

PS tagged gamma energy reconstruction

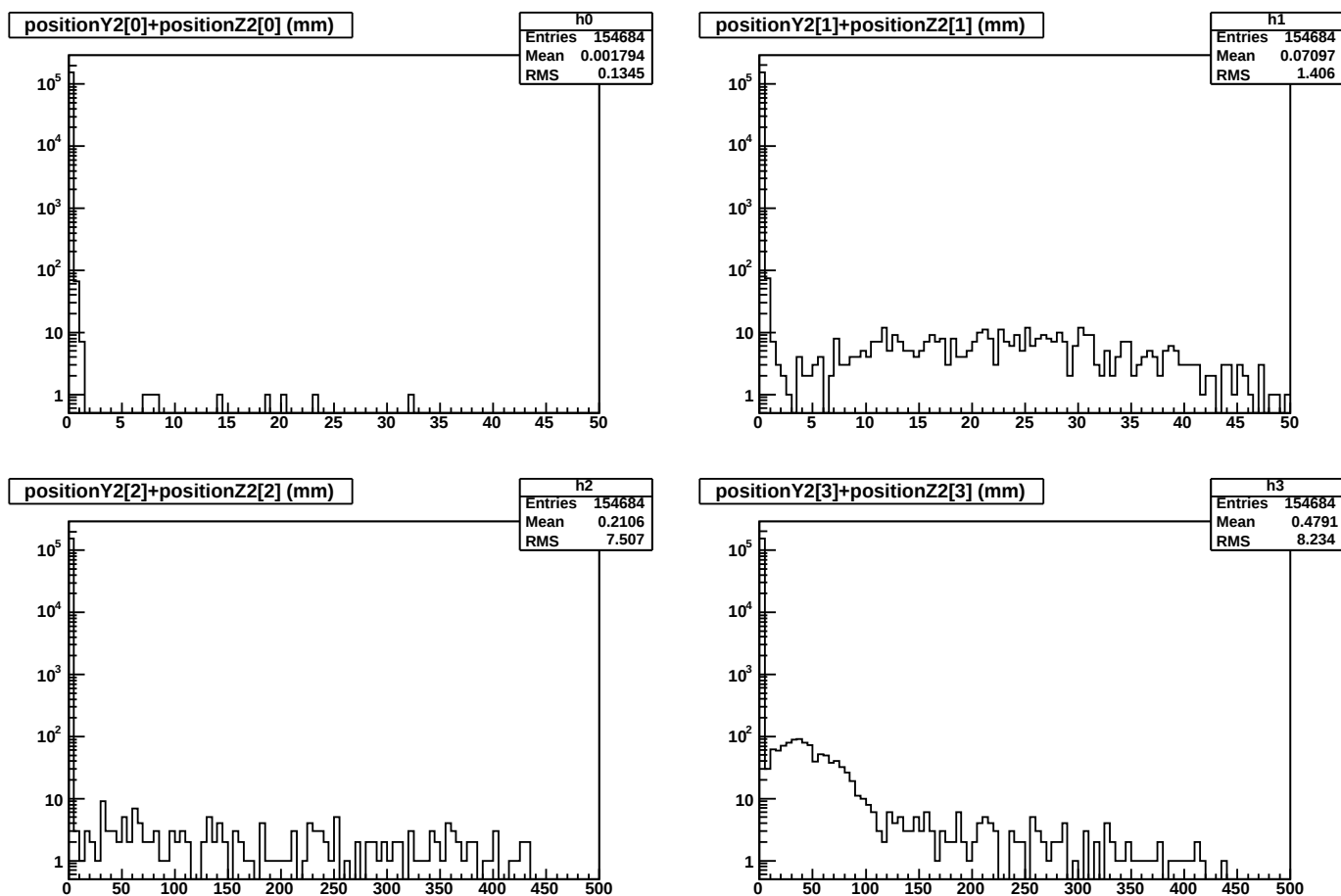
- Quality cuts
- E_{rec} from MPtuple_PS_tag.cpp
- Trajectory fit

Simulation

- BeamtestRelease v1r0801p0
- beamtest06 (April 6th)
- 1 GeV electron beam :
divergence 0.12 deg (hor.) and 0.14 deg (vert)
1% energy dispersion
- modified geometry : bigger Si chambers

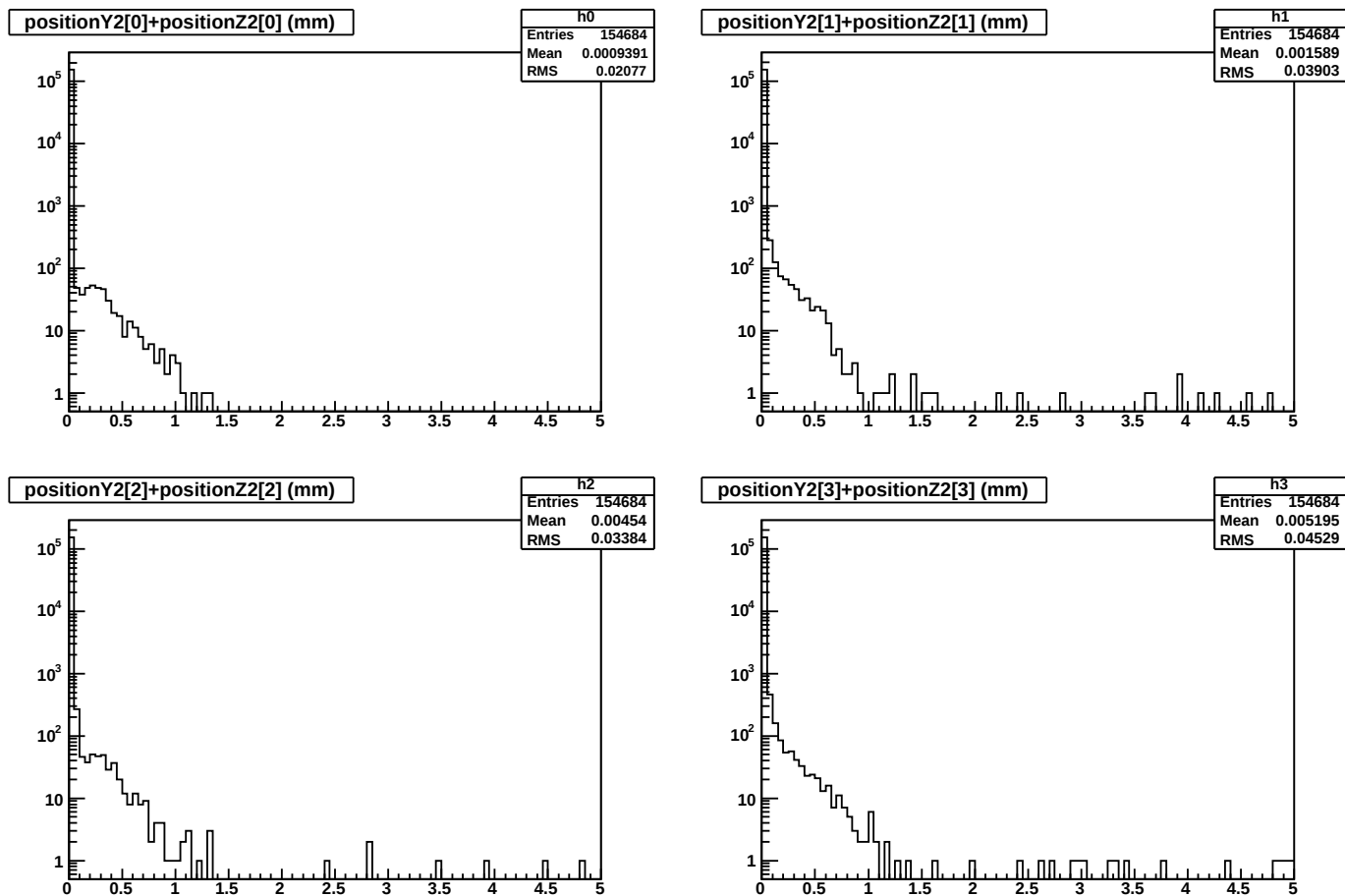
Quality cuts (1)

- there is some energy in all the Si chambers
- looking at the RMS along Y and Z



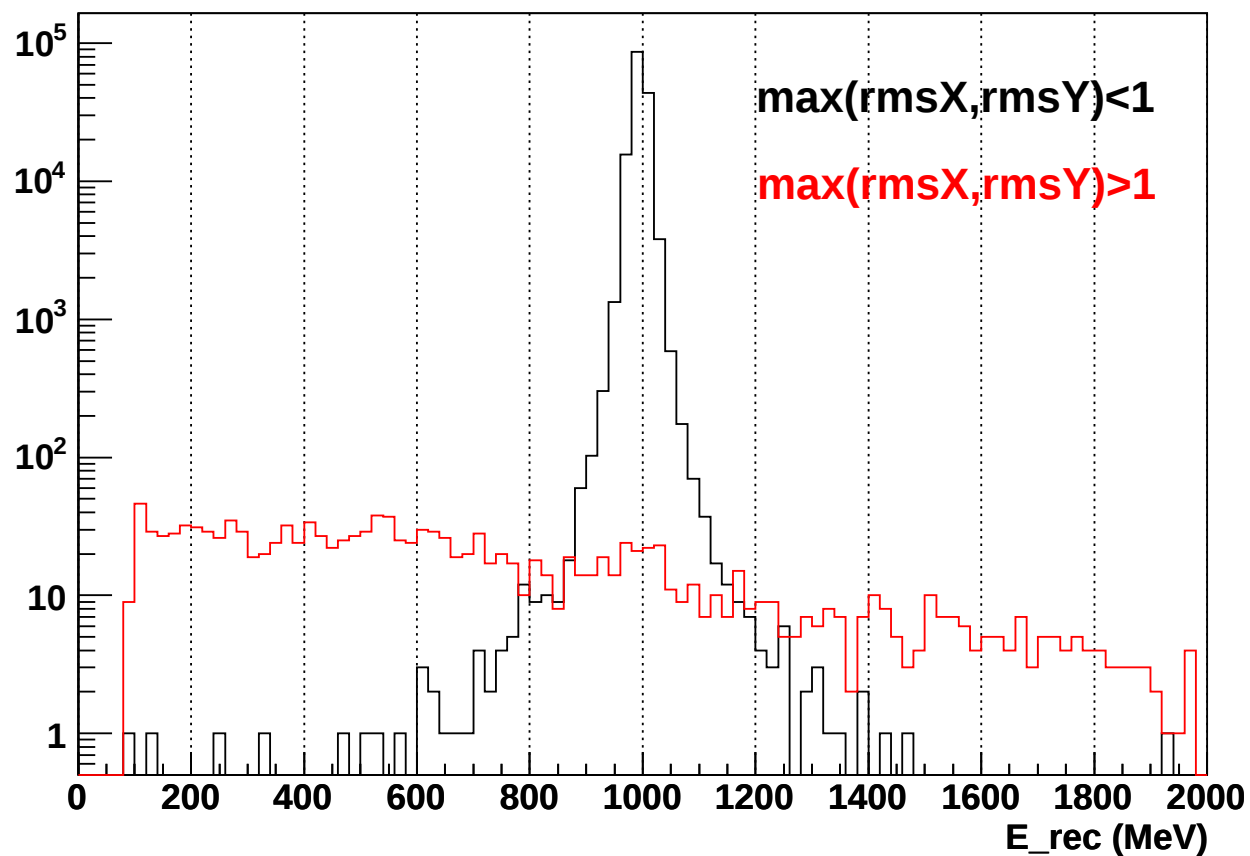
Quality cuts (2)

- zoom between 0 and 5 mm



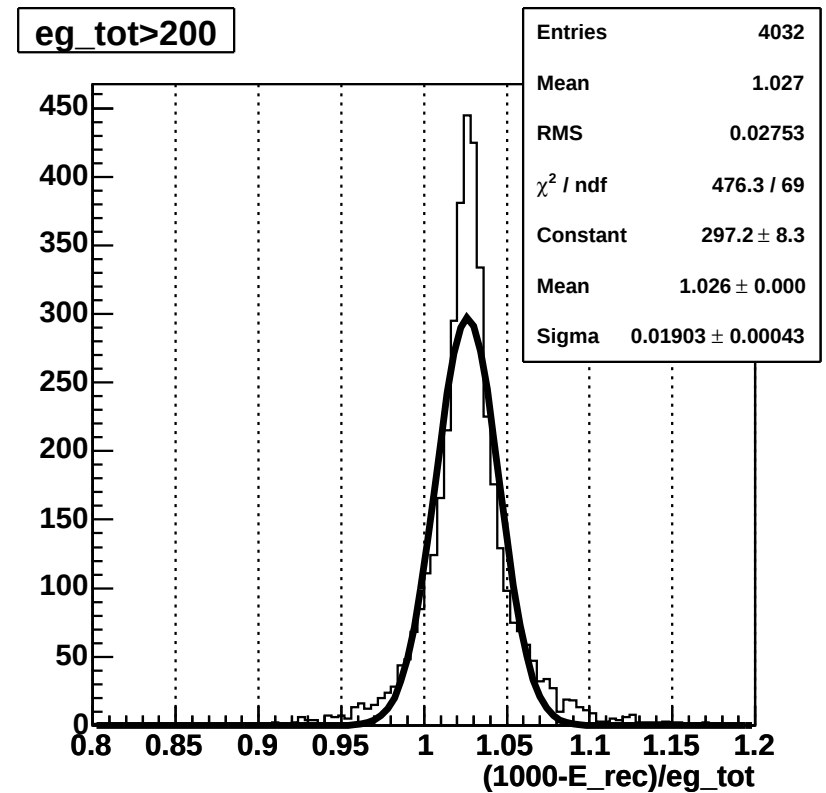
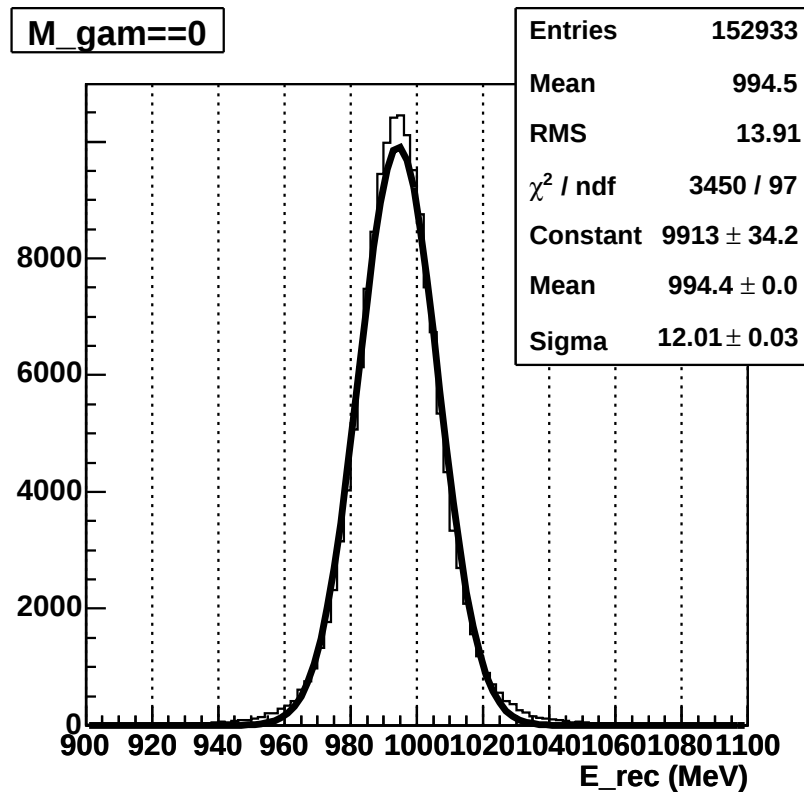
Quality cuts (3)

- requiring that $\max_{i=0,4}(rms_X^i + rms_Y^i) < 1$ mm to get rid of delta ray e^- (and dump backsplash ?)



E_{rec} results

- $E_{rec}(\text{MeV}) = 300 \times 1 \text{ T} \times 0.5 \text{ m} / (2 \tan(\theta/2))$
- -0.5% bias for E_{rec}
- $+2.5\%$ for $E_{gam} = 1000 - E_{rec}$

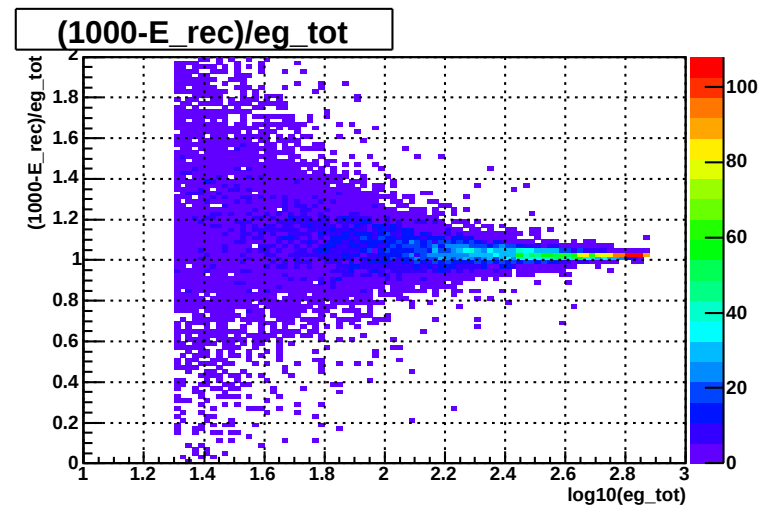
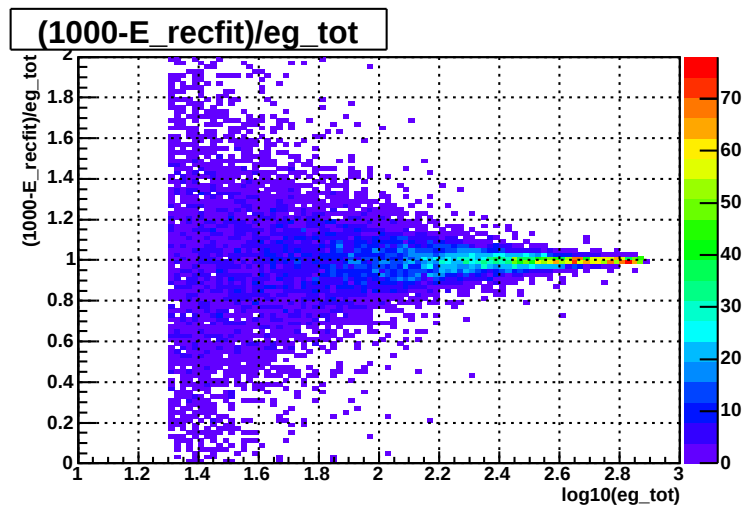
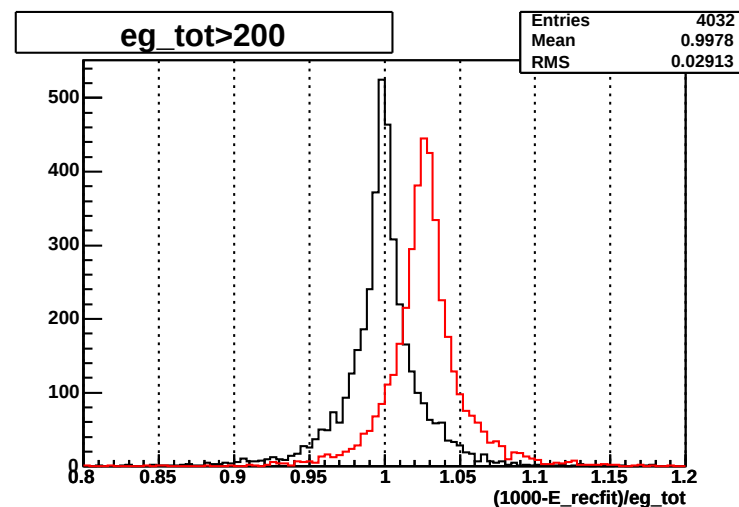
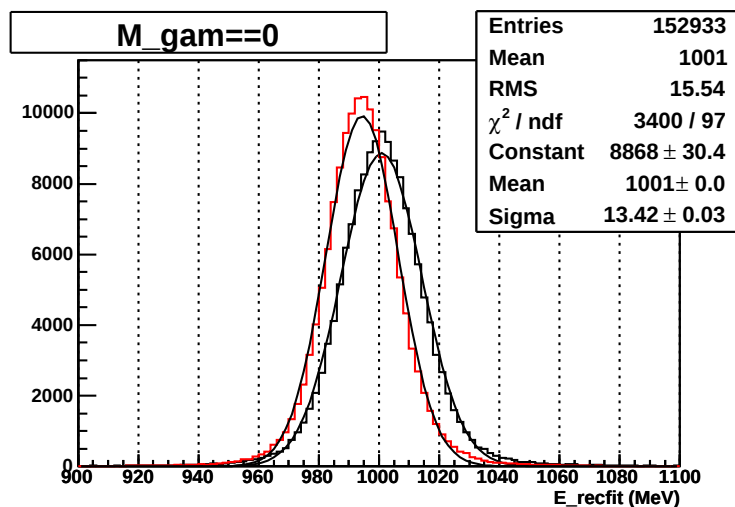


simple χ^2 minimization

- retrieve the positions in the 2 first Si chambers
- determine the position when entering the magnet
- compute the trajectory inside the magnet
- determine the positions in the 2 last Si chambers : $P2_{fit}(E)$ and $P3_{fit}(E)$
- minimize
$$\chi^2 = dist^2(P2, P2_{fit}(E)) + dist^2(P3, P3_{fit}(E))$$

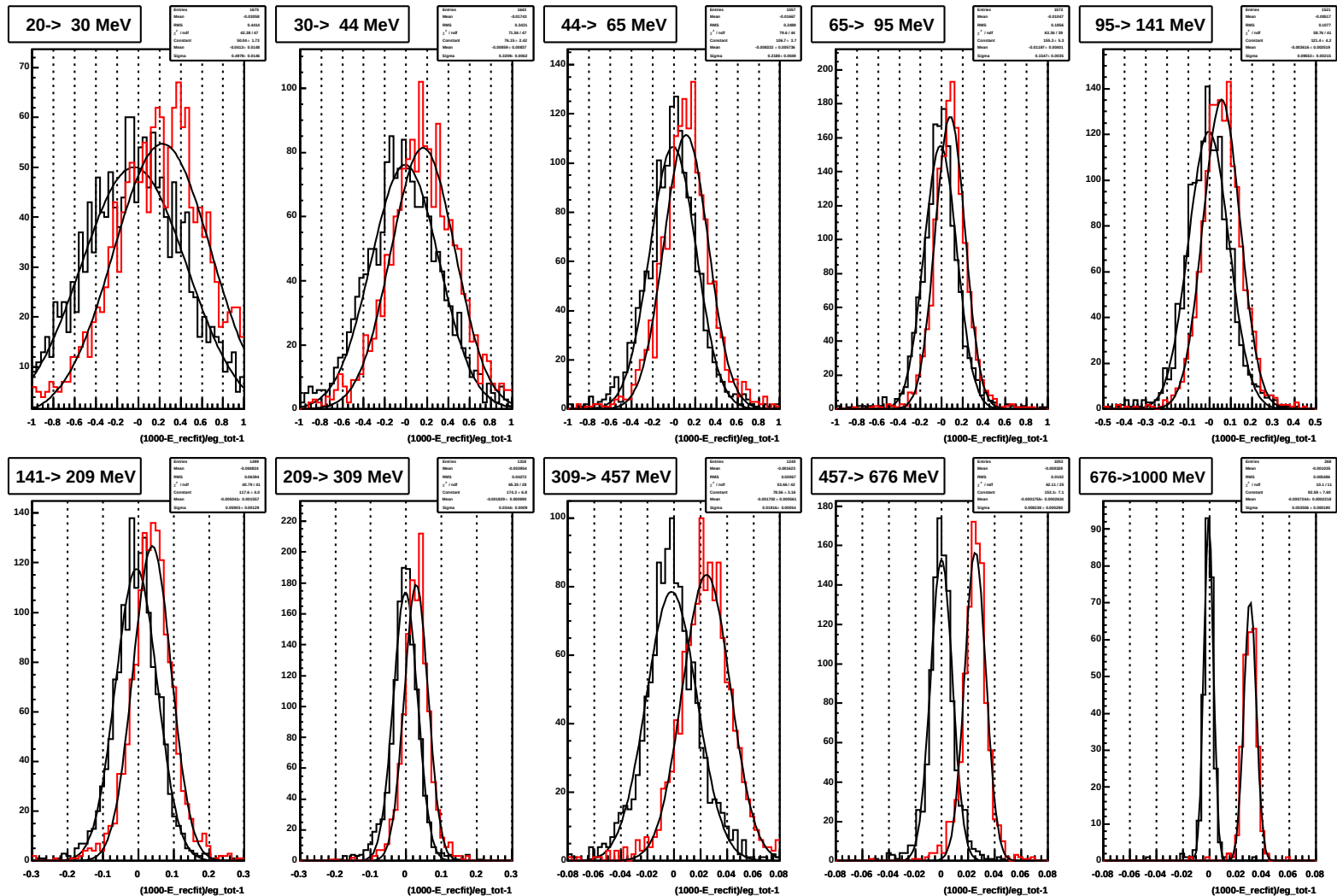
E_{recfit} results

● E_{recfit} (in black) : no bias anymore



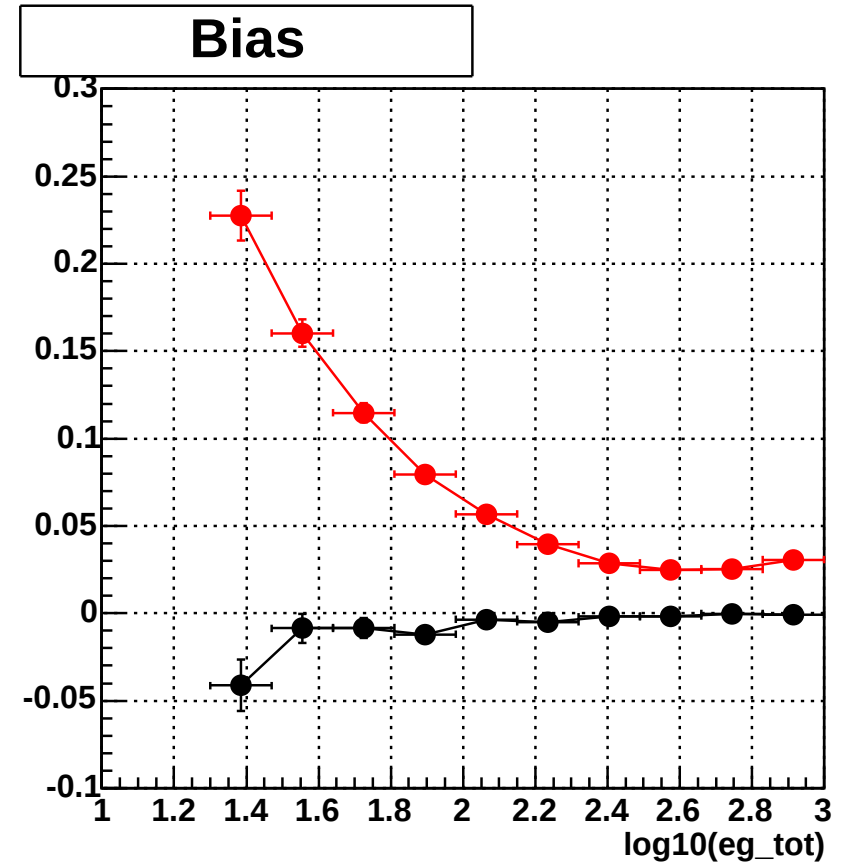
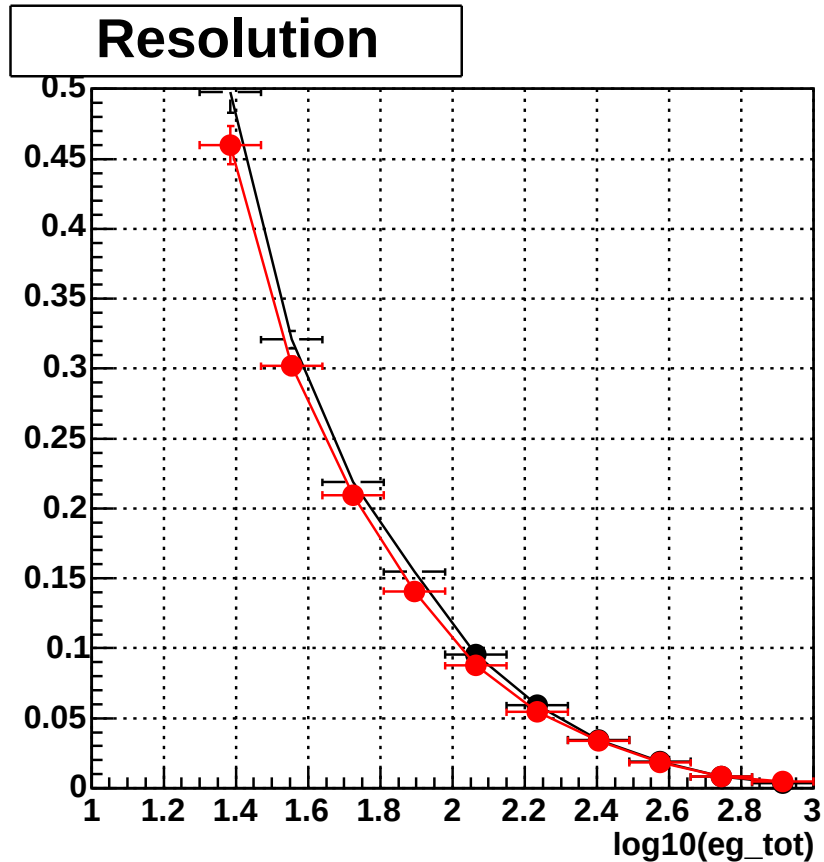
Tagged gamma energy reconstruction

● from E_{recfit} (black) and E_{rec} (red)



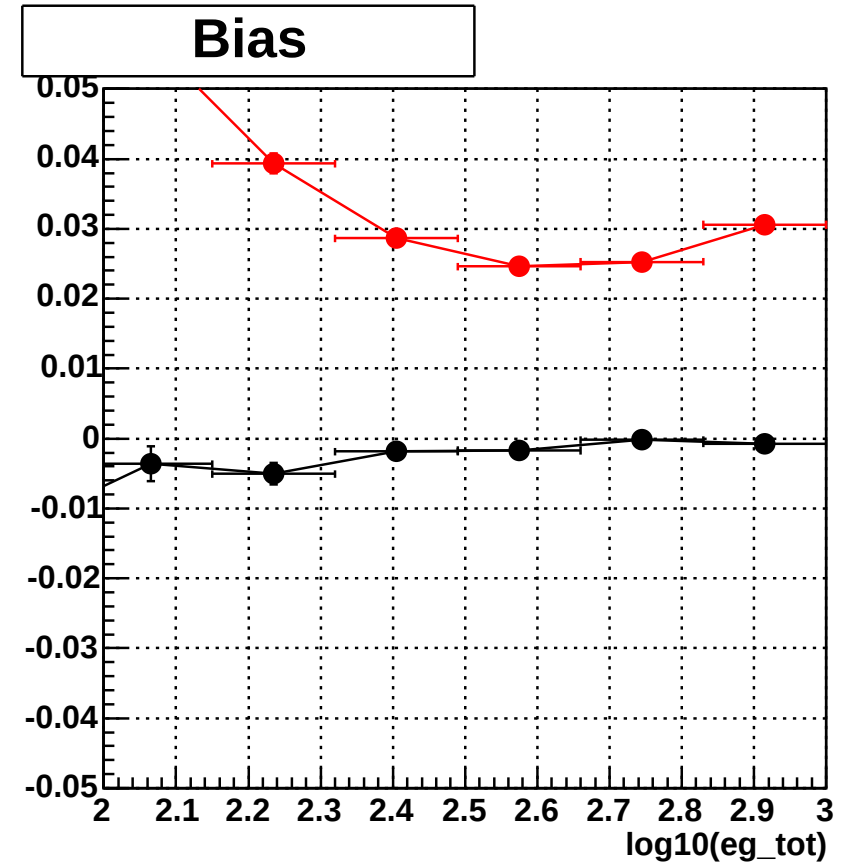
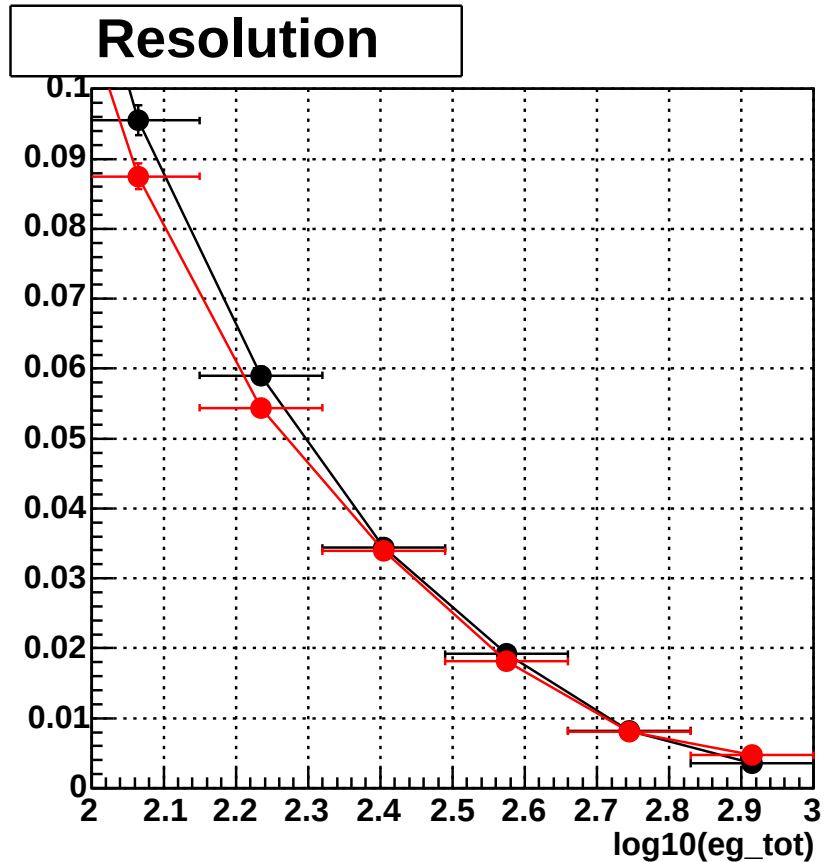
Tagged gamma energy resolution

● from E_{recfit} (black) and E_{rec} (red)



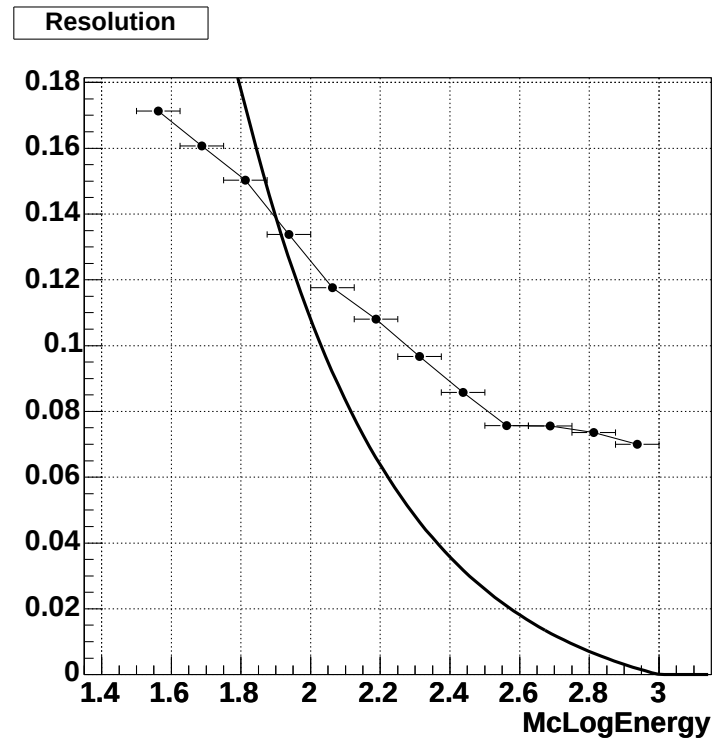
Tagged gamma energy resolution

● zoom above 100 MeV



New energy limitation at PS

- on-axis CU energy resolution (points) and tagged gamma energy resolution (function)
- with a 1 GeV electron beam : it is possible to check the CU energy resolution down to 150 MeV



Preliminary today's news

- I have added the Si chamber containers (0.012 mm Al) :
no change
- I have just discovered that the simulation was made in
vacuum... :
within air, the resolution for unradiating electrons is 1.5%
instead of 1.2%
- next ? :
try to improve the fit taking into account multiple scattering
- question :
is it realistic ? (accuracy of the position of the Si chambers
in the experimental hall, etc...)