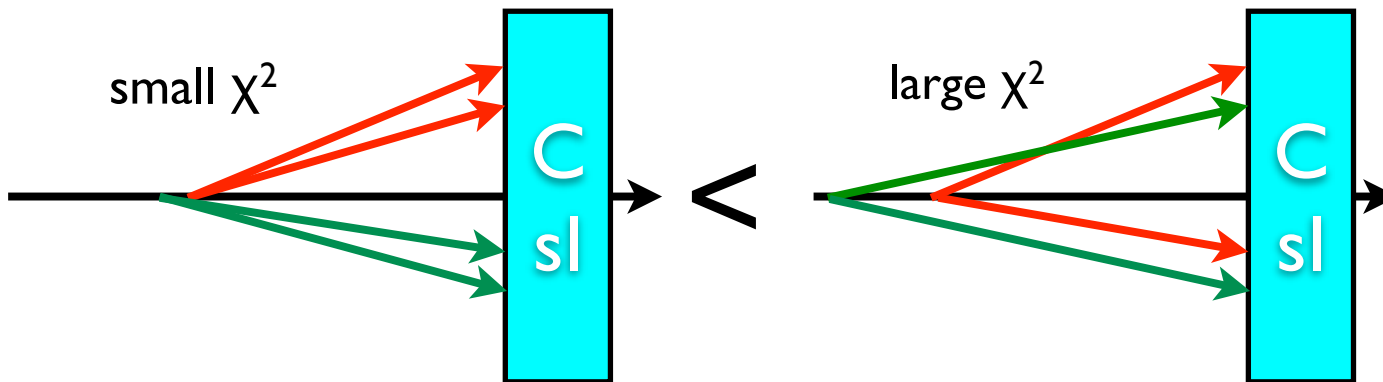
- combination of gammas is decided by chi square

$$\chi^2 = \sum_{i=1}^3 \frac{[Z_{K_L} - Z_{\text{vtx}}(i)]^2}{\sigma_{\text{vtx}}(i)}$$

- KL vertex x,y = center-of-gravity of gammas

$$\begin{pmatrix} X_{K_L} \\ Y_{K_L} \\ Z_{K_L} \end{pmatrix} = \begin{pmatrix} s \times \sum_{i=1}^6 E_{\gamma} X_{\gamma}(i) / E_{K_L} \\ s \times \sum_{i=1}^6 E_{\gamma} Y_{\gamma}(i) / E_{K_L} \\ Z_{K_L} \end{pmatrix}$$

$$s = \frac{Z_{K_L} - Z_{\text{target}}}{Z_{CsI} - Z_{\text{target}}},$$



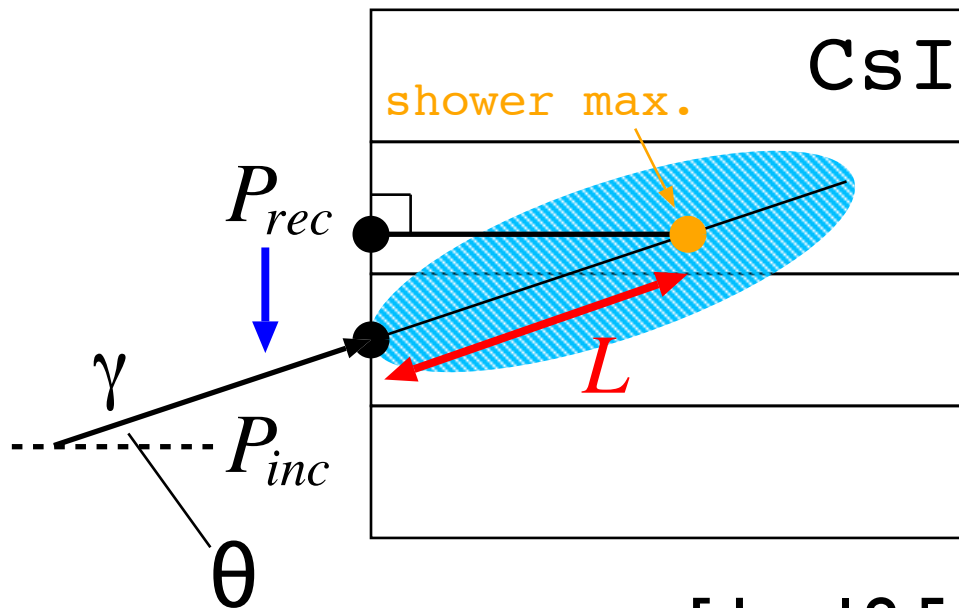
class RecKlong provide the function for KL reconstruction

`std::vector<Klong> recK2pi0( const std::list<Pi0>& pi0list)`

`std::vector<Klong> recK3pi0( const std::list<Pi0>& pi0list)`

# flow of analysis

## 5. position correction



- correct cluster position
- center-of-energy position ( $P_{rec}$ ) is different from gamma incident position ( $P_{inc}$ ) due to shower length

$$P_{inc} = P_{rec} (1 - L \sin \theta)$$

[  $L = 18.5(6.49 + 0.992 \log(E[\text{GeV}]))$  mm from MC ]

the function is defined in class E14GNAnaFunction

**void correctPosition(Gamma& gamma)**

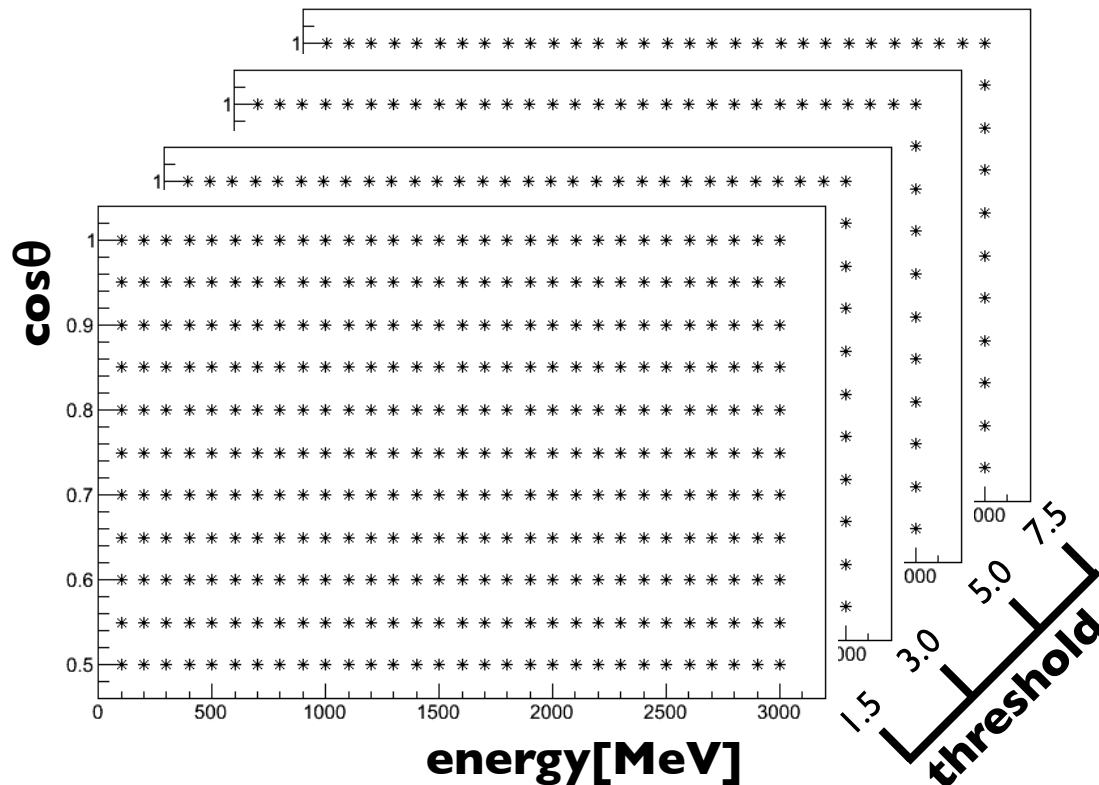
Note: this can't be done before vertex reconstruction because angle of gamma is needed.

# flow of analysis

## 6. energy correction (2nd)

- energy correction including angle dependence.

points on correction table



**$f = f(E, \cos\theta, \text{threshold})$**

is derived from correction table

➡ each point have value  $f_i$   
estimated by MC

To calculate  $f$  from these point,

- spline interpolation for  $E$
- linear interpolation  
for  $\cos\theta$  and threshold

the function is defined in class E14GNAnaFunction

**void correctEnergyWithAngle(Gamma& gamma)**

# flow of analysis

## 7. re-reconstruction

- repeat  $\pi^0(K_L)$  reconstruction again with corrected positions and energies.

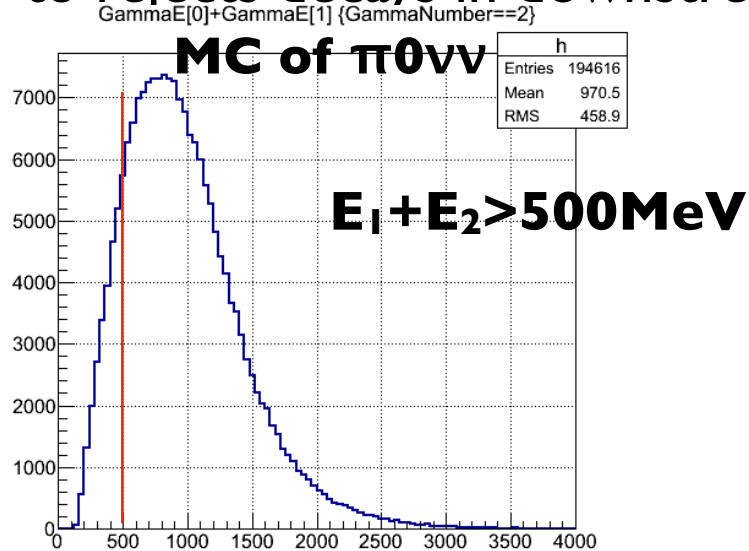
you can find some examples in

[\\$E14\\_TOP\\_DIR/examples/e14gnana/e14g?ana/src/user\\_rec.cc](#)

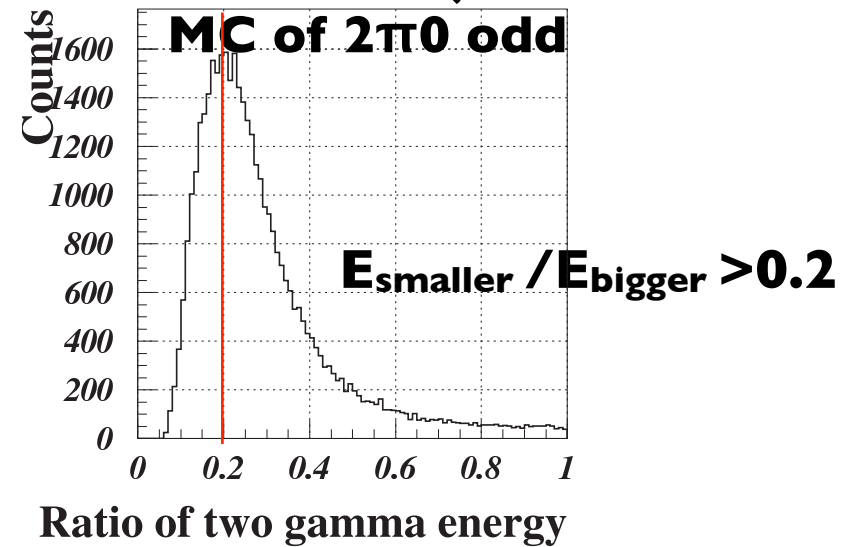


- sum of energy

to rejects decays in downstream

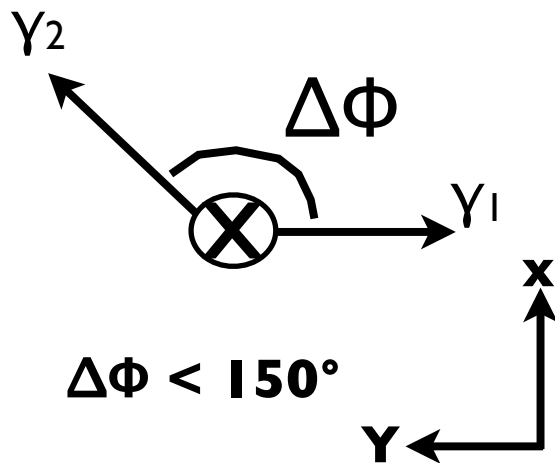


- E ratio to reject  $2\pi^0$  odd

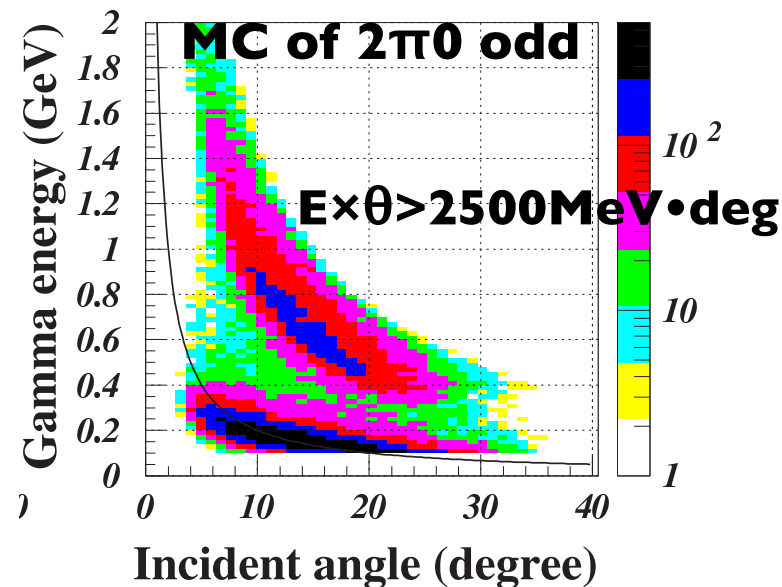


- collinear

to reject  $K \rightarrow 2\gamma$



- E- $\theta$  to reject  $2\pi^0$  odd

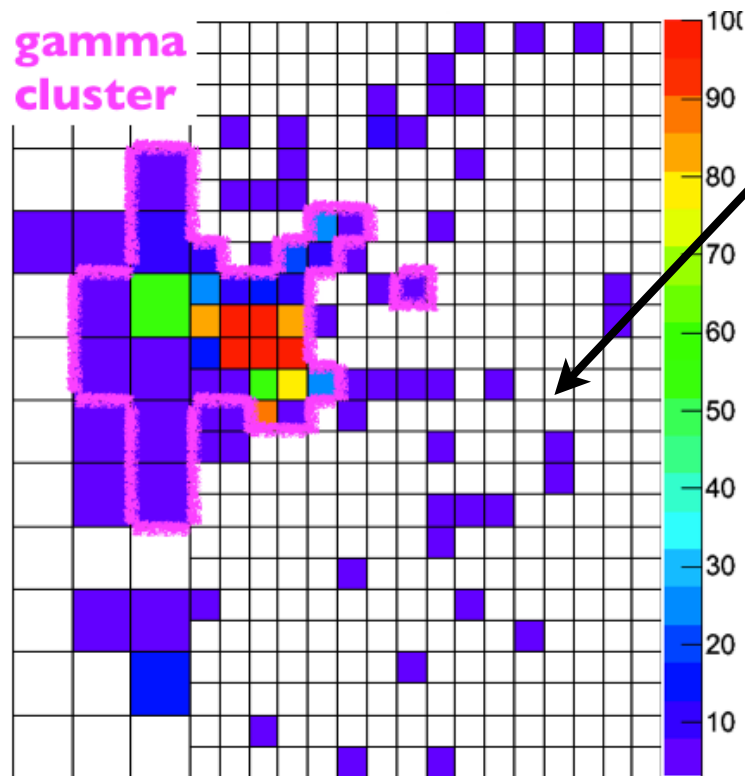


you can find a example of how to apply these cuts in

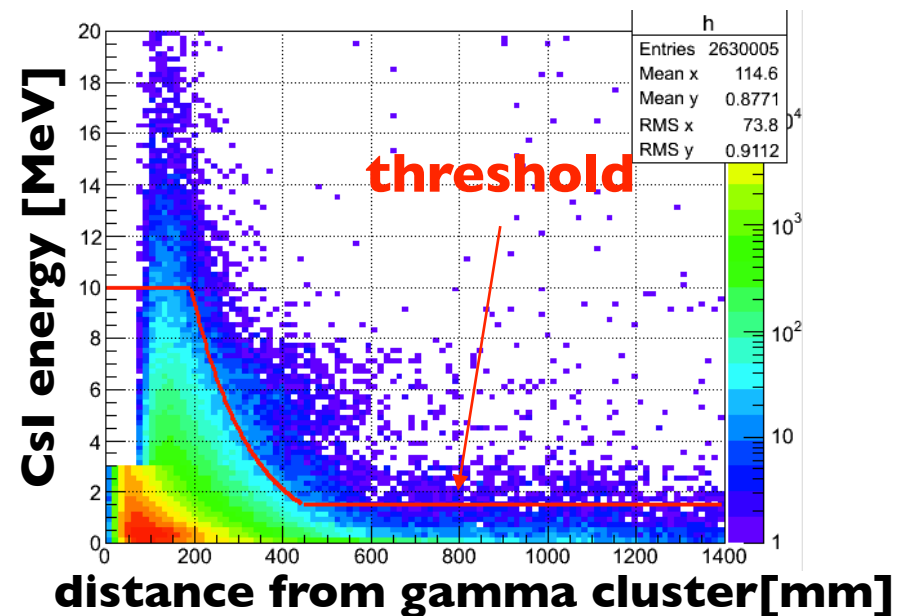
[\\$E14\\_TOP\\_DIR/examples/e14gnana/e14g2ana/src/user\\_cut.cc](#)

# flow of analysis

## \*9. Csl veto cut



Csl hits that were not included in gamma clusters should have energy less than a threshold.



threshold is a function of distance from gamma cluster  
(derived from MC, signal acceptance loss = 10%)

you can find some example of how to apply this cut in

[\\$E14\\_TOP\\_DIR/examples/e14gnana/e14g2ana/src/user\\_cut.cc](#)

# flow of analysis

## \*10. veto cut

- \* not prepared standard cut values for veto detector.  
It's responsible for the one who analyze data.

you can find some example in

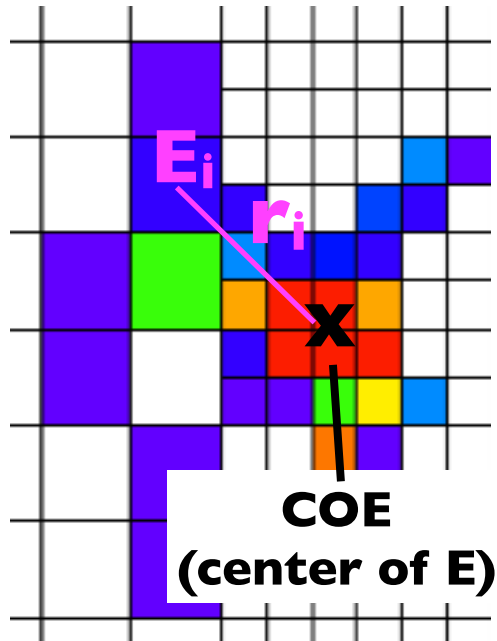
[\\$E14\\_TOP\\_DIR/examples/el4gnana/el4g2ana/src/user\\_cut.cc](#)

# flow of analysis

## \* I I. shower shape cut

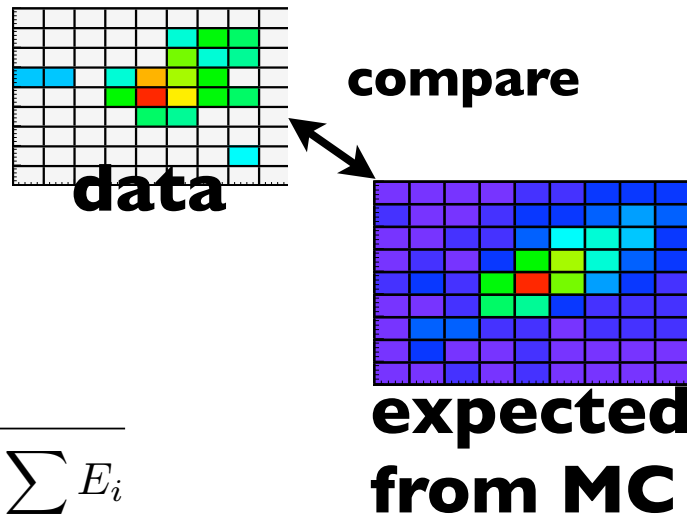
- qualify gamma cluster from its shape.

- RMS

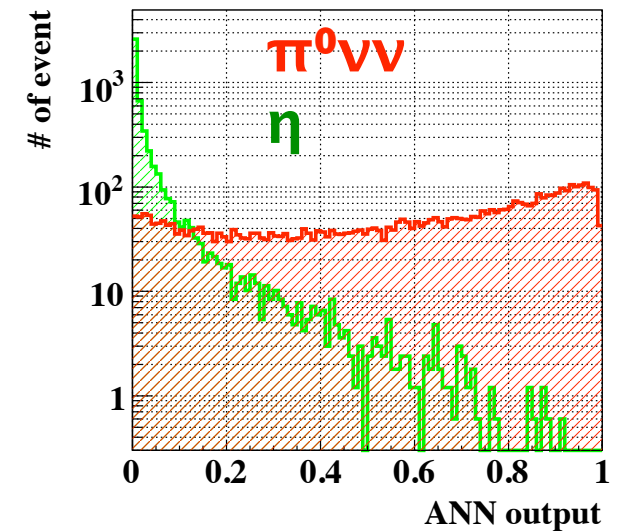


$$RMS = \sqrt{\sum_i (E_i \times (r_i - COE)^2) / \sum_i E_i}$$

- $\text{shape}\chi^2$   
compare with MC  
expectation



- ANN  
neural net to  
distinguish  $\eta$  events



RMS : automatically calculated in cluster finding

$\text{shape}\chi^2$  : class ShapeChi2 provides. `void shapeChi2(Gamma& gamma)`

$\text{shapeANN}$  : class ShapeANN provides. `void shapeANN(Gamma& gamma)`

# data container class

☆ classes of data container for analysis

- Cluster (**libcluster**) :  $\$E14\_TOP\_DIR/sources/ana/rec/cluster$ 
  - cluster informations
- Gamma (**libgamma**) :  $\$E14\_TOP\_DIR/sources/ana/rec/gamma$ 
  - gamma-cluster informations
- Pi0 (**libpi0**) :  $\$E14\_TOP\_DIR/sources/ana/rec/pi0$ 
  - reconstructed  $\pi^0$  informations
- Klong (**libklong**) :  $\$E14\_TOP\_DIR/sources/ana/klong$ 
  - reconstructed  $K_L$  informations

# Klong

**Pi0**

**Gamma**

position  
momentum  
etc...

**Cluster**

COE position  
E deposit  
time  
Csl ID  
Csl E  
Csl time  
threshold

vertex  
momentum  
inv. mass  
etc...

**Gamma  
Cluster**

**Pi0**

**Gamma  
Cluster**

**Gamma  
Cluster**

**Pi0**

**Gamma  
Cluster**

**Gamma  
Cluster**

vertex  
momentum  
inv. mass  
etc...

# summary

- E14 library developing version has various functions for ...
  - clustering
  - gamma identification
  - $\pi^0$  or KL reconstruction
  - energy correction
  - position correction
  - shower shape cut
  - ...etc

I summarized the functions of e14lib along standard analysis flow.

- Some example executables are prepared in `$E14_TOP_DIR/examples/e14gnana`. They will help you to learn how to use e14lib for analysis.

# e l 4lib web page

- e l 4lib developing version is often updated.  
You can get the information of e l 4lib from the web page  
<https://justice.hep.sci.osaka-u.ac.jp/trac/e l 4lib>  
( username: koto, password:[usual one] ).
  - from “timeline” tab, you can see the changes in recent updates.
- \* some information in this web page is a bit old. Please ignore  
“installation” part.

backup

flow chart  
of example programs

# eI4clustering

output from gsim4

(GsimDigiData)

data conversion to simple array form

DigiReader::getCsiDigi(n,ID[],E[],Time[])

(array ( n,ID[],E[],Time[] ) )

clustering

ClusterFinder::findCluster(n,ID[],E[],Time[])

(std::list<Cluster>)

data conversion to simple array form

EI4GNAnaDataContainer::setData(std::list<Cluster>)

(array)

output to root file

# eI4g2ana

output from eI4clusteing

↓ (array)

data conversion to `std::list<Cluster>`

`EI4GNAnaDataContainer::setData(std::list<Cluster>&)`

↓ `std::list<Cluster>`

gamma finding & 1st energy correction

`GammaFinder::findGamma(std::list<Cluster>, std::list<Gamma>&)`

↓ `std::list<Gamma>`

pi0 reconstruction

`Rec2g::recPi0WithConstM(std::list<Gamma>)`

↓ `std::list<Pi0>`

position correction & 2nd energy correction → reconstruction

`EI4GNAnafunction::correctPosition(Pi0&)`

`EI4GNAnafunction::correctEnergyWithAngle(Pi0&)`

`Rec2g::recPi0WithConstM(std::list<Gamma>)`

`std::list<Pi0>` ← apply kinetic cuts, veto cuts, shape cut

data conversion to simple array form

`EI4GNAnaDataContainer::setData(std::list<Pi0>)`



output to  
root file

# eI4g4(6)ana

output from eI4clusteing

↓ (array)

data conversion to `std::list<Cluster>`

`EI4GNAnaDataContainer::setData(std::list<Cluster>&)`

↓ `std::list<Cluster>`

gamma finding & 1st energy correction

`GammaFinder::findGamma(std::list<Cluster>, std::list<Gamma>&)`

↓ `std::list<Gamma>`

KL reconstruction

`RecKlong::recK2pi0(std::list<Gamma>)`

↓ `std::vector<Klong>`

position correction & 2nd energy correction → reconstruction

`EI4GNAnafunction::correctPosition(Klong&)`

`EI4GNAnafunction::correctEnergyWithAngle(Klong&)`

`RecKlong::recK2pi0(std::list<Gamma>)`

`std::vector<Klong>` ↓ ← apply kinetic cuts, veto cuts, shape cut

data conversion to simple array form

`EI4GNAnaDataContainer::setData(std::vector<Klong>)`

→ output to root file

demonstration (I)  
estimate # of  $K_L \rightarrow \pi^0 \nu \nu$   
events using MC

**(how to use analysis tools)**

# generate simulation

```
Terminal — ssh — 80x24
[sato@xenys example]$
[sato@xenys example]$ mkdir KLpi0nunuAna make working directory
[sato@xenys example]$ cd KLpi0nunuAna/
[sato@xenys KLpi0nunuAna]$ mkdir gsimoutput
[sato@xenys KLpi0nunuAna]$ cp $E14_TOP_DIR/examples/gsim4test/e14.mac .
[sato@xenys KLpi0nunuAna]$ emacs e14.mac & copy a macro file and edit it
[1] 24782 (→see next page)
[sato@xenys KLpi0nunuAna]$ $E14_TOP_DIR/bin/gsim4test e14.mac gsimoutput/KLpi0nu
nu.root 100000 0 0 output file
#event random seed

Geant4 version Name: geant4-09-00 (29-June-2007)
Copyright : Geant4 Collaboration
Reference : NIM A 506 (2003), 250-303
WWW : http://cern.ch/geant4

illegal code for meson PDG code=10000130
Particle KLpi0nunu has a strange PDGEncoding
illegal code for meson PDG code=100130
Particle KLpienu has a strange PDGEncoding
illegal code for meson PDG code=200130
Particle KLpimunu has a strange PDGEncoding
illegal code for meson PDG code=300130
```



**start MC**

# edit the macro file

\* configuration of MC are written in a macrofile.

examples of macrofile are found in \$E14\_TOP\_DIR/examples/gsim4test

- edit e14.mac as follows

```
/GsimPhysicsListFactory/QGSP_BERT/buildAndRegister
/GsimPrimaryGeneratorActionFactory/GsimGeneralParticleSource/buildAndRegister gps
#/gps/particle KLpi0pi0
/gps/particle KLpi0nunu
/gps/energy 1 GeV
/gps/position 0 0 -1 m
/gps/direction 0 0 1
#/control/execute oglcini.mac
#/GsimEventAction/setVisualizationMode 1
#/GsimEventAction/setAccumulationNumber -1
/GsimTrackingAction/setBriefTrackStore true
```

**change a kind of particle to KLpi0nunu**

**comment out viewer settings**

• • • • •

```
#/GsimSpectrumFactory/GsimE14Spectrum/buildAndRegister
#/GsimSpectrum/GsimE14Spectrum/addSpectrum Special
/GsimSpectrumFactory/GsimE14SurveySpectrum/buildAndRegister
/GsimSpectrum/GsimE14SurveySpectrum/addSpectrum Special
```

```
/GsimDetectorManager/update
/run/initialize
#/run/beamOn 20
```

**change momentum spectrum  
from E14Spectrum ( a bit old )  
to E14SurveySpectrum.**

# clustering

```
Terminal — ssh — 80x24
[info] [Thu Aug 2 12:24:27 2012] [MemSize = 68242 | RES = 57681 | SHR
= 7347] ### 100 % of events processed.
[info] [Thu Aug 2 12:24:27 2012] [MemSize = 68242 | RES = 57684 | SHR
= 7350] number of event = 100000 User=4850.18s Real=4857.25s Sys=6.15s
```

Graphics systems deleted.

Visualization Manager deleting...

```
[sato@xenys KLpi0nunuAna]$ mkdir clusteringoutput
```

```
[sato@xenys KLpi0nunuAna]$ $E14_TOP_DIR/bin/e14clustering gsimoutput/KLpi0nunu.r
oot clusteringoutput/KLpi0nunu.root ← make cluster from gsim4 output file
```

```
input file: gsimoutput/KLpi0nunu.root
```

```
output file: clusteringoutput/KLpi0nunu.root
```

```
[info] DigiReader::addDetector: ID 13:os.
```

```
[info] read /net/fs/powval3/koto/e14lib/installed20120731/e14/share/csimap/cs
i_dettree.root as CsiMap-data file.
```

```
[info] CsiMap:: # of crystals == 2716
```

```
Error in <TTree::SetBranchStatus>: unknown branch -> CC01.
```

```
Error in <TTree::SetBranchStatus>: unknown branch -> CC02.
```

```
Error in <TTree::SetBranchStatus>: unknown branch -> os.
```

```
of entry in input tree ==100000
```

```
start loop analysis for 100000 events...
```

```
0%
```

```
1%
```

**start  
clustering  
program**

# g2ana ( reconstruct pi0 )

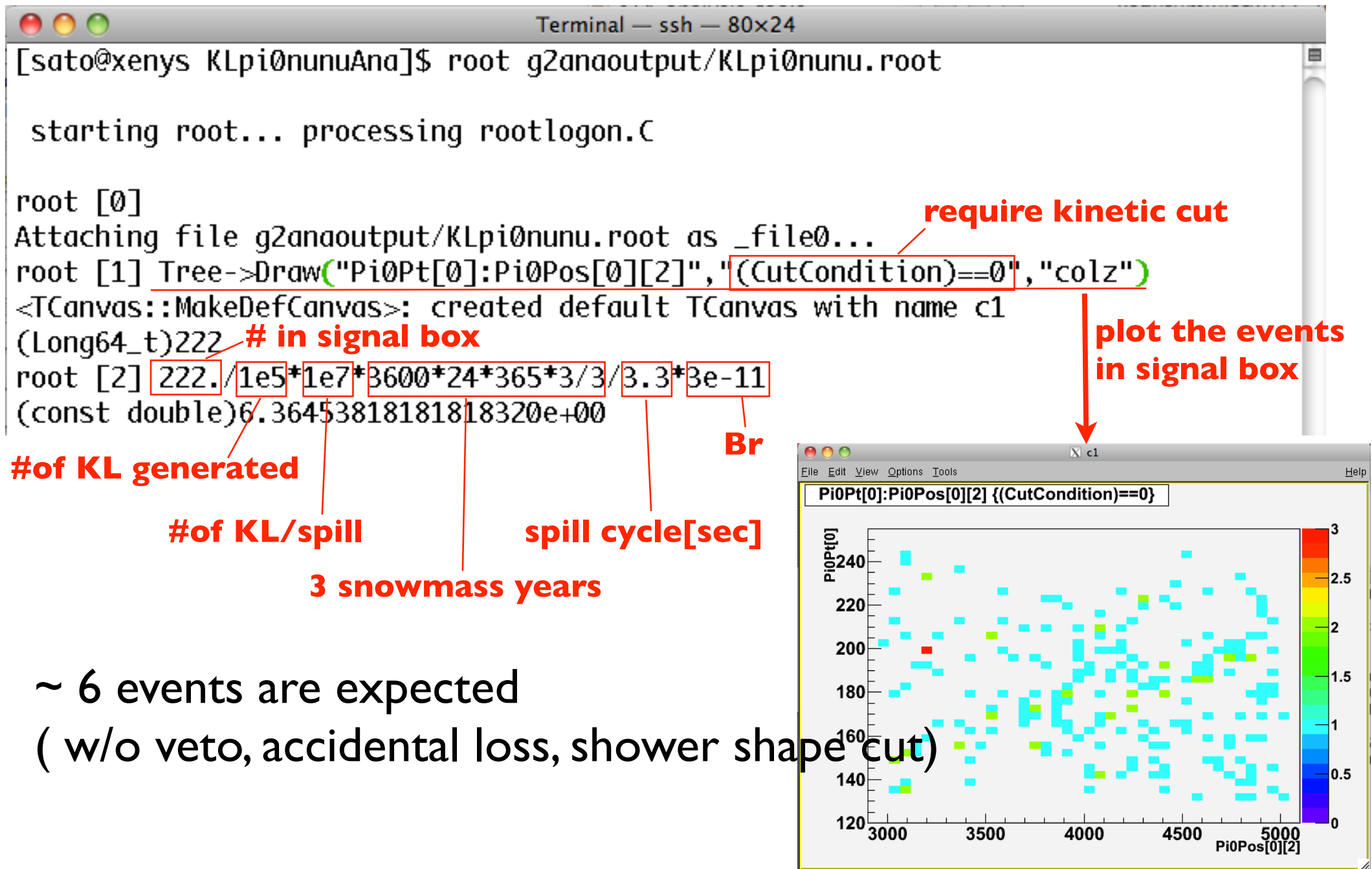
```
Terminal — ssh — 80x24
97%
98%
99%
finish!
[sato@xenys KLpi0nunuAna]$ mkdir g2anaoutput
[sato@xenys KLpi0nunuAna]$ $E14_TOP_DIR/bin/e14g2ana clusteringoutput/KLpi0nunu.
root g2anaoutput/KLpi0nunu.root ←reconstruct pi0 from e14clustering output file
input file: clusteringoutput/KLpi0nunu.root
output file: g2anaoutput/KLpi0nunu.root
start loop analysis
of entry : 100000
0%
[info] CsiMap:: # of crystals == 2716
read /net/fs/powval3/koto/e14lib/installed20120731/e14/share/gamma/mapFineCsi.f.
dat
.
60%
70%
80%
90%
[sato@xenys KLpi0nunuAna]$
```

**e14clustering** ↓

**start reconstruction program** ↓

**reconstruction program** ↓

# result



demonstration (2)  
compile your code  
with el4library

consider you want to compile test.cc that use class Cluster.

```
$ mkdir test make working directory
```

```
$ cd test/
```

```
$ emacs test.cc& edit test.cc
```

```
$ cp $E14_TOP_DIR/examples/e14gnana/e14clustering/REQUIREMENTS .
```

```
$ cp $E14_TOP_DIR/examples/e14gnana/e14clustering/Makefile . copy Makefile and REQUIREMENT
```

```
$ emacs Makefile edit Makefile
```

```
$ make
```

```
Making dependency for file test.cc ...
```

```
Compiling test.cc ...
```

```
Using granular libraries ...
```

```
Linking test ...
```

```
... Done!
```

```
$ bin/test execute
```

```
hello, cluster work as expected
```

test.cc

```
#include <iostream>
#include <cluster/Cluster.h>

int main() {
 Cluster cluster;
 std::cout<<"hello, cluster"<<std::endl;
}
```

REQUIREMENT

```
#config
gnana
csimap
cluster
gamma
pi0
klong
rec2g
E14Fsim
GsimData
GsimPersistency
```

**name of libraries  
are defined in  
REQUIREMENT**

Makefile

```
-*- Makefile -*-

#MAINS = e14clustering
MAINS = test ← edit MAINS
PACKAGE = $(shell basename `pwd`)
INSTALL_LIBS = no
#INSTALL_BINS = yes ← change yes⇒no
INSTALL_BINS = no

include $(E14_CONFIG_DIR)/Makefile.g4
```