

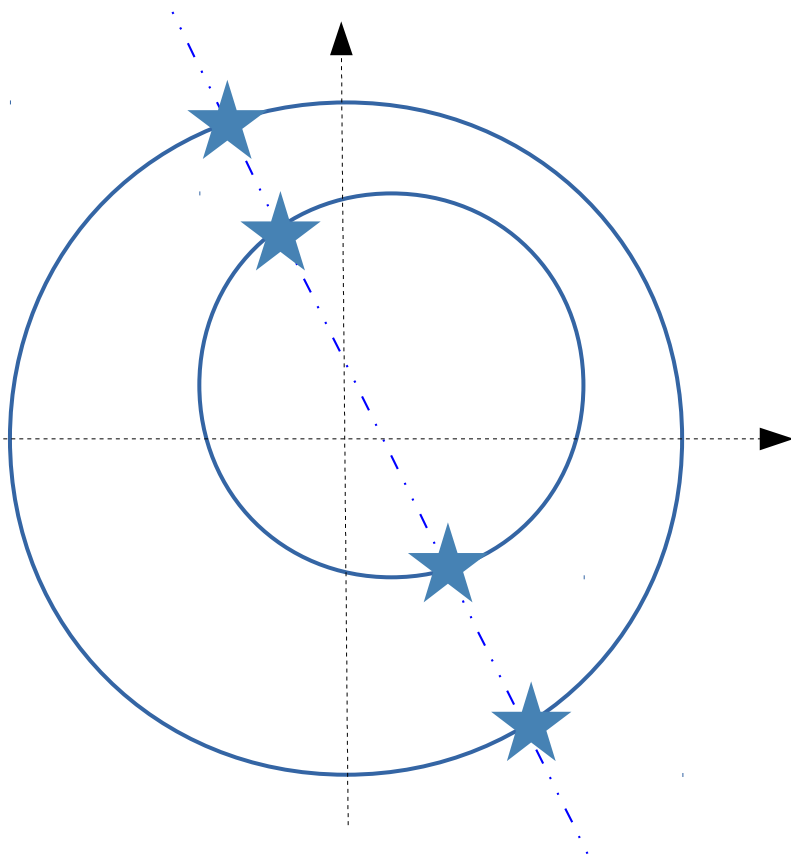
Update on QA procedure and use of CgemLineFit in it

Lia Lavezzi

2020-10-29

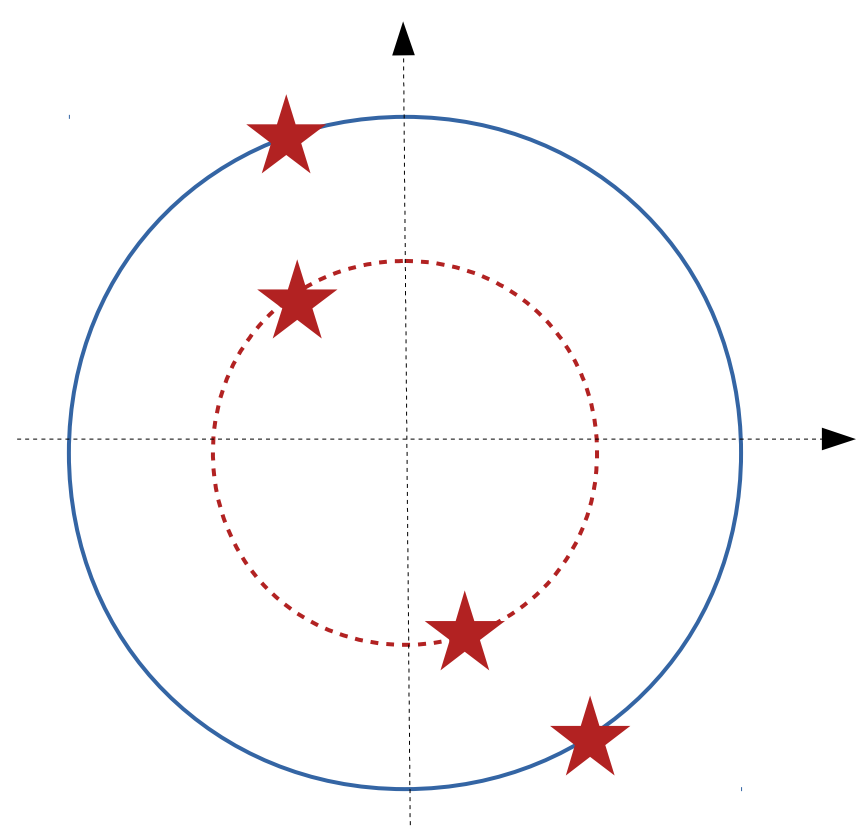
Fixed CgemCosmicRayQA

- usage of alignment functions
- computation of intersections of the fit track to the planes
- computation of residual distros



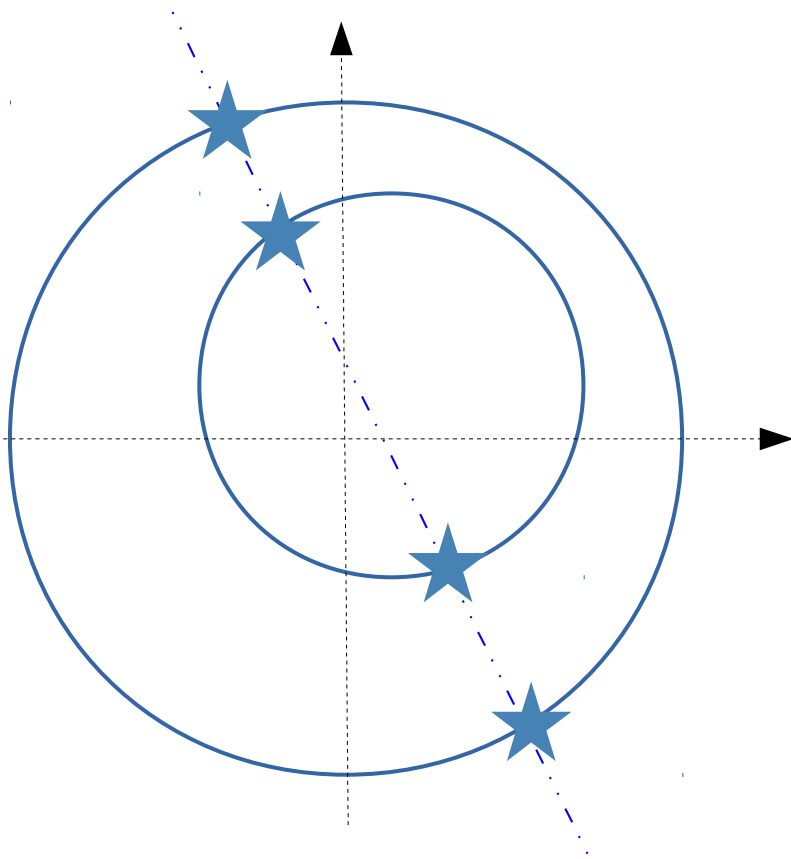
REALITY

- The dotted line is the real cosmic ray
- The blue stars are the positions where the cosmic ray is detected on the layers as positioned in the real setup



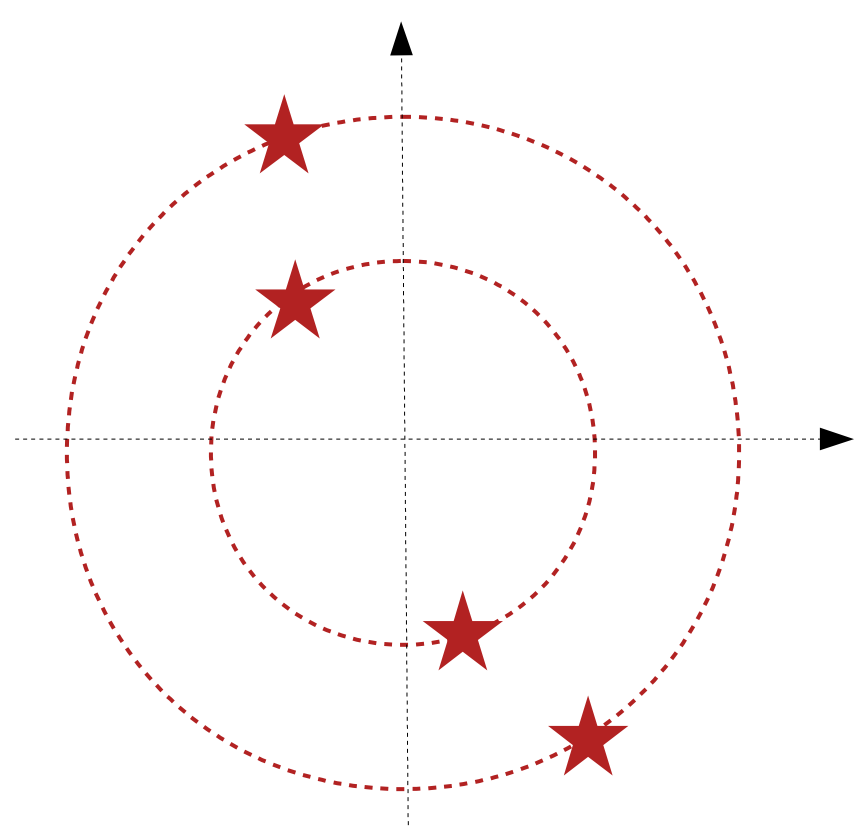
EXPERIMENTAL POINTS before alignment

- The red stars are the measurements on the layers, they are not aligned since the position of the layers is not corrected
- These are the points without alignment



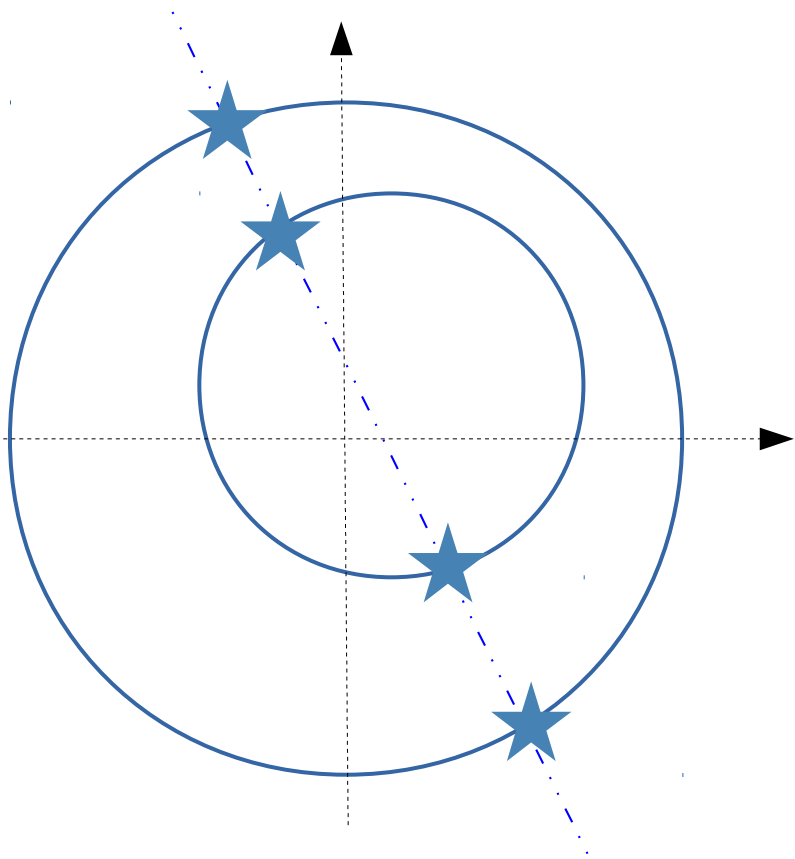
**GLOBAL (aligned)
reference frame**

- Track from fit
- Intersections of the fitted track on the layers, coordinates x, y, z
- Point Of Closest Approach (POCA) of the fitted track to $(0,0,0)$

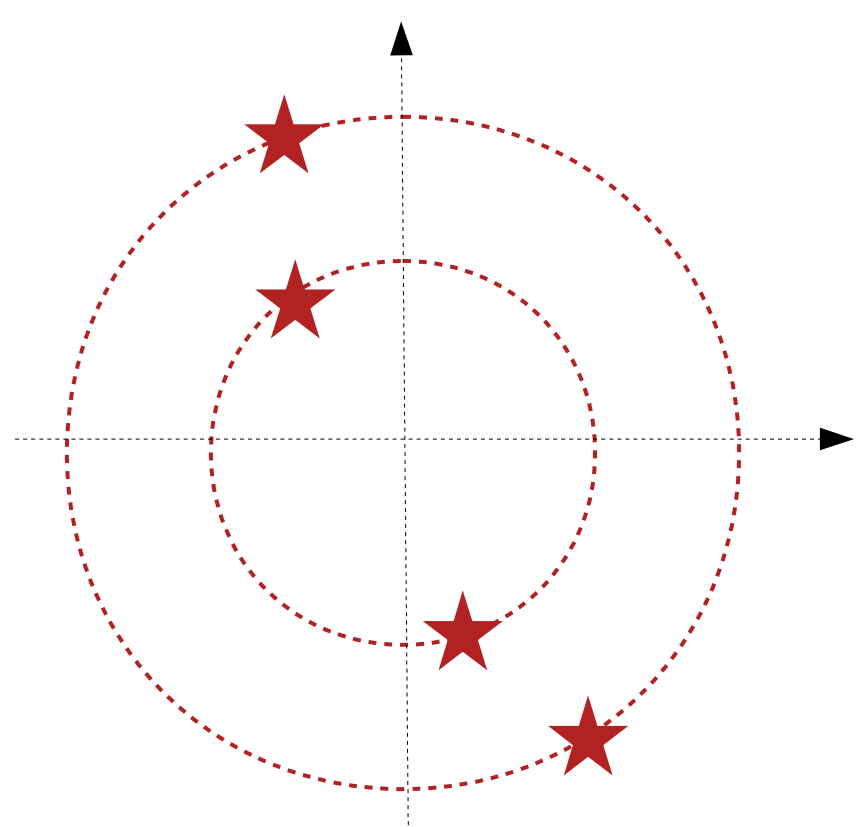


**LOCAL (not-aligned)
reference frame**

- Experimental hits
- Reco clusters 1D
- Reco clusters 2D
- Intersections of the fitted track on the layers, coordinates f, v



**GLOBAL (aligned)
reference frame**



**LOCAL (not-aligned)
reference frame**



CgemGeoAlign::point_transform

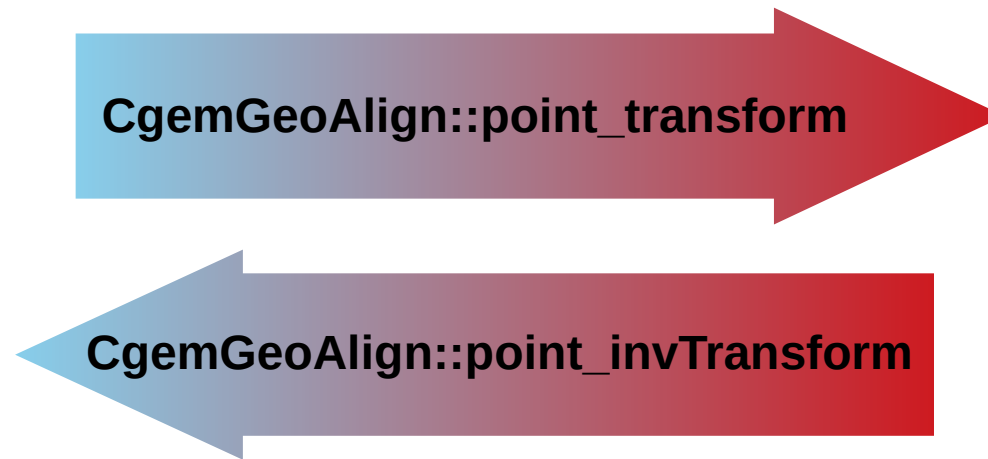


CgemGeoAlign::point_invTransform

*Thankx to
Aiqiang*

GLOBAL (aligned)
reference frame

LOCAL (not-aligned)
reference frame



MY PREVIOUS VERSION WAS WRONG

- I was computing the residuals all in the GLOBAL frame
- porting there the experimental positions with point_invTransform
- BUT I was finding the intersections of the fitted line to **non aligned** layers (LOCAL) with my **own** code → WRONG

NEW VERSION IS FIXED

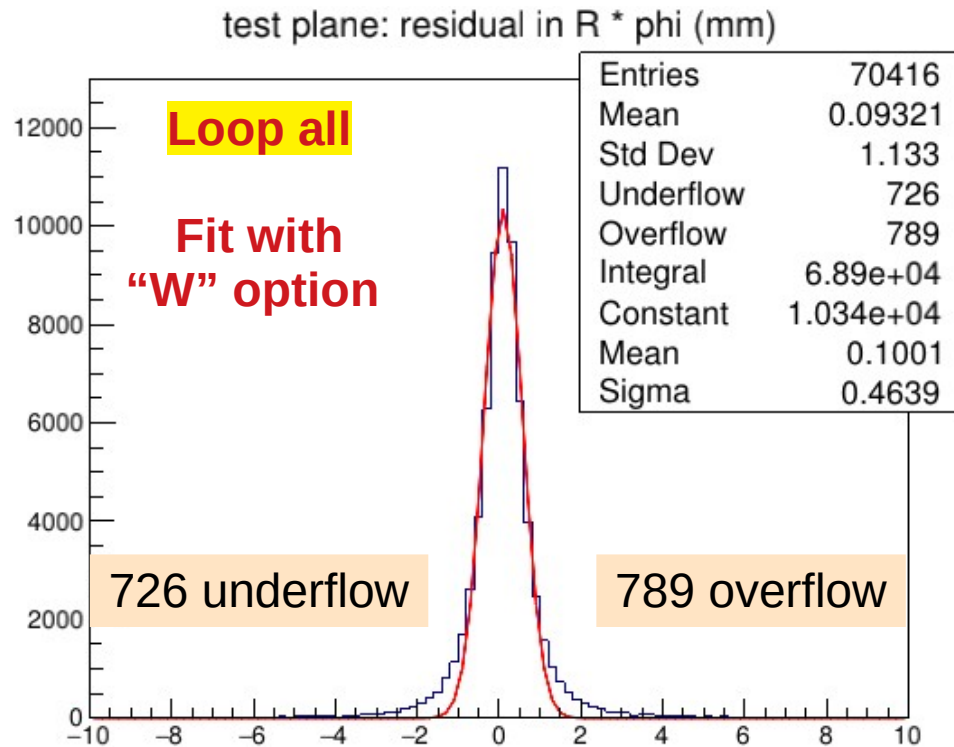
- I compute the residuals in the LOCAL frame
- I compute the intersections in the GLOBAL aligned frame with the **dedicated** function

```
bool gotit;  
  
if(align_flag==true) gotit = midplane->getPointAligned(layerid, linefit, posup, posdown, phivup, phivdown);  
else gotit = midplane->getPointIdealGeom(layerid, linefit, posup, posdown, phivup, phivdown);
```

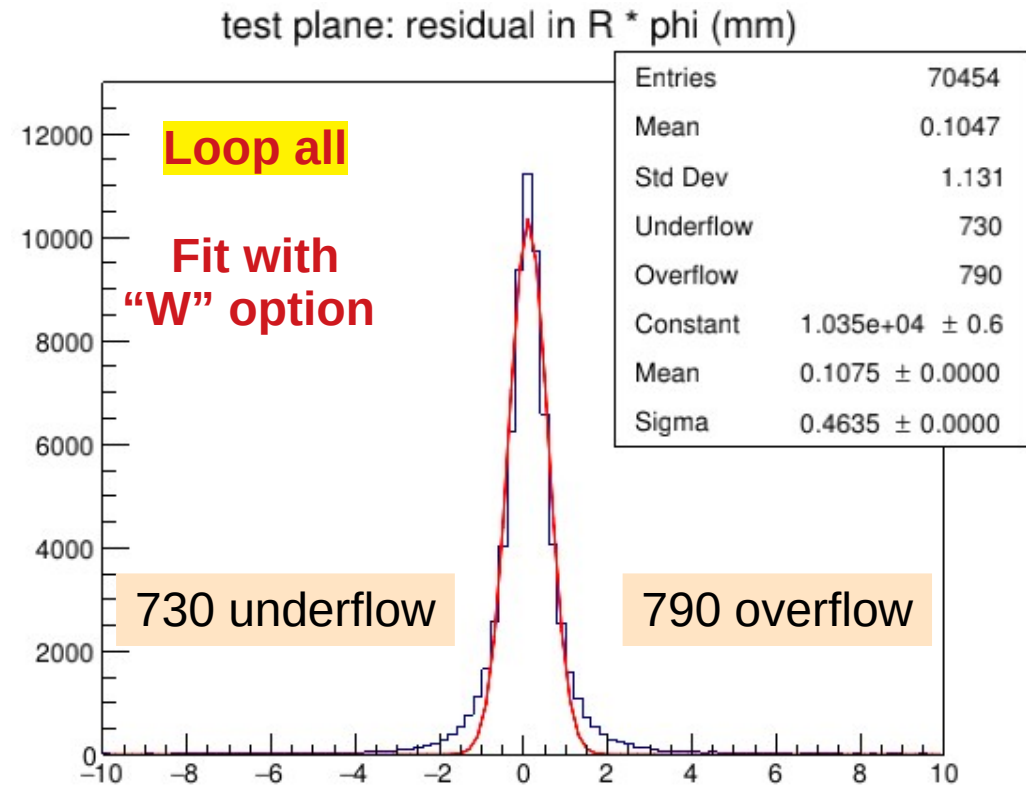
- Use ϕ and v from the LOCAL frame

residual $r^*\phi$ - L1 bottom

PREVIOUS VERSION (WRONG)



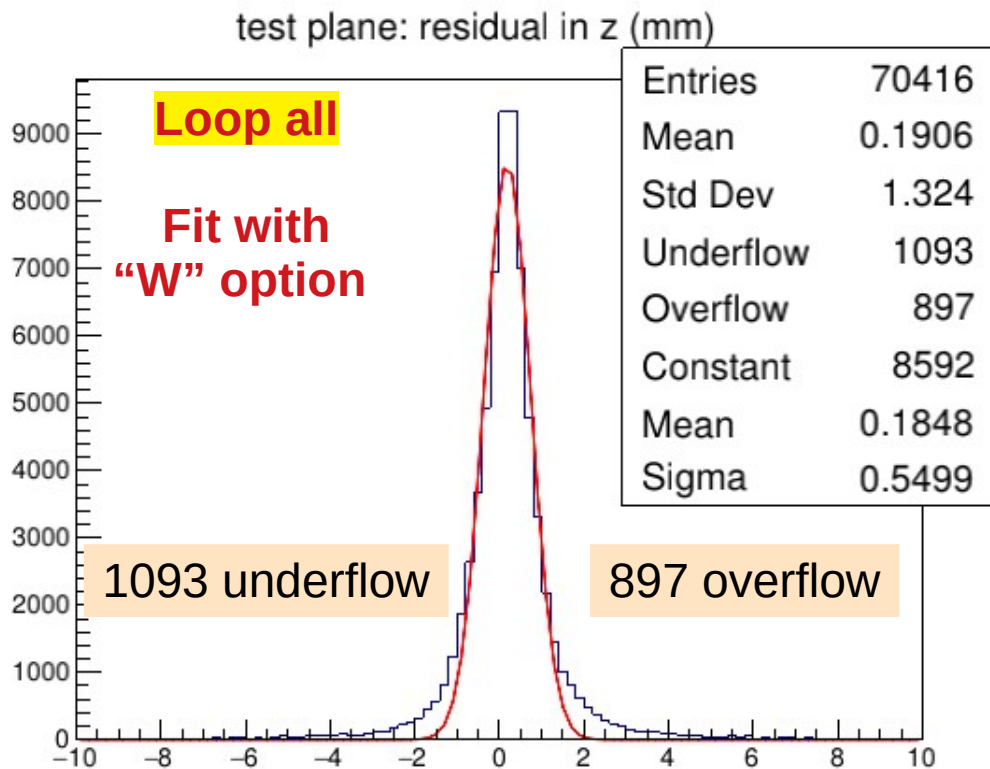
NEW VERSION (FIXED)



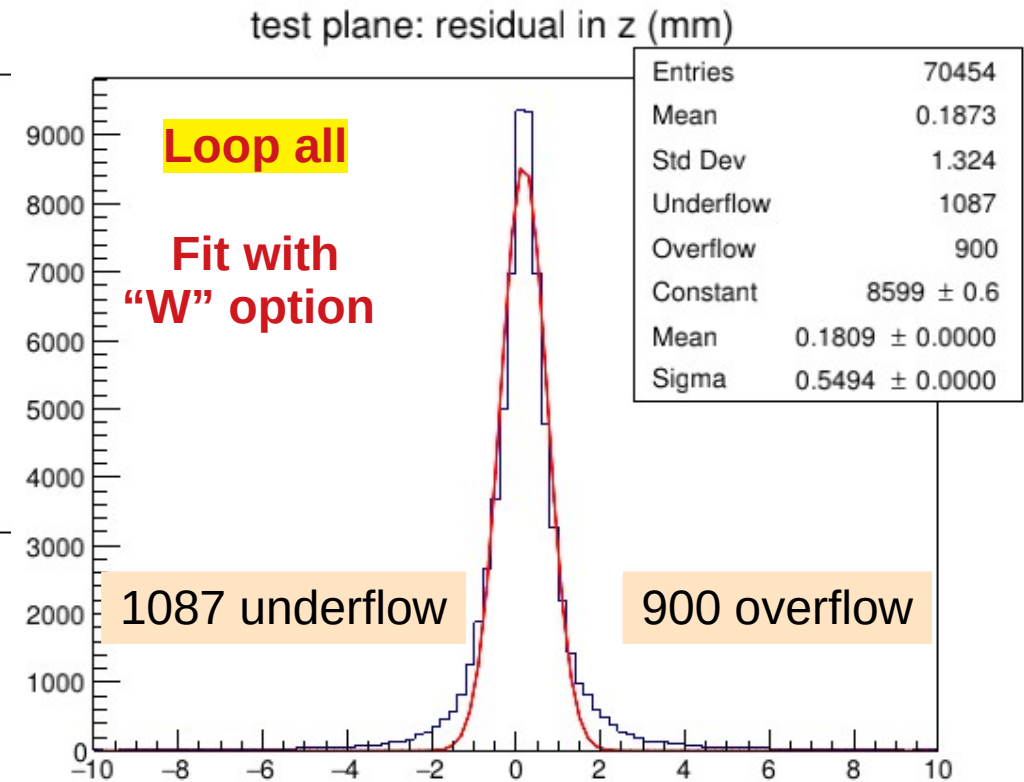
There is no actual difference in the results

residual z - L1 bottom

PREVIOUS VERSION (WRONG)



NEW VERSION (FIXED)



There is no actual difference in the results

Test 1 - use Loop All

- Use it on every test plane
- Tested alignment of residual distros
- Tested chi2 cut
- Tested cut on cluster charge

Loop All on all planes

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll** algorithm
- Test all the layers**, one by one
- L1 TOP and L2 BOTTOM have problems

L1 BOTTOM

n FITTED track	120502
n VALID track (after cuts)	75089
efficiency	0.8520955133
background component	0.0238117434

L1 TOP

n FITTED track	0
n VALID track (after cuts)	0
efficiency	-nan
background component	-nan

L2 BOTTOM

n FITTED track	6402
n VALID track (after cuts)	3581
efficiency	0.8553476682
background component	0.01396258028

L2 TOP

n FITTED track	126655
n VALID track (after cuts)	81455
efficiency	0.8554907618
background component	0.01419188509

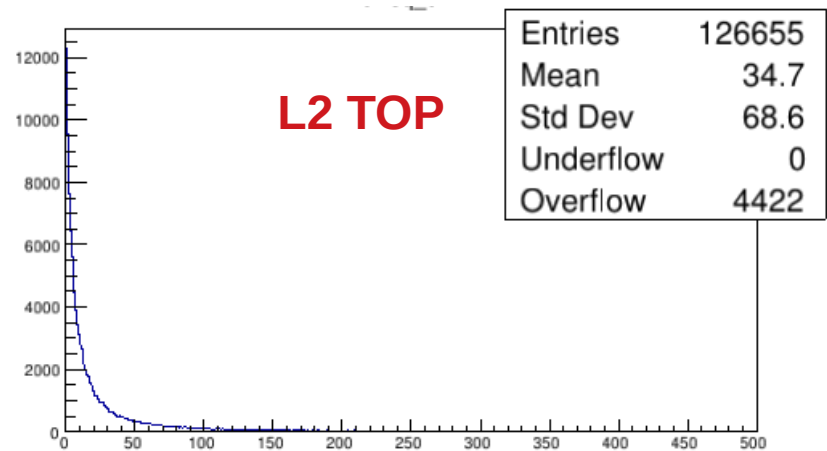
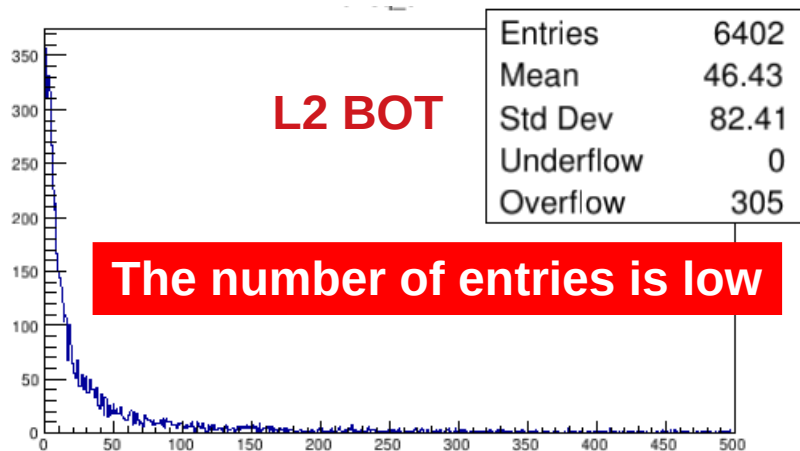
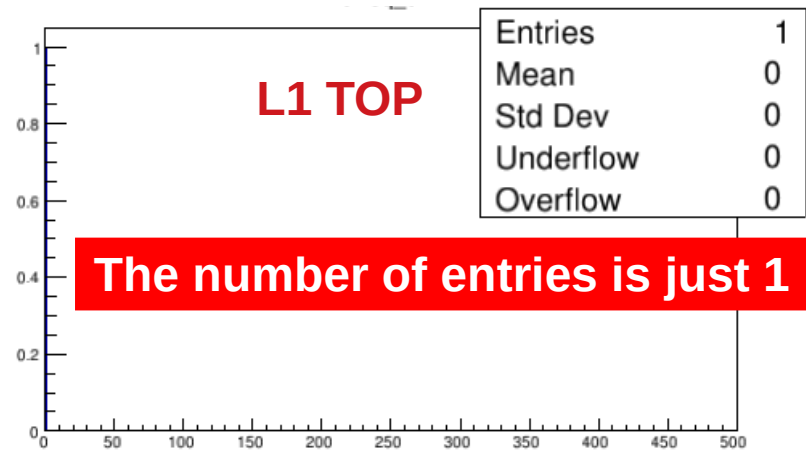
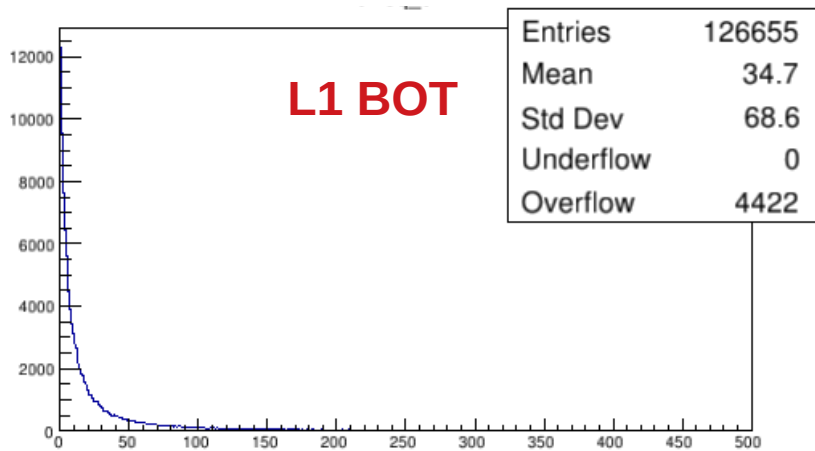
Loop All on all planes

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll algorithm
- Test all the layers, one by one
- L1 TOP and L2 BOTTOM have problems: **the LoopAll fit does not work fine here!**

χ^2 DISTRIBUTIONS

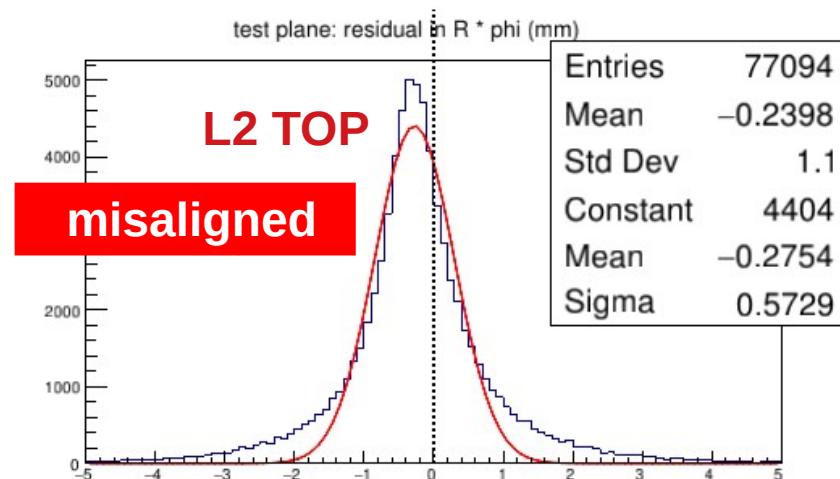
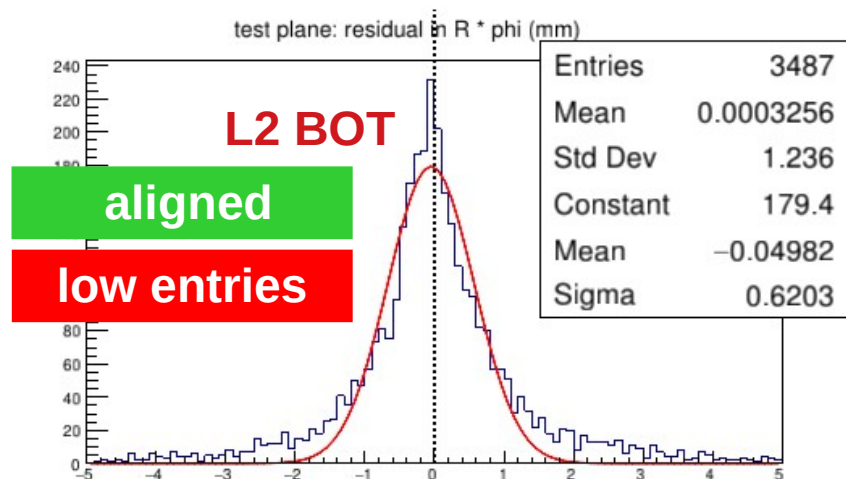
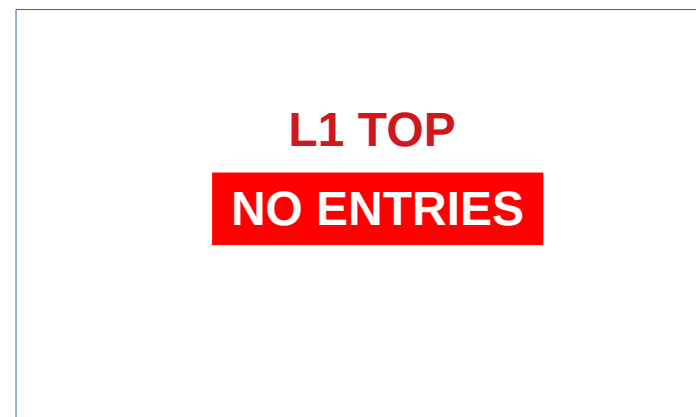
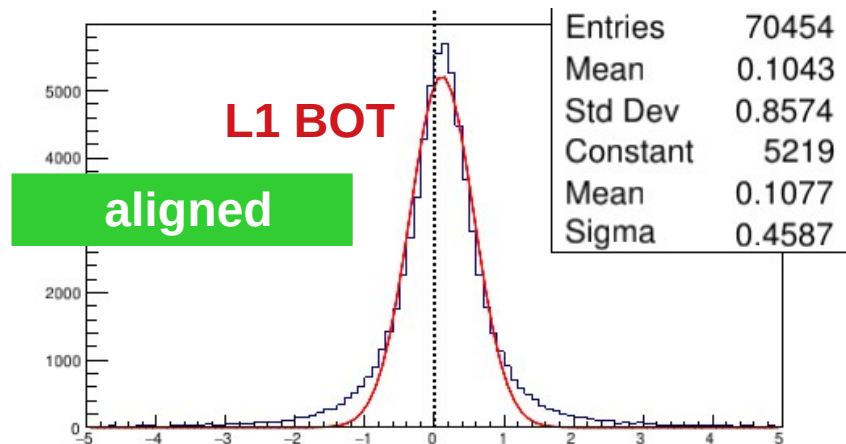


The residual distributions

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll algorithm
- Test all the layers, one by one
- Chi2 cut = 20



R * ϕ RESIDUAL DISTRO

chi2 CUT scan

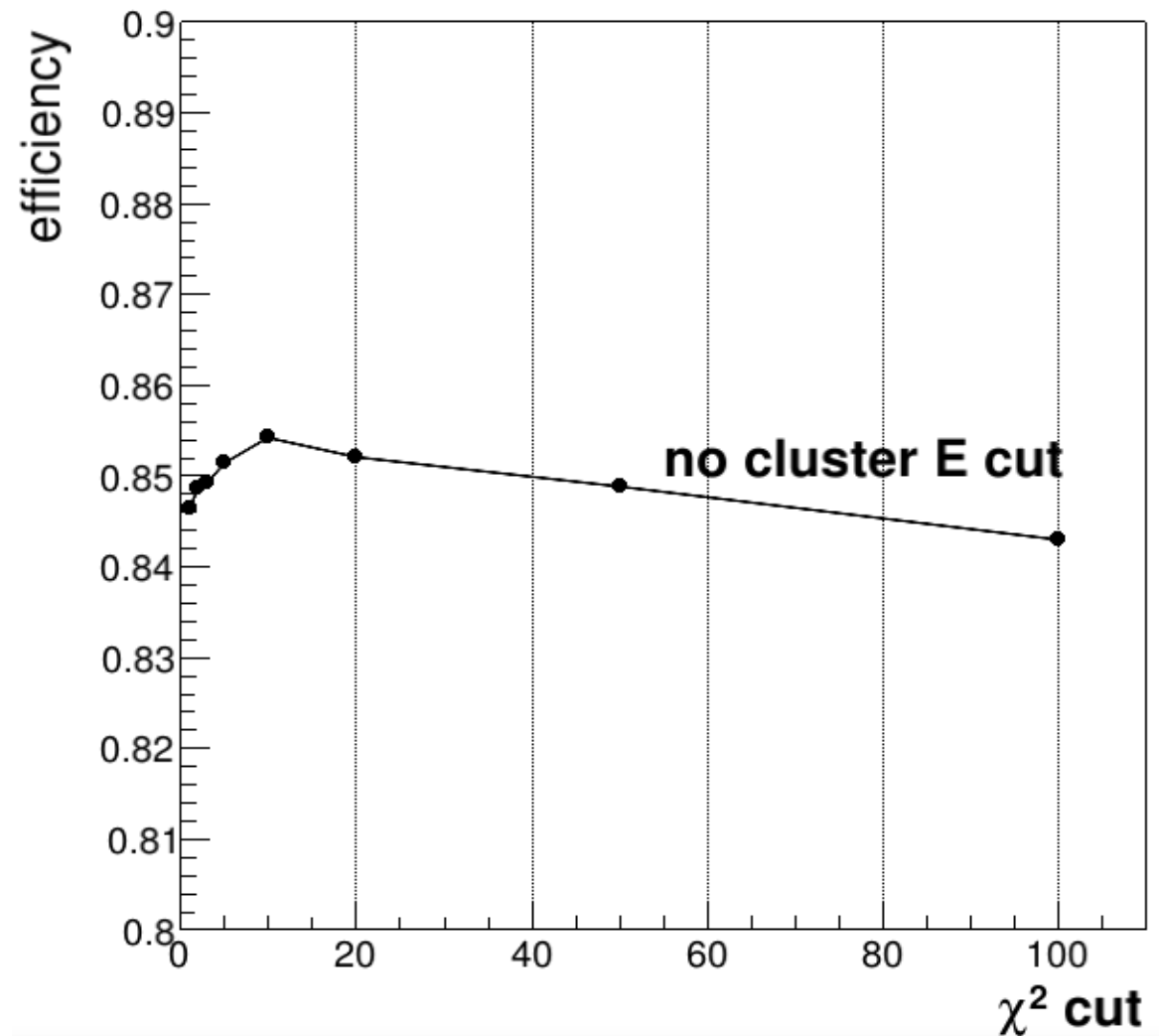
- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll** algorithm
- Run a **chi2 scan** for **L1 BOT**

Chi2 scanned in the values
[1,2,3,5,10,20,50,100]

There is not a real and relevant
difference

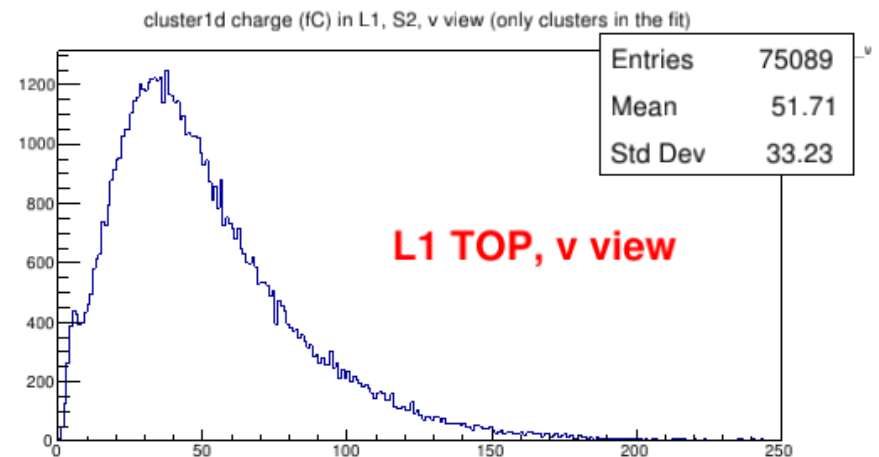
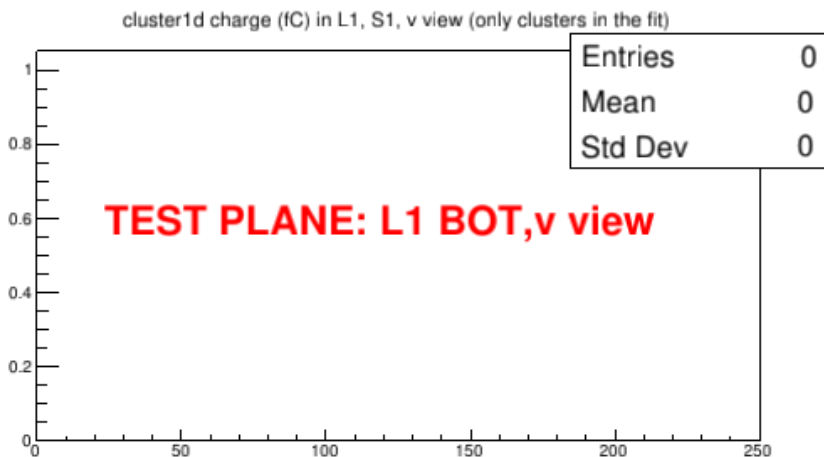
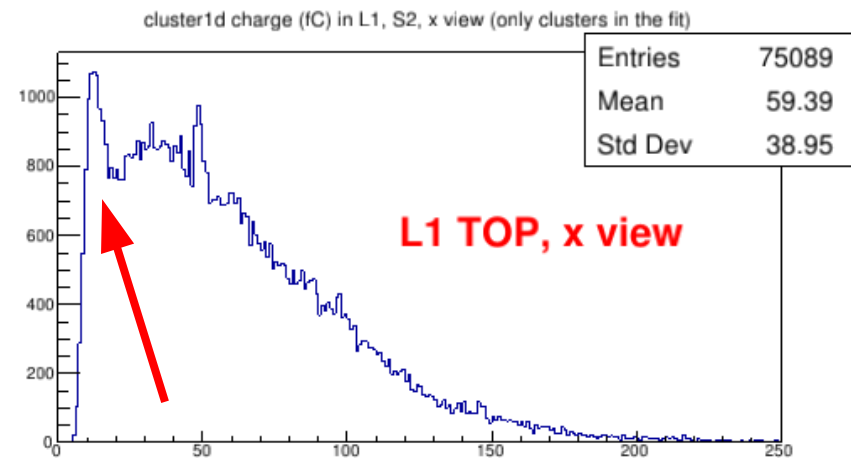
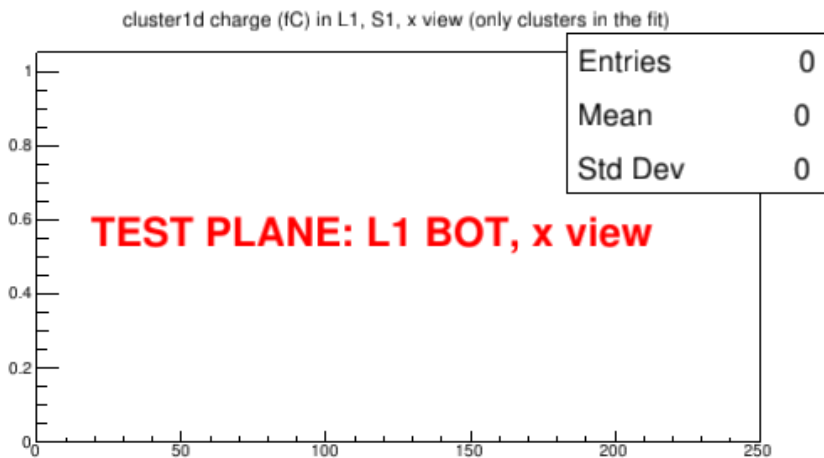


cluster energy

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll** algorithm
- See the energy distribution of the clusters used for fitting**

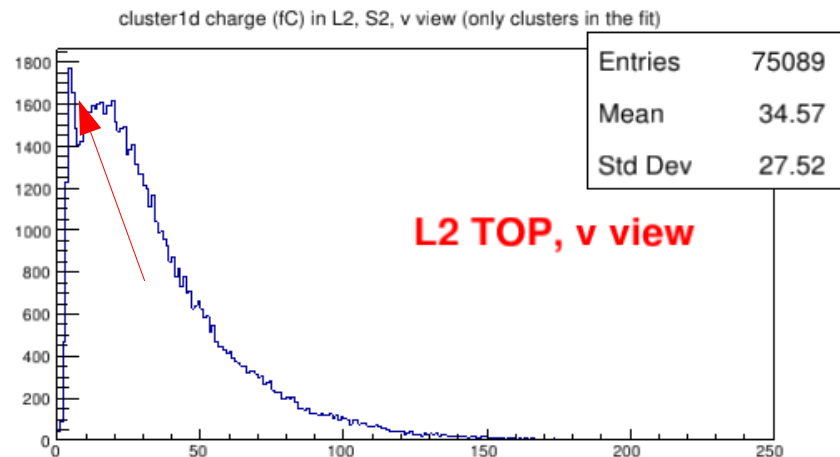
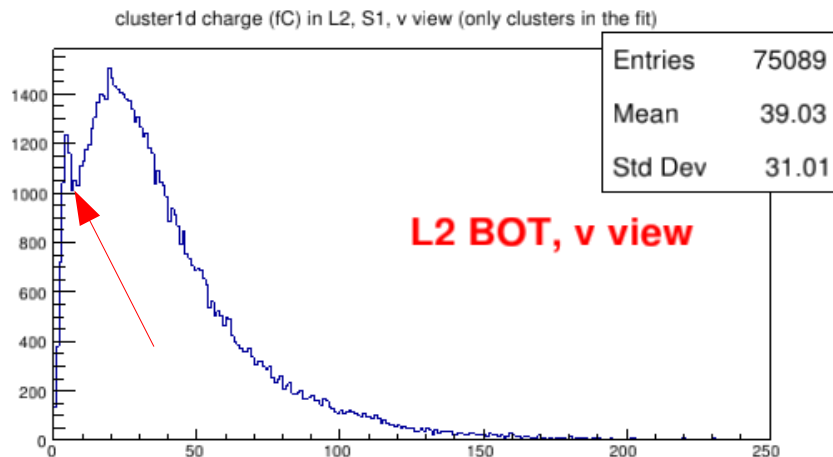
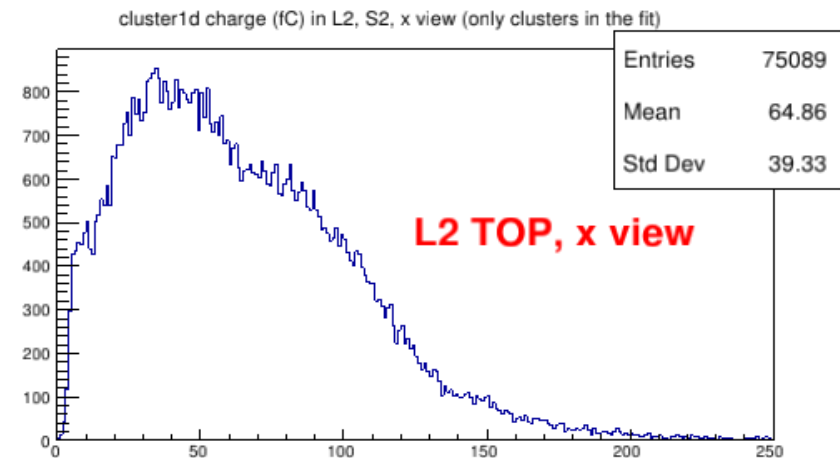
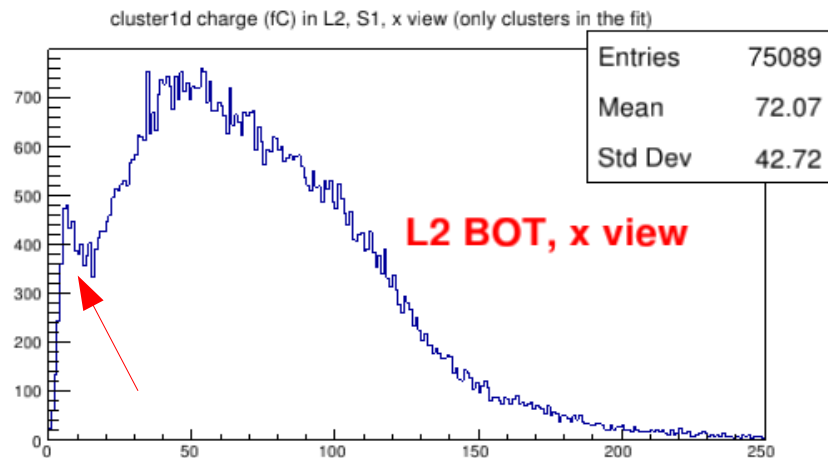


cluster energy

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll** algorithm
- See the energy distribution of the clusters used for fitting**



cluster energy

- Cuts used in my old standalone were:

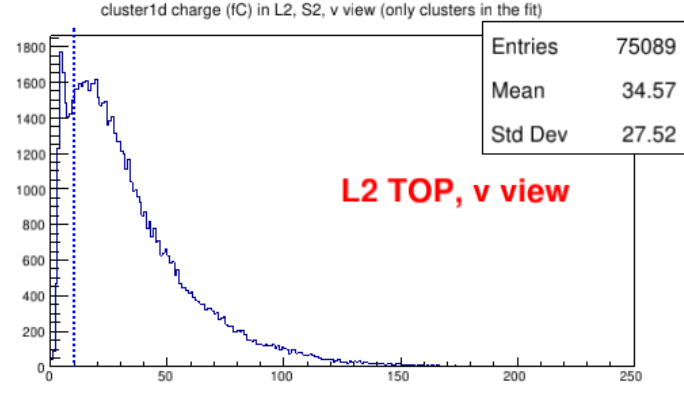
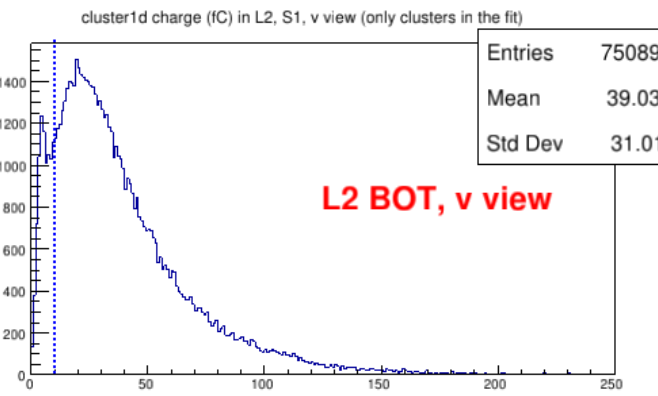
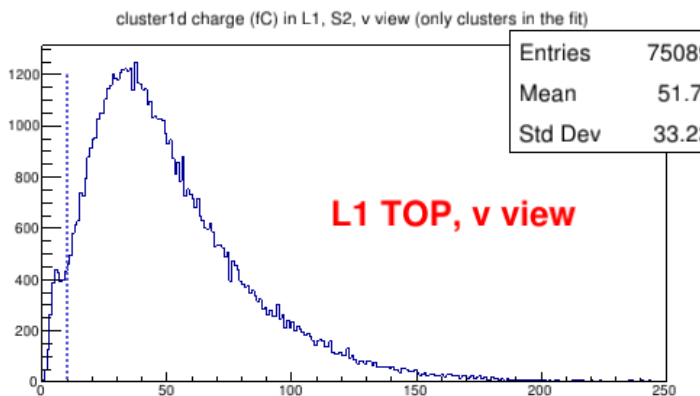
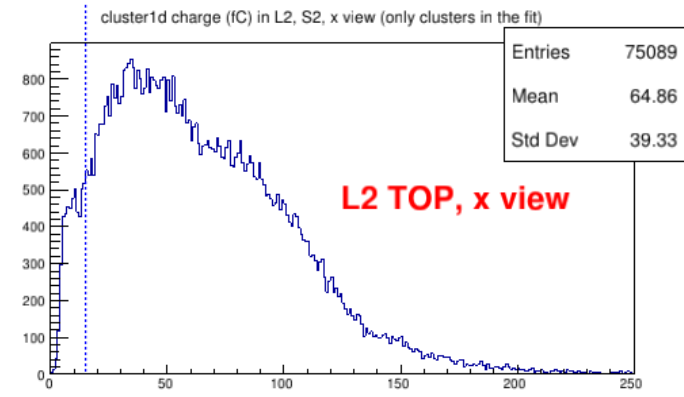
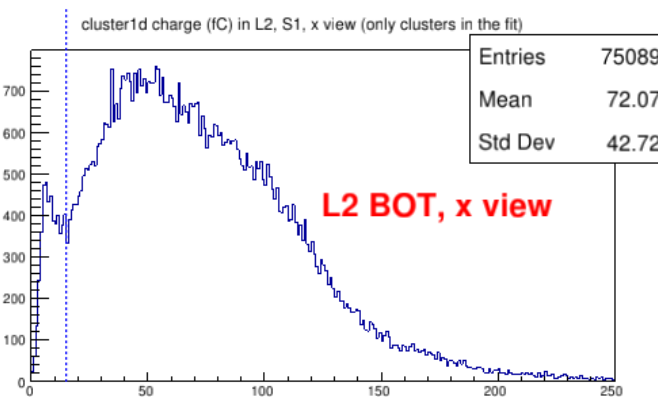
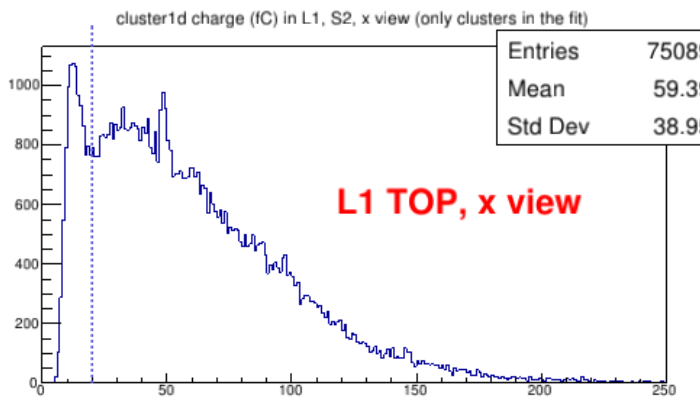
L1, x view, $Q_{\text{CLUSTER}} > 20 \text{ fC}$
 L2, x view, $Q_{\text{CLUSTER}} > 15 \text{ fC}$

L1, v view, $Q_{\text{CLUSTER}} > 10 \text{ fC}$
 L2, v view, $Q_{\text{CLUSTER}} > 10 \text{ fC}$

L1 TOP

L2 BOT

L2 TOP

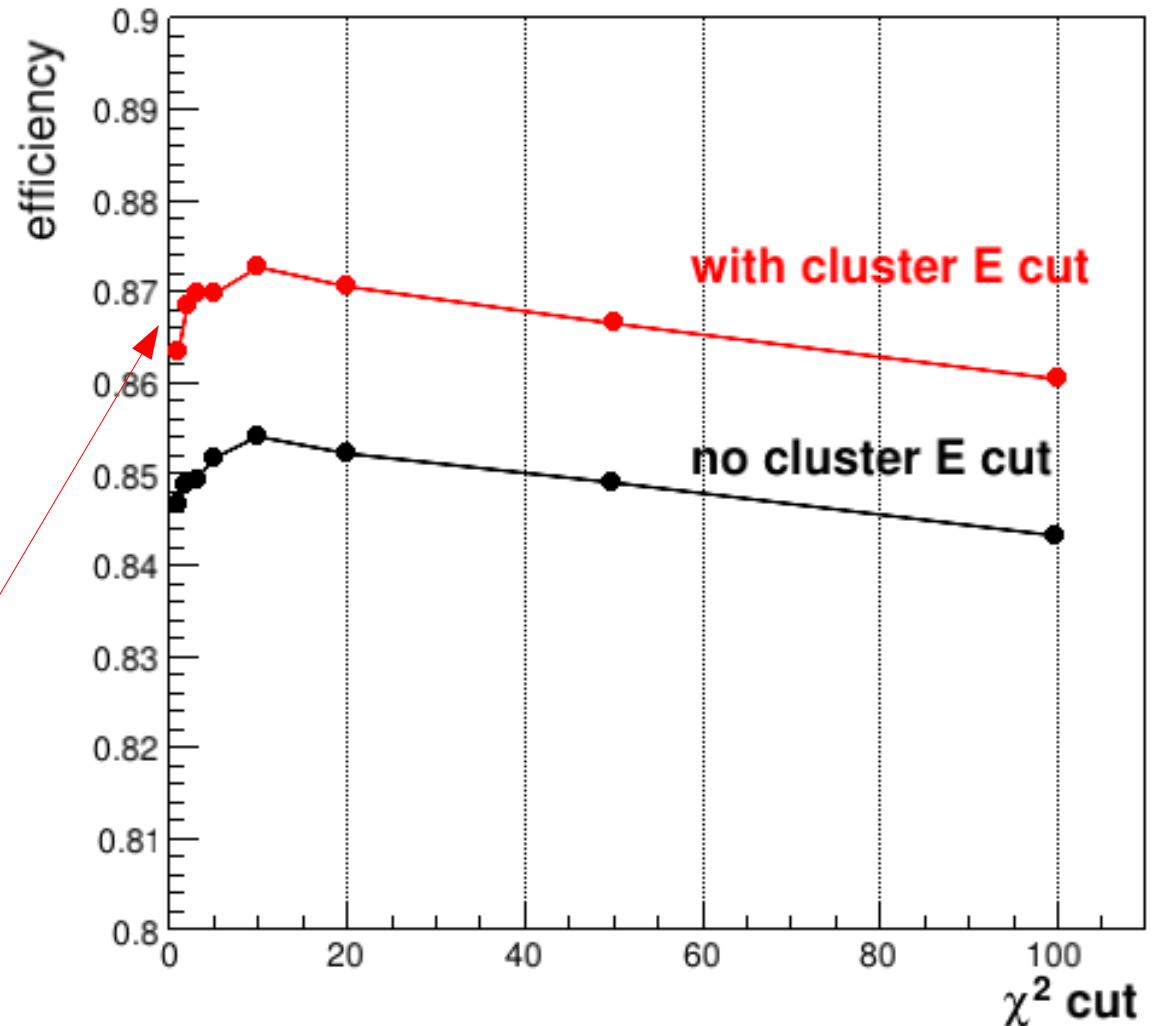


chi2 CUT scan

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll** algorithm
- Run a chi2 scan for L1 BOT**



Chi2 scanned in the values [1,2,3,5,10,20,50,100]

There is not a real and relevant difference

Also by cutting away the events where one of the tracker cluster does not pass the energy cut there is no change by chi2 cut

Test 2 - use Loop All + Max Q

- Use it on every test plane
- Tested alignment of residual distros

Test 2 – Loop All + Max Charge

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll + MaxCharge** algorithm
- Test all the layers**, one by one
- LoopAll+MaxChage has a fitting procedure which is fine for all the planes**

L1 BOTTOM

n FITTED track	153160
n VALID track (after cuts)	86453
efficiency	0.8354018947
background component	0.02558615664

L1 TOP

n FITTED track	151818
n VALID track (after cuts)	86609
efficiency	0.8479026429
background component	0.02145273586

L2 BOTTOM

n FITTED track	150586
n VALID track (after cuts)	86040
efficiency	0.8389702464
background component	0.01813110181

L2 TOP

n FITTED track	153433
n VALID track (after cuts)	88779
efficiency	0.8349834983
background component	0.01954290992

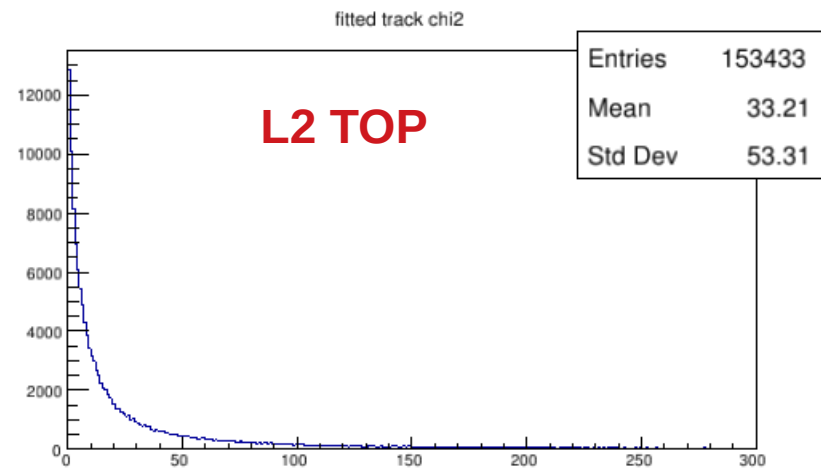
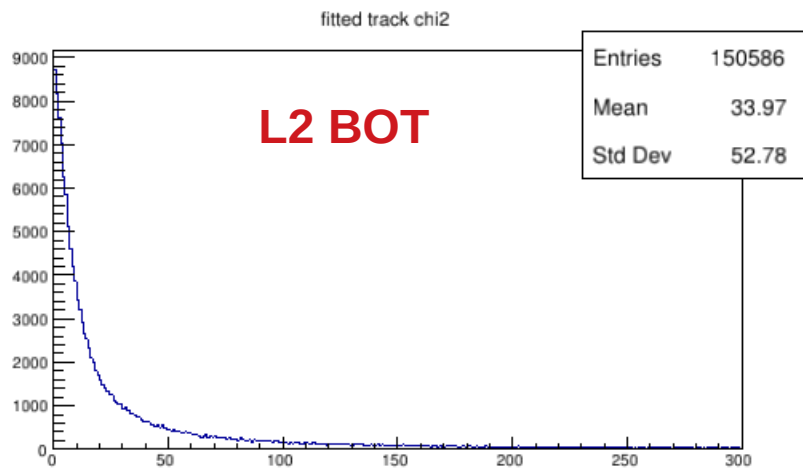
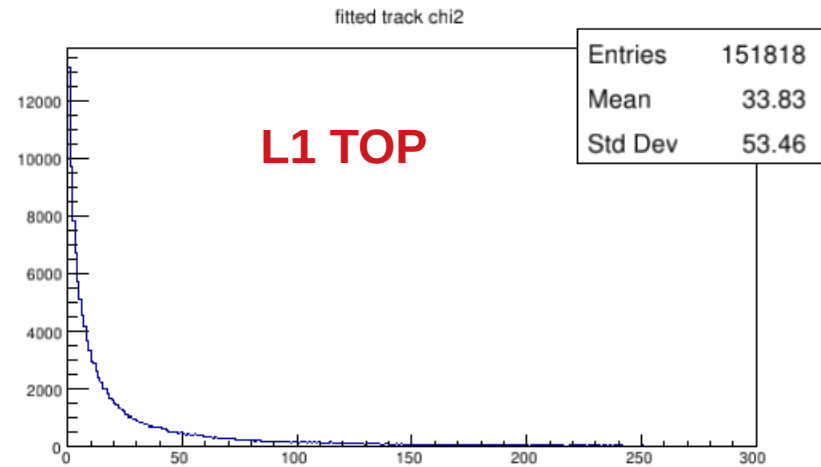
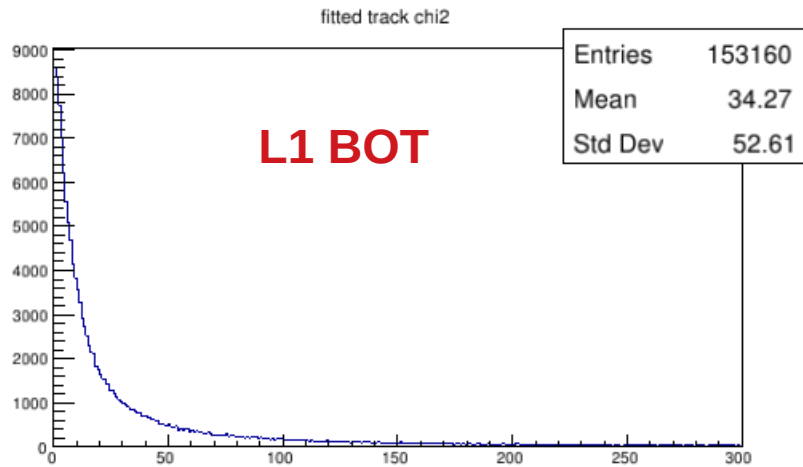
Loop All on all planes

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll + MaxCharge** algorithm
- Test all the layers, one by one**

χ^2 DISTRIBUTIONS



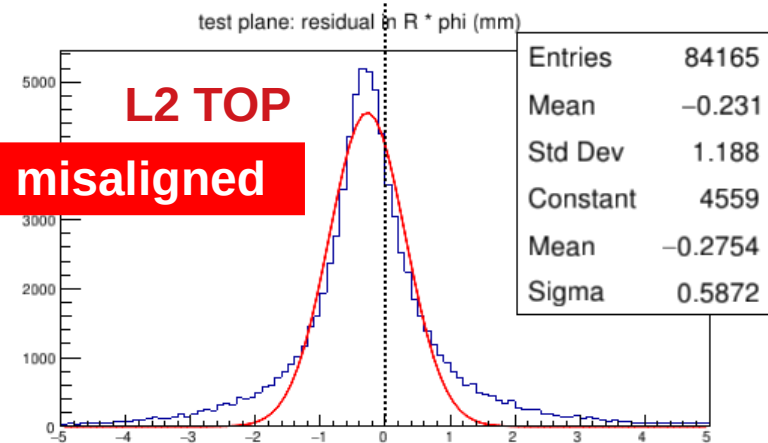
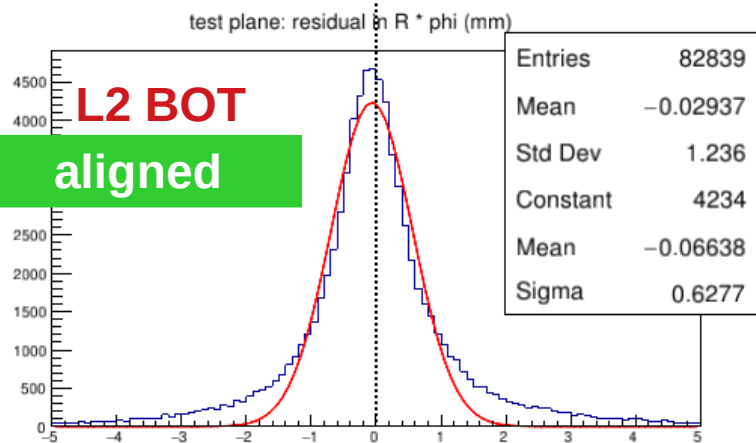
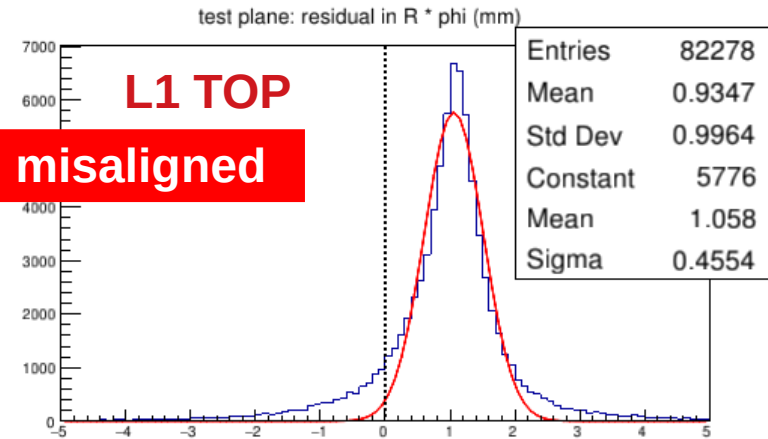
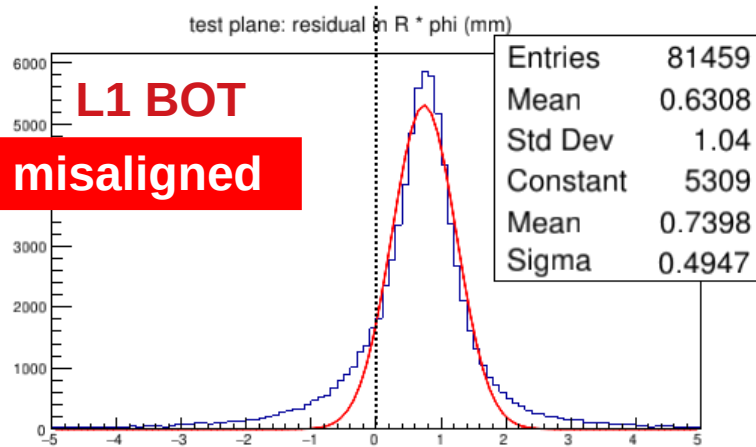
Test 2 – Loop All + Max Charge

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

- LoopAll + MaxCharge algorithm
- Test all the layers, one by one
- LoopAll+MaxChage: what about the misalignment?**

R * ϕ RESIDUAL DISTRO



Test 3 – use the 4 layers all at once

Number of fitted tracks

- Alignment **on**, parameters are:

Elements	DeltaX(mm)	DeltaY(mm)	DeltaZ(mm)	RX(rad)	RY(rad)	RZ(rad)
layer1	0.0889837	0.0000000	1.8009692	0.0000000	0.0000000	-0.0092171
layer2	0.0000000	-0.0000000	-0.0000000	0.0000000	0.0000000	0.0000000
layer3	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

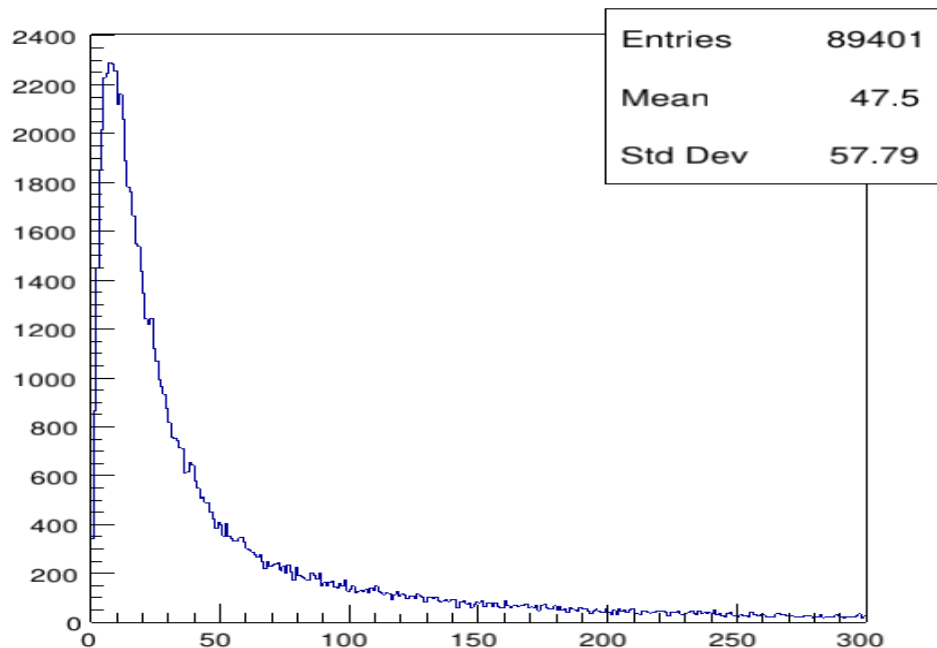
LoopAll

n FITTED track 89401
n not passing E cut 15052
n VALID track (after cuts) 20737

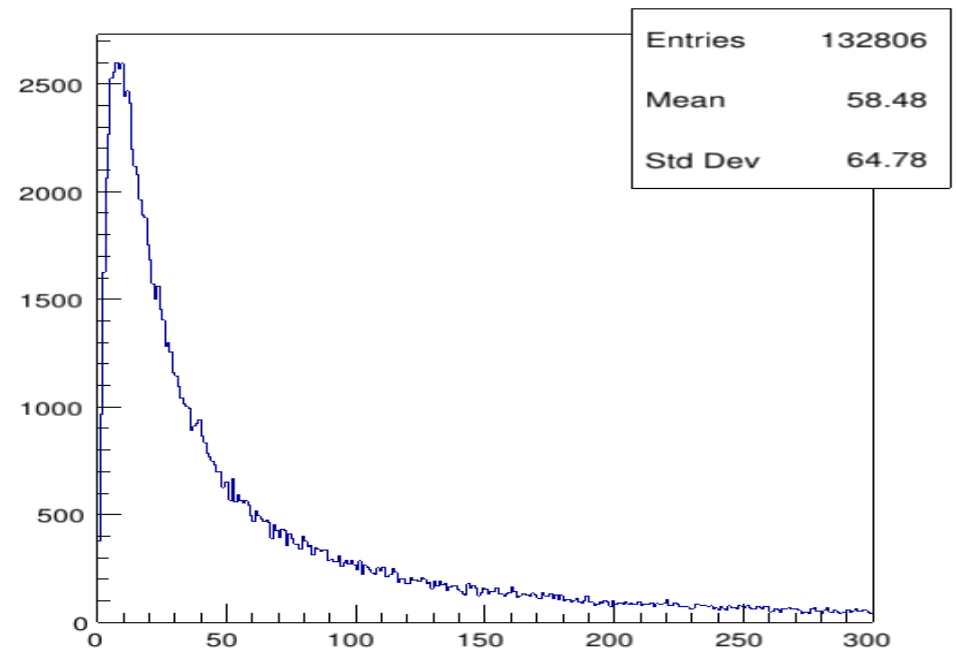
LoopAll + MaxQ

n FITTED track 132806
n not passing E cut 17535
n VALID track (after cuts) 23822

fitted track chi2

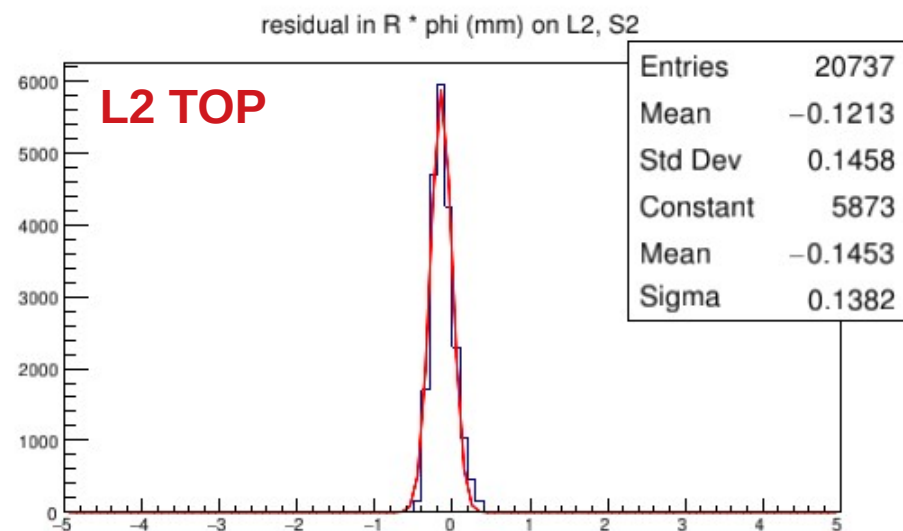
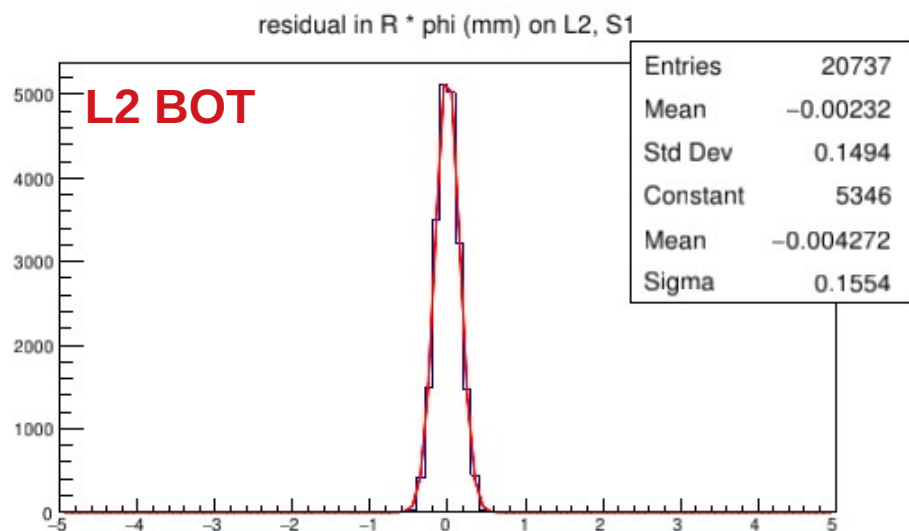
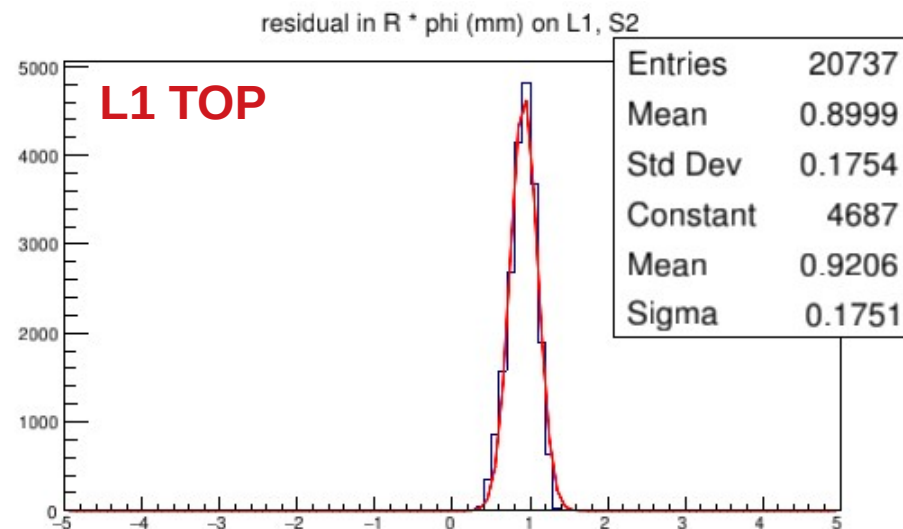
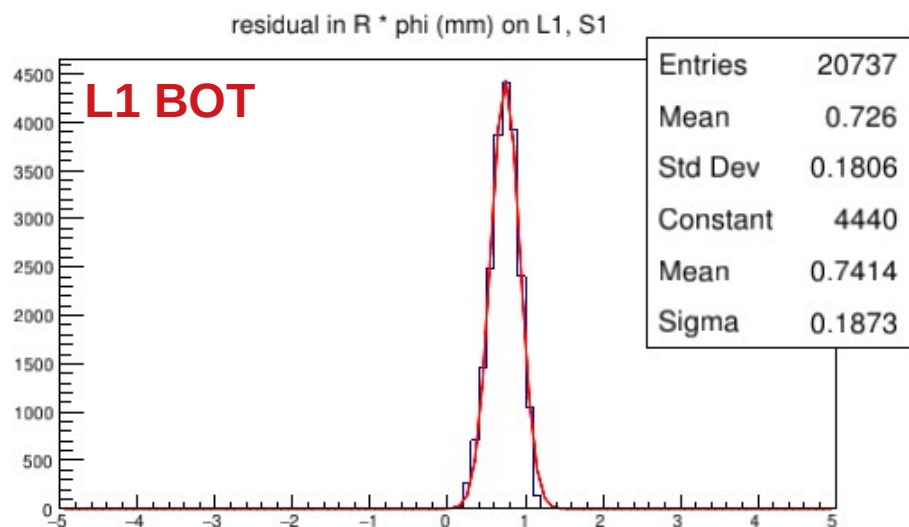


fitted track chi2



R * phi - LoopAll

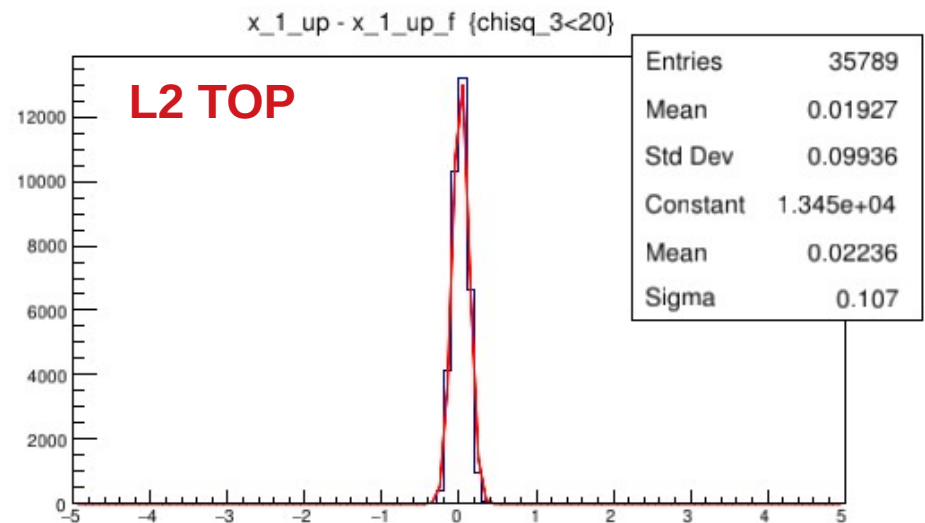
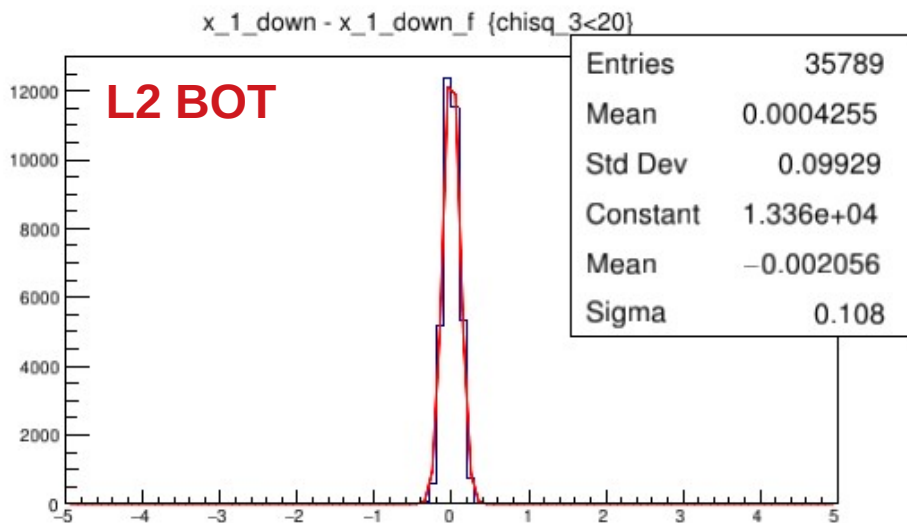
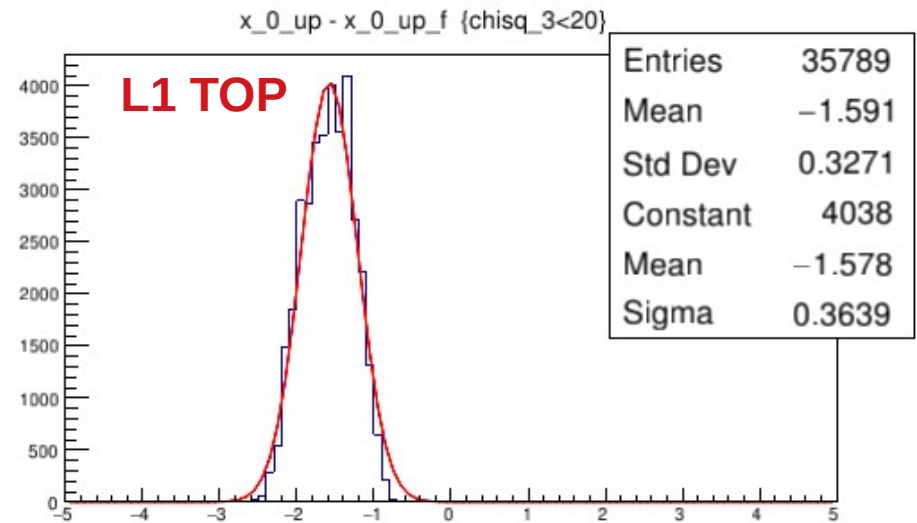
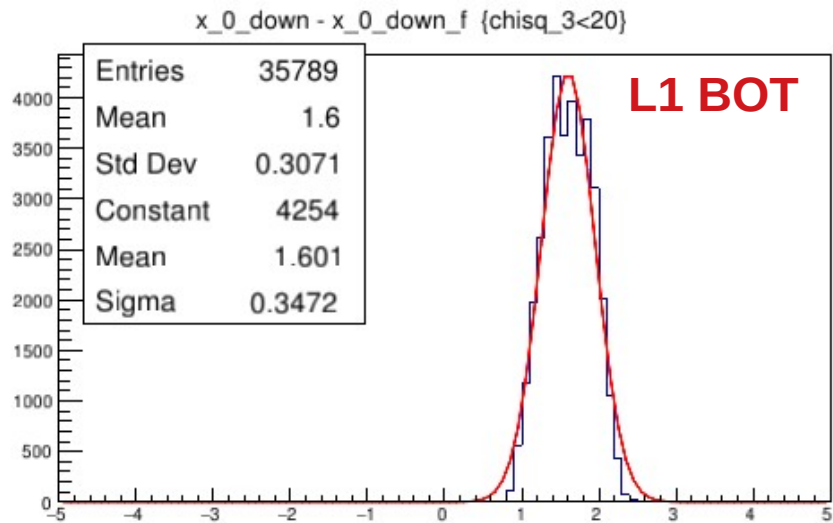
- Alignment **on**
- Chi2 cut = 20
- Energy cut on clusters



phi - LoopAll ~ HongPeng (?)

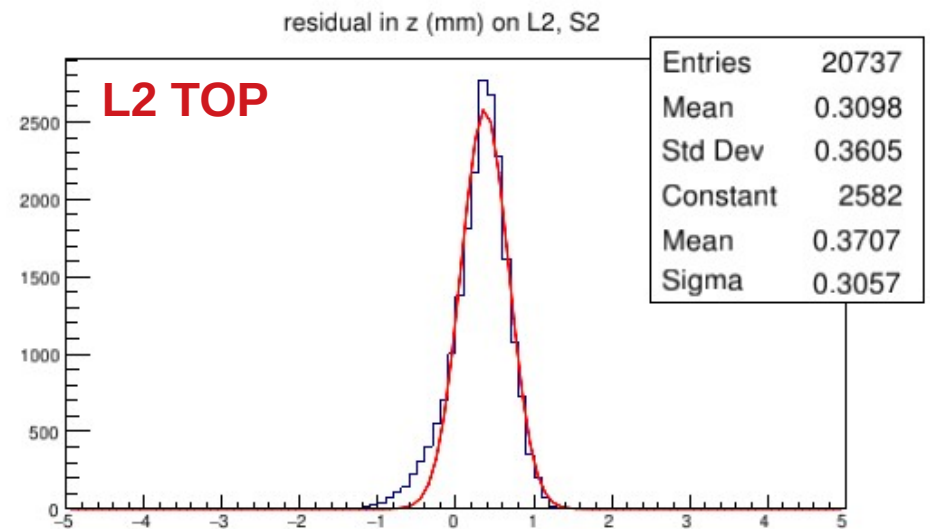
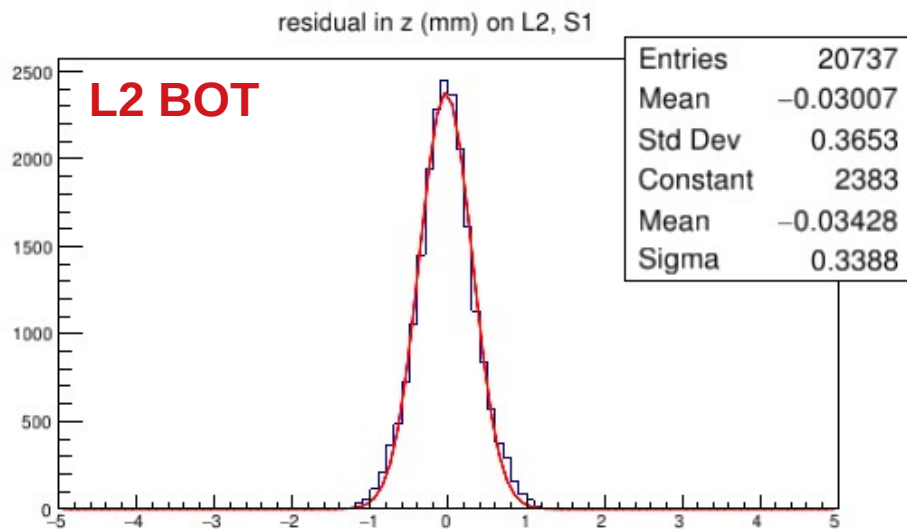
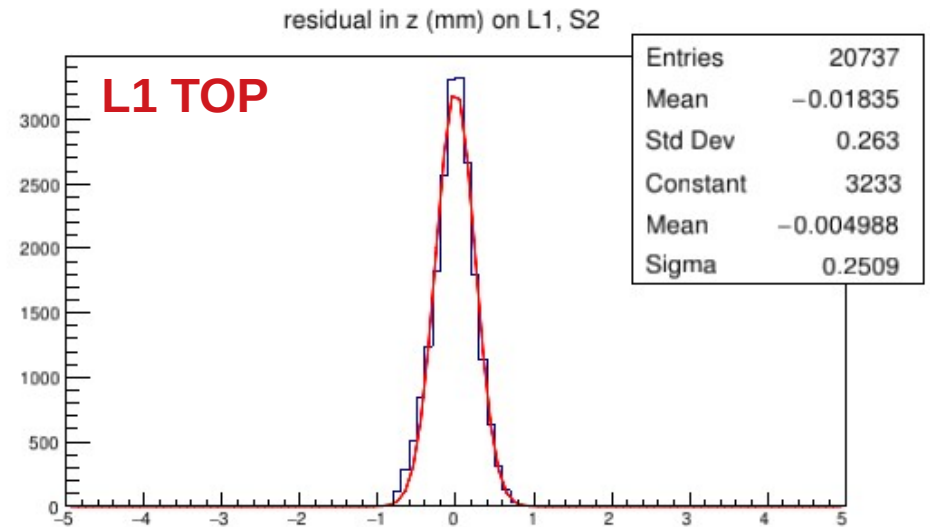
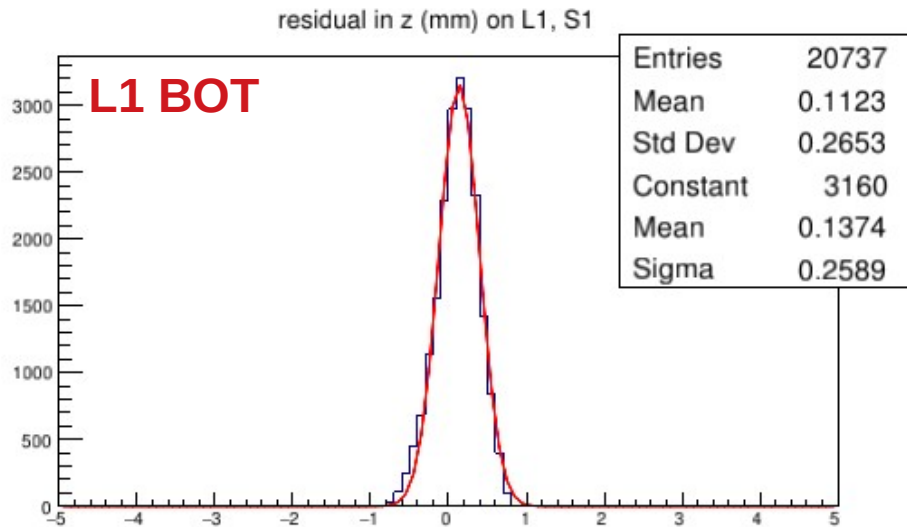
- Alignment on
- Chi2 cut = 20
- Energy cut on clusters

x_0_down - x_0_down_f



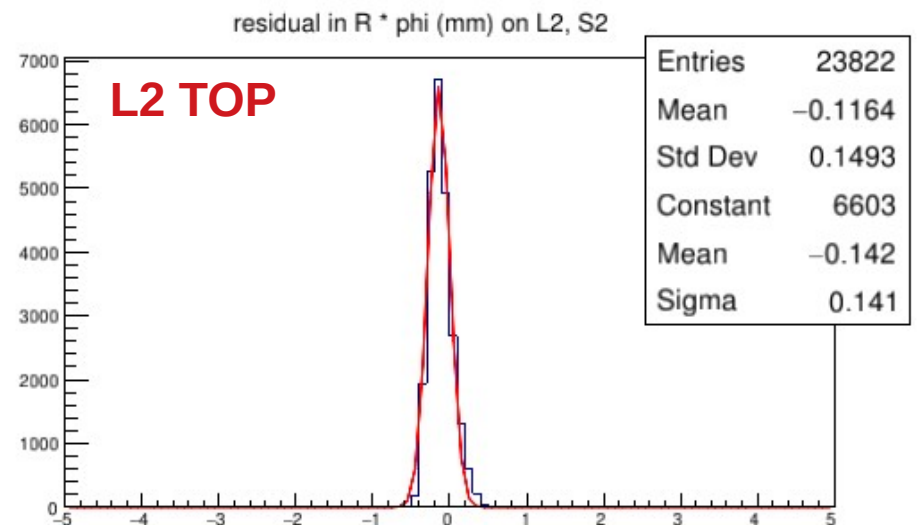
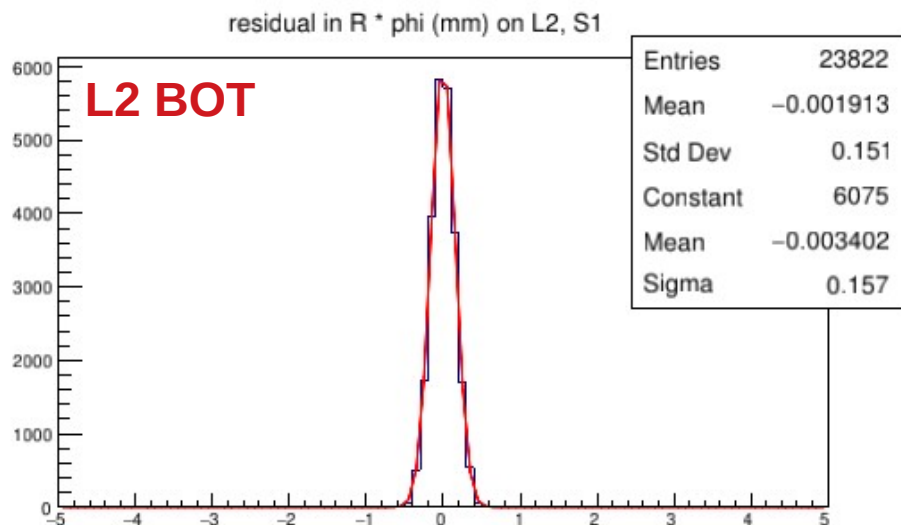
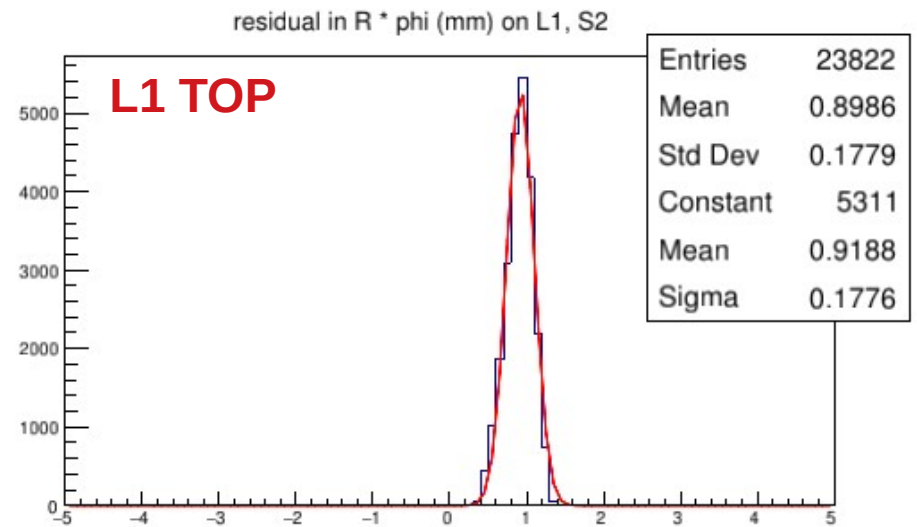
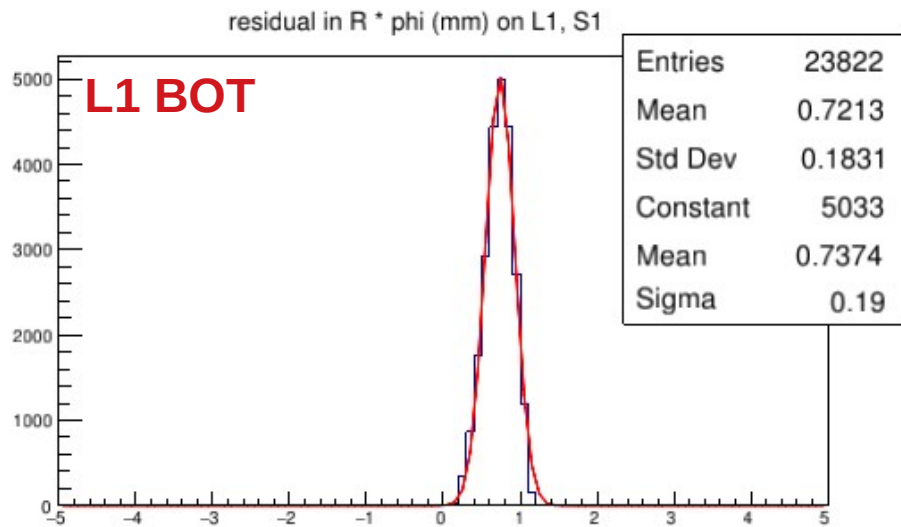
z - LoopAll

- Alignment on
- Chi2 cut = 20
- Energy cut on clusters



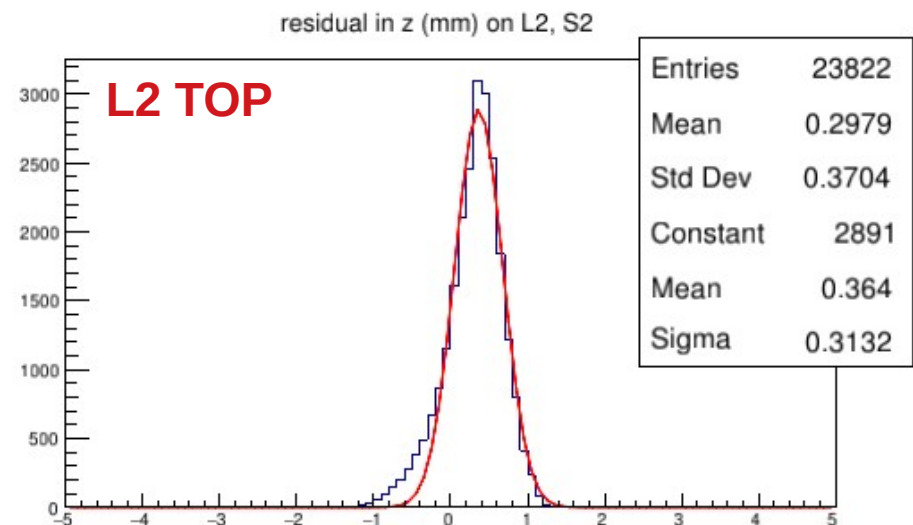
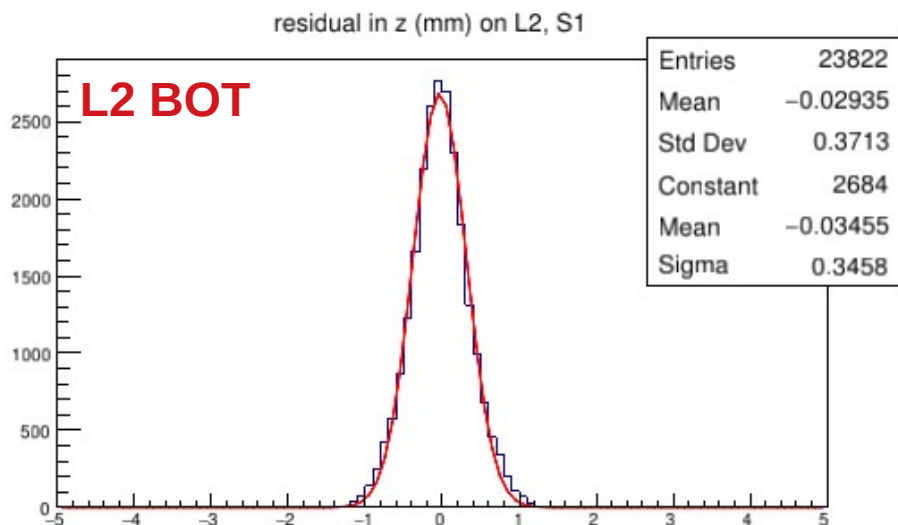
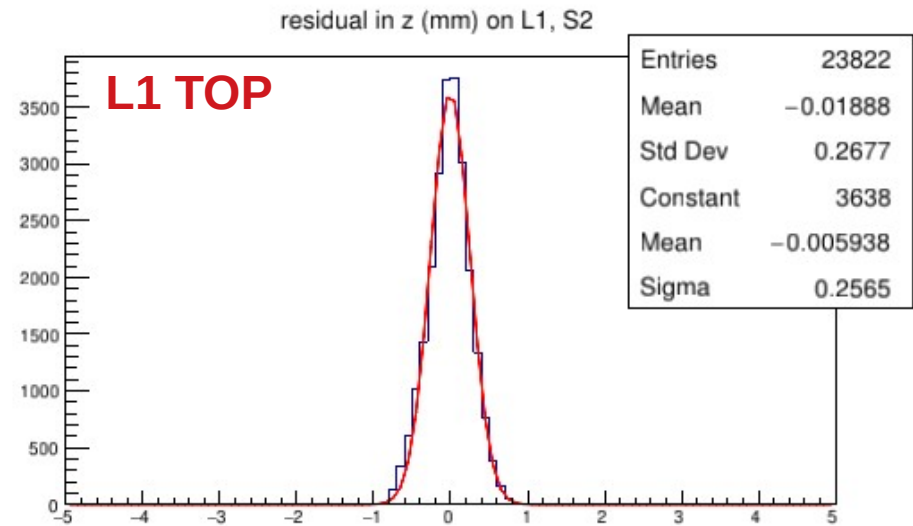
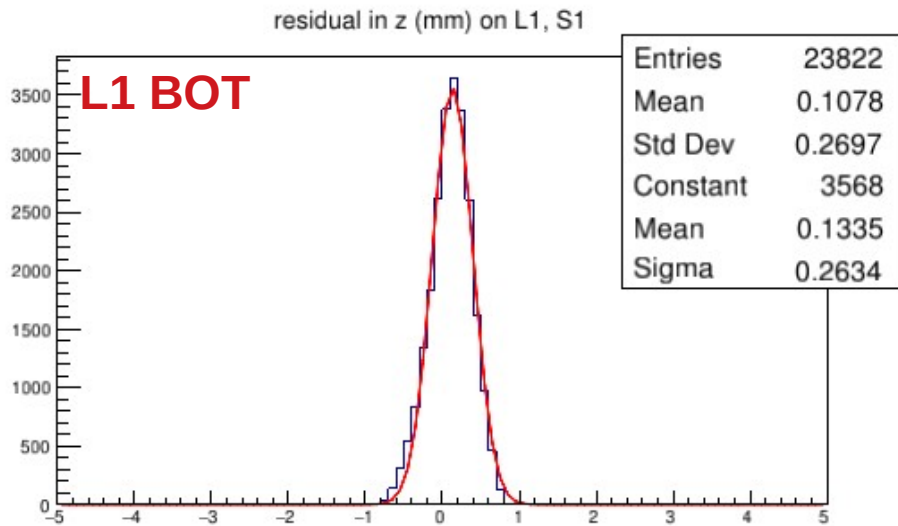
R * phi - LoopAll+MaxQ

- Alignment on
- Chi2 cut = 20
- Energy cut on clusters



z - LoopAll+MaxQ

- Alignment on
- Chi2 cut = 20
- Energy cut on clusters



Conclusions

- uploaded the CgemLineFit-00-00-15 rev. with the fix in CgemCosmicRayQA
- Problem in LoopAll algorithm
- Alignment not taken into consideration properly?
- Can someone independently check the residuals so that we are sure everything is fine please?