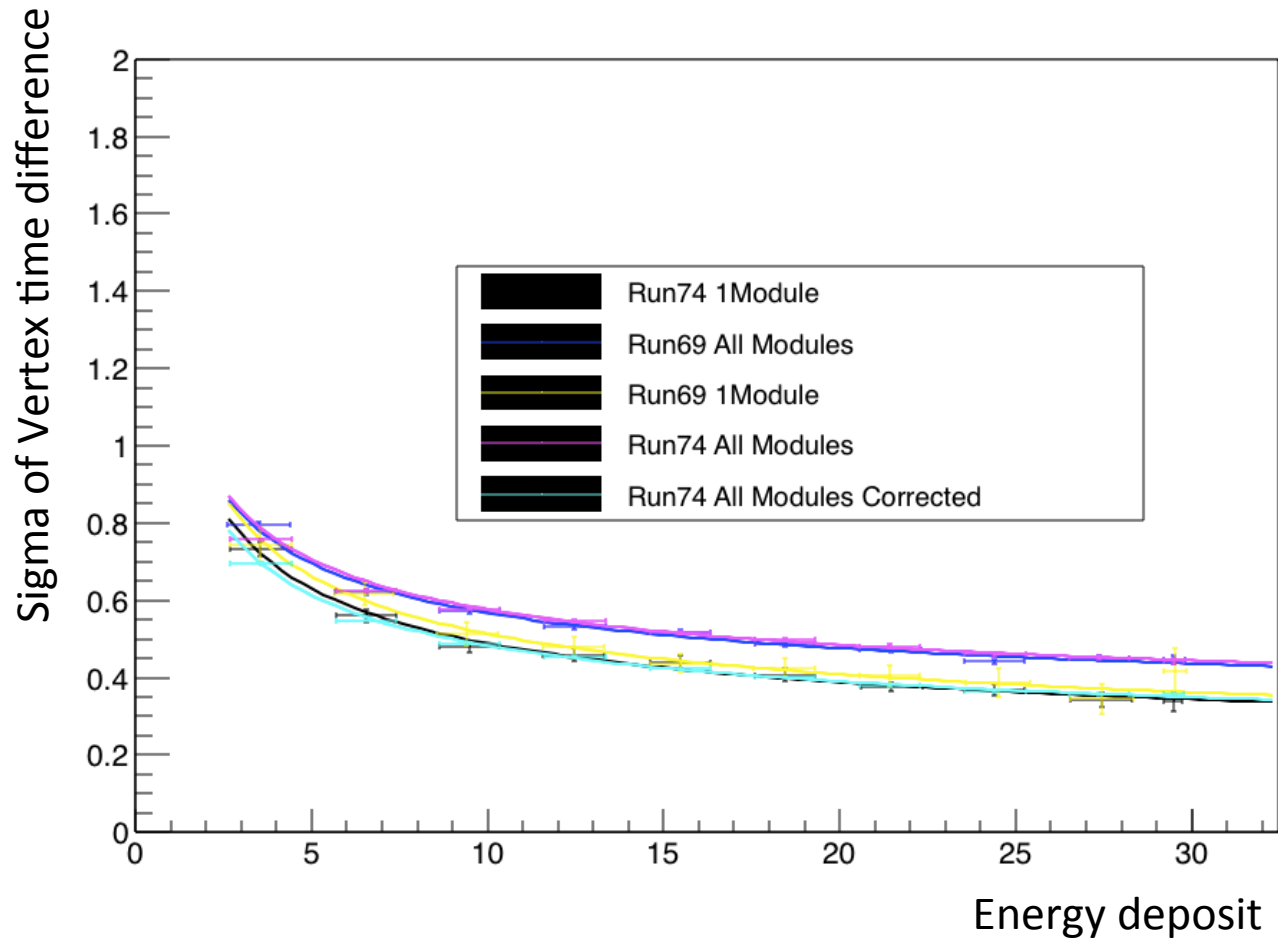


Report_170517

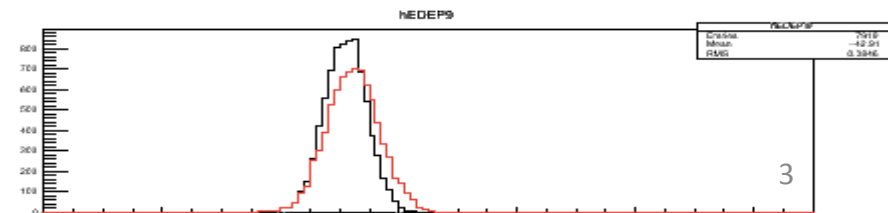
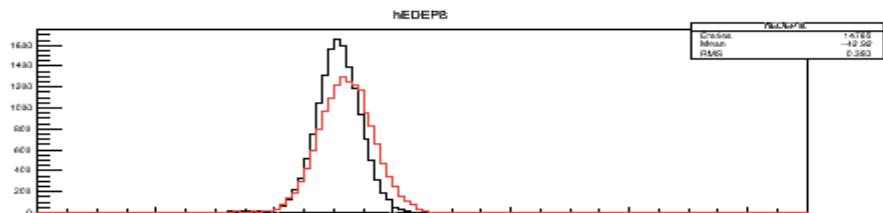
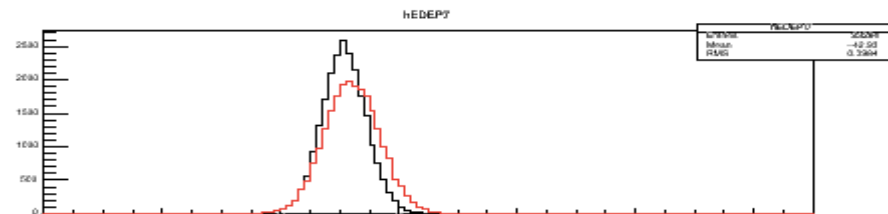
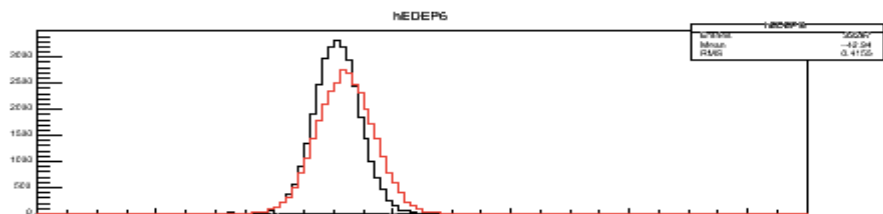
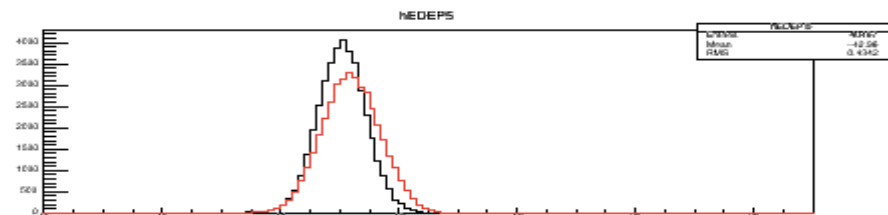
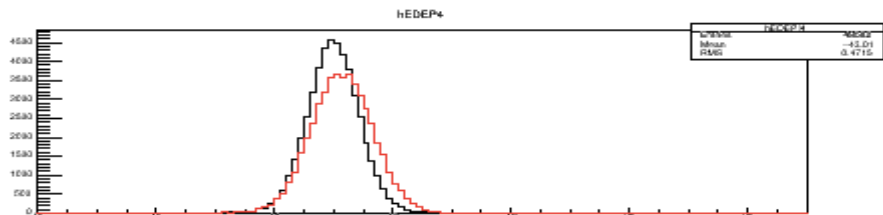
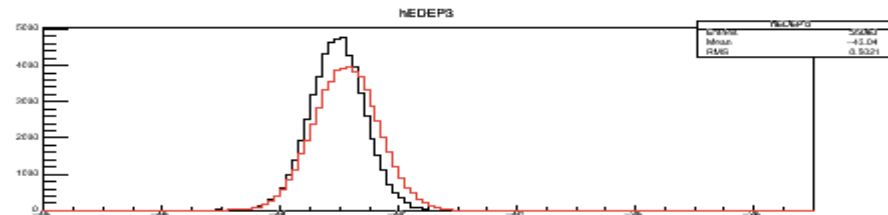
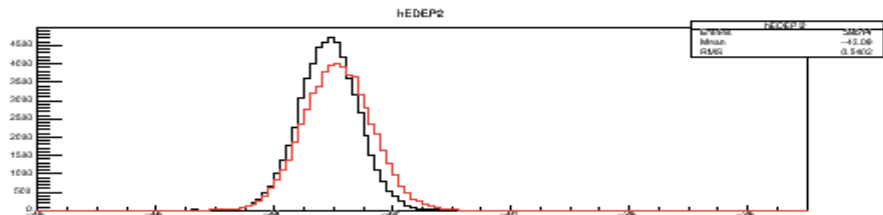
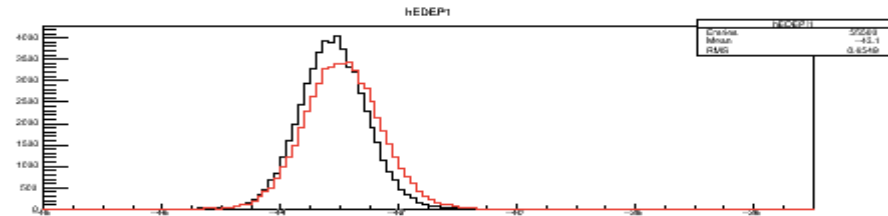
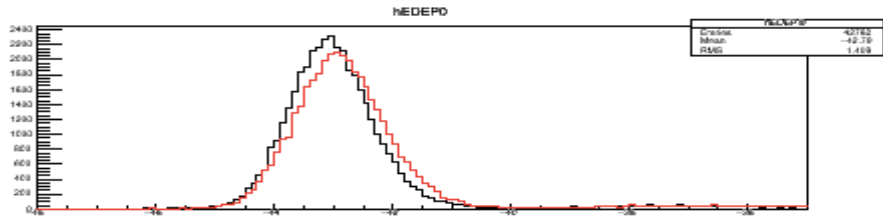
Timing Resolution

- Module Correction check

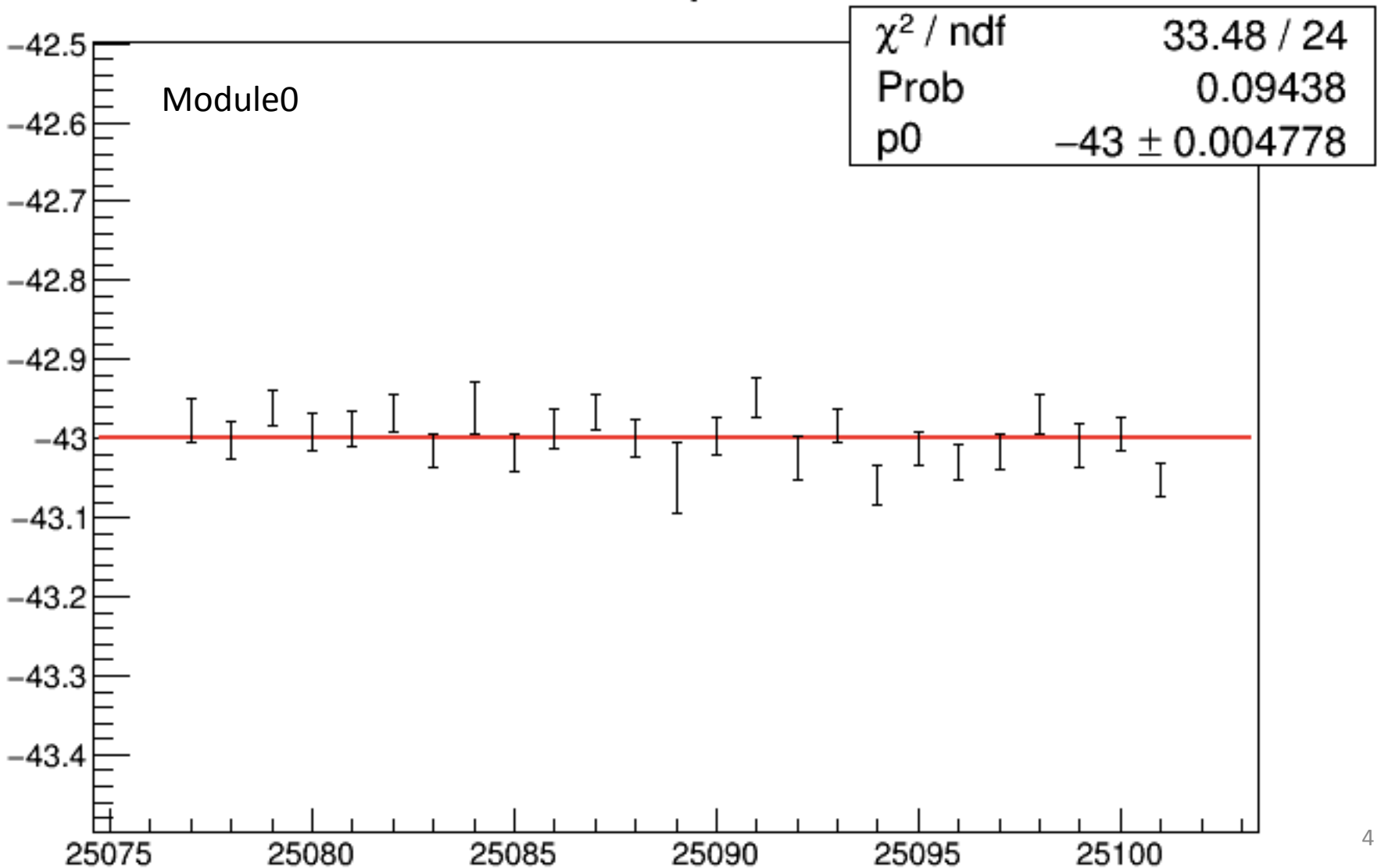


Raw Histograms

Black : After Correction (Run74)
Red : Before Correction (Run74)



Module stability



Sampling Fraction simulation

- Change thickness of IMB lead sheet to 0.9mm, 0.8mm respectively. (From 1.0mm)
- Check IMB && OMB status

Radiation length check

- Data Selection
 - IMB Hit && OMB Hit
- Check Ratio of Energy deposited

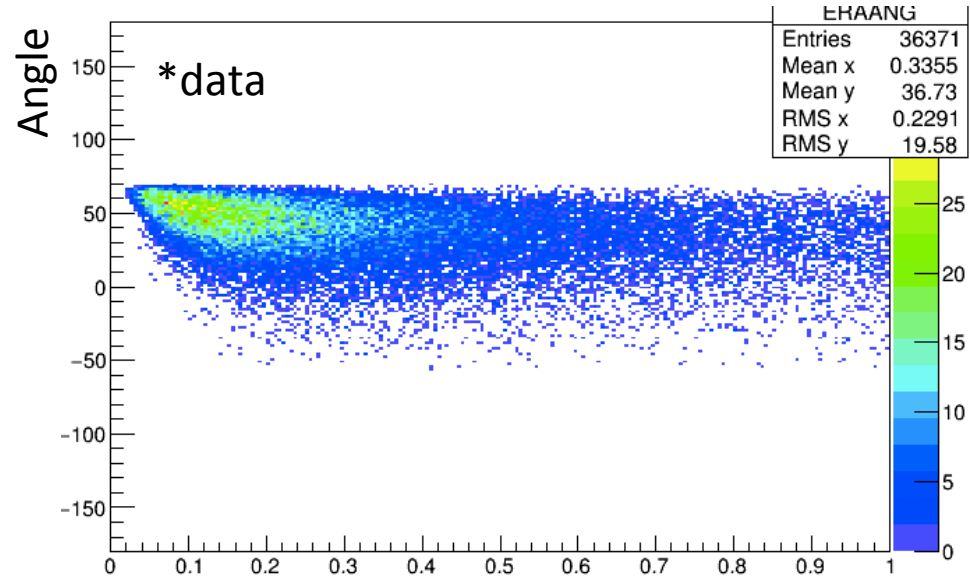
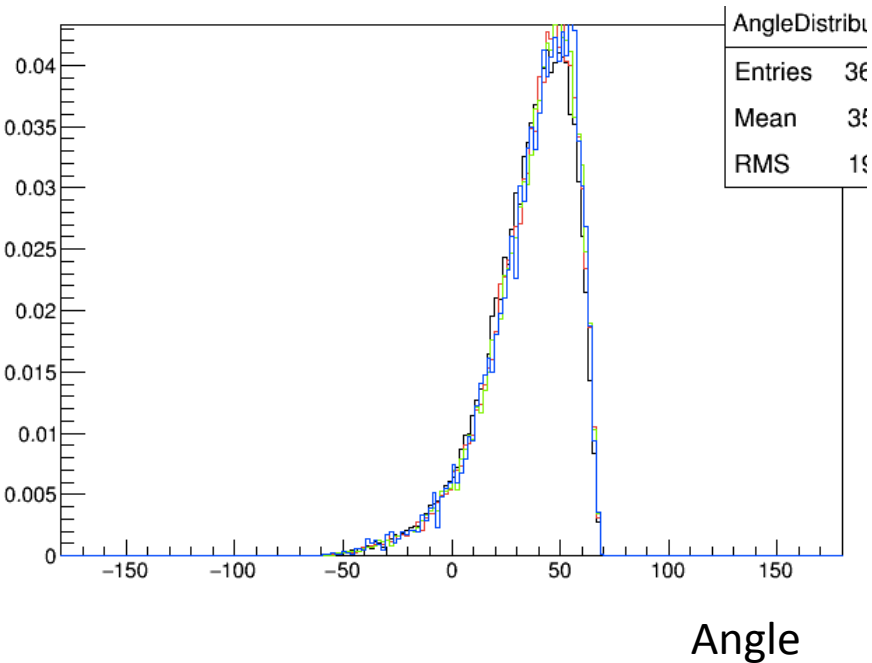
$$\text{Ratio} = \frac{E_{\text{OMB}}}{E_{\text{IMB}}} = \frac{\text{SAMP}_{\text{OUT}}}{\text{SAMP}_{\text{IN}}} \times \frac{E_{\text{INCE}}(2^{-(X-x)/X})}{E_{\text{INCE}}(1-2^{-(X-x)/X})}$$

Capital X : Radiation Length
Small x : starting point of interaction

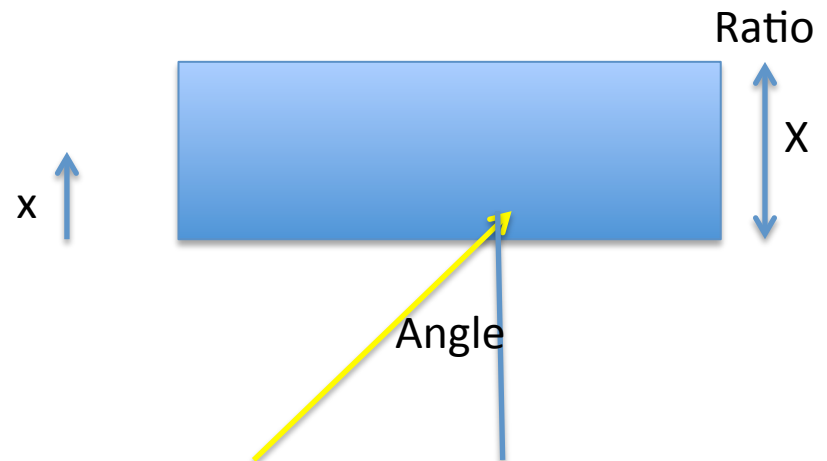
Full deposition is assumed

Distribution of x : exponential with slope X
X is related to thickness of Lead sheet

Angle dependency



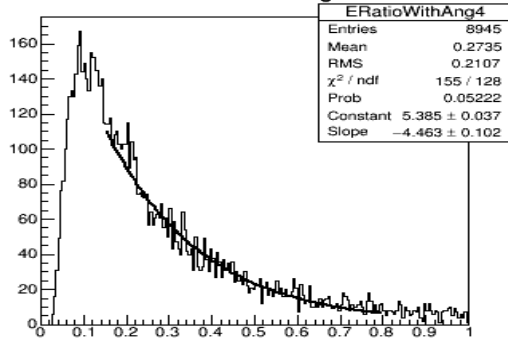
Black : data
 Red : Orig MC
 Green : 0.9mm MC
 Blue : 0.8mm MC



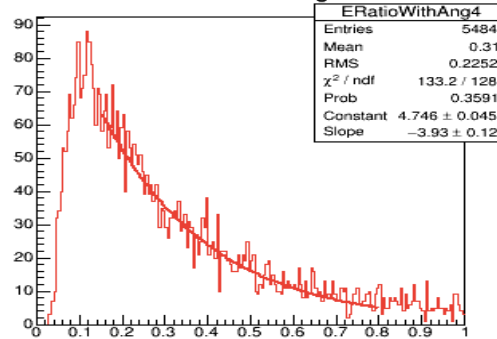
Ratio

With angle selection ($48 < \theta < 60$)

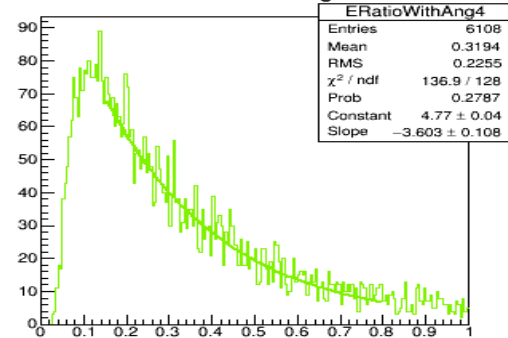
ERatioWithAng4



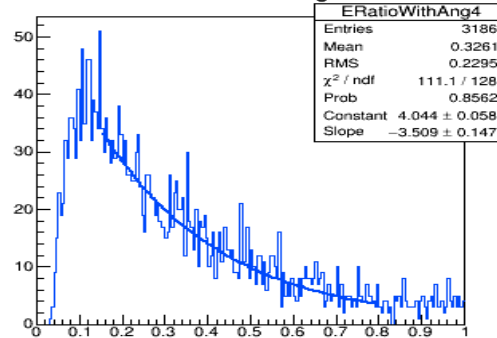
ERatioWithAng4



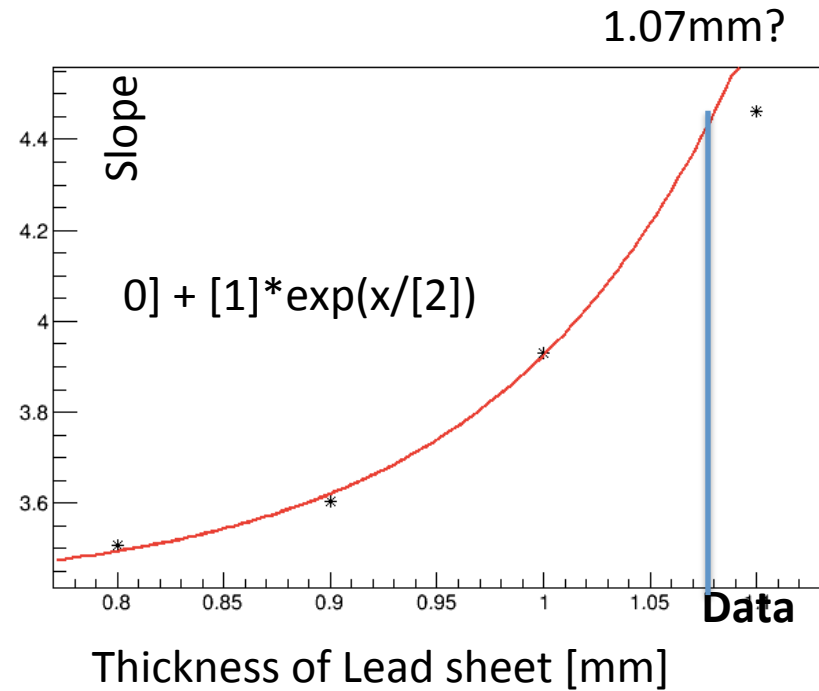
ERatioWithAng4



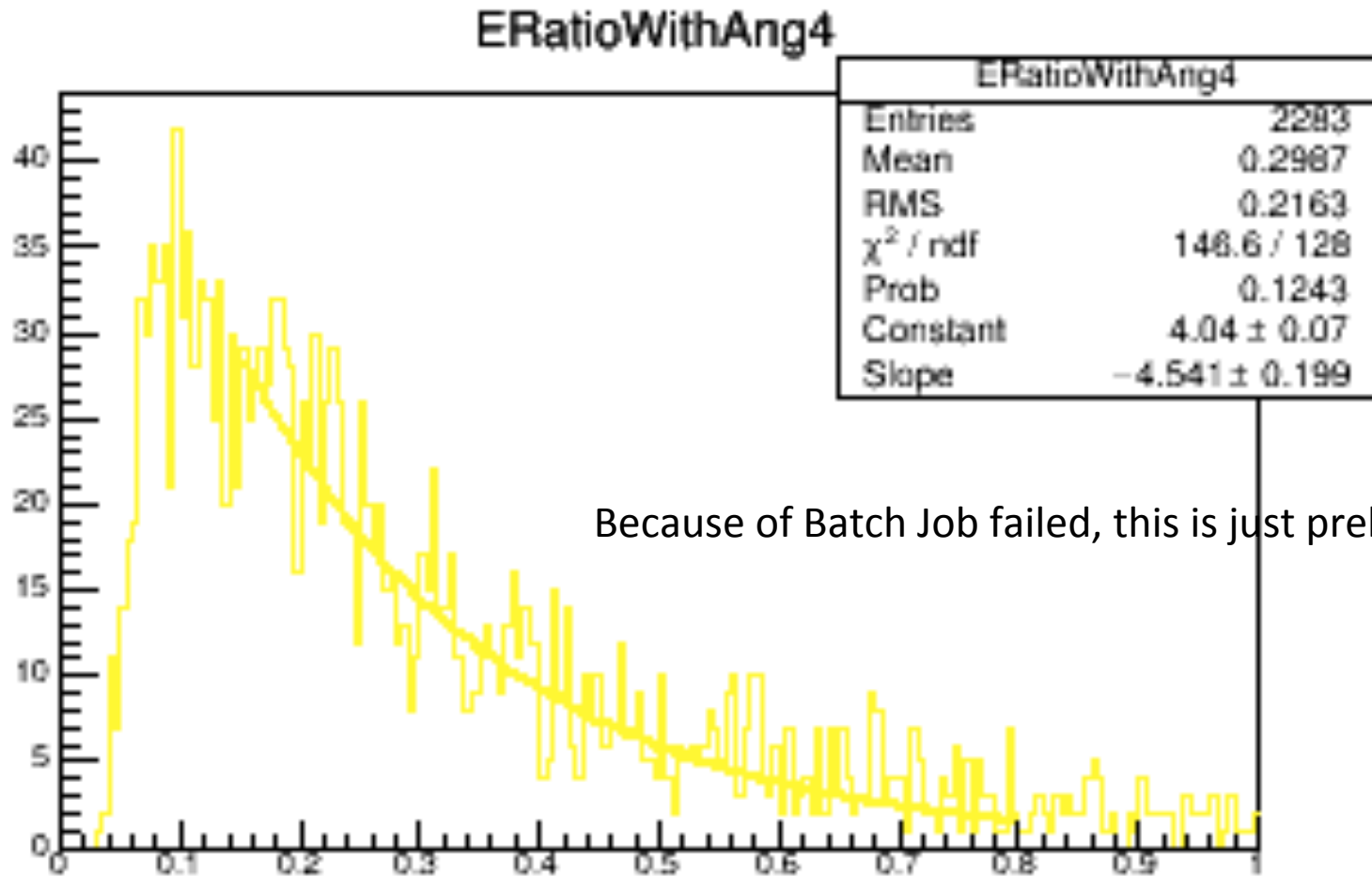
ERatioWithAng4



Exponential Fitting

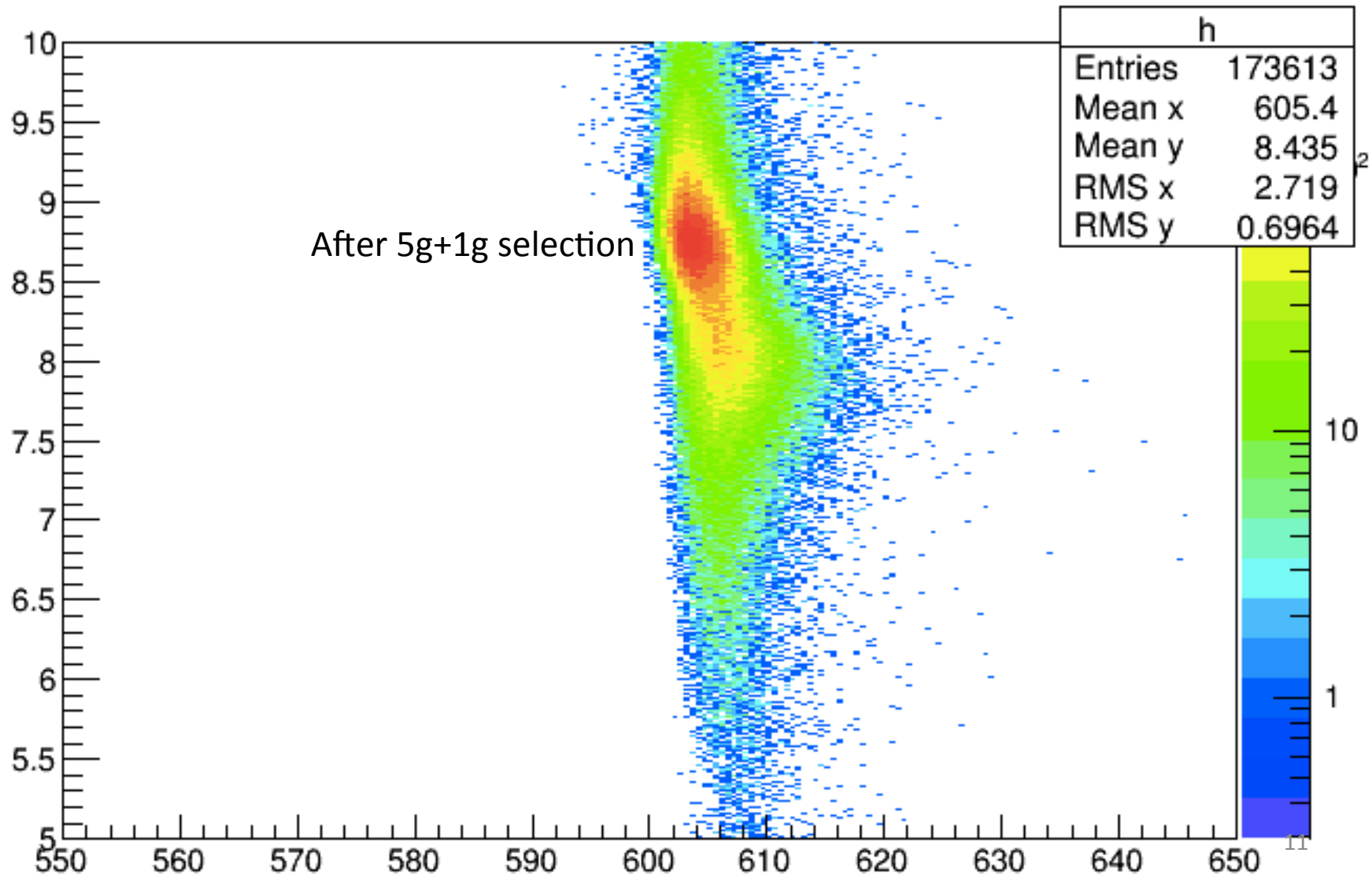


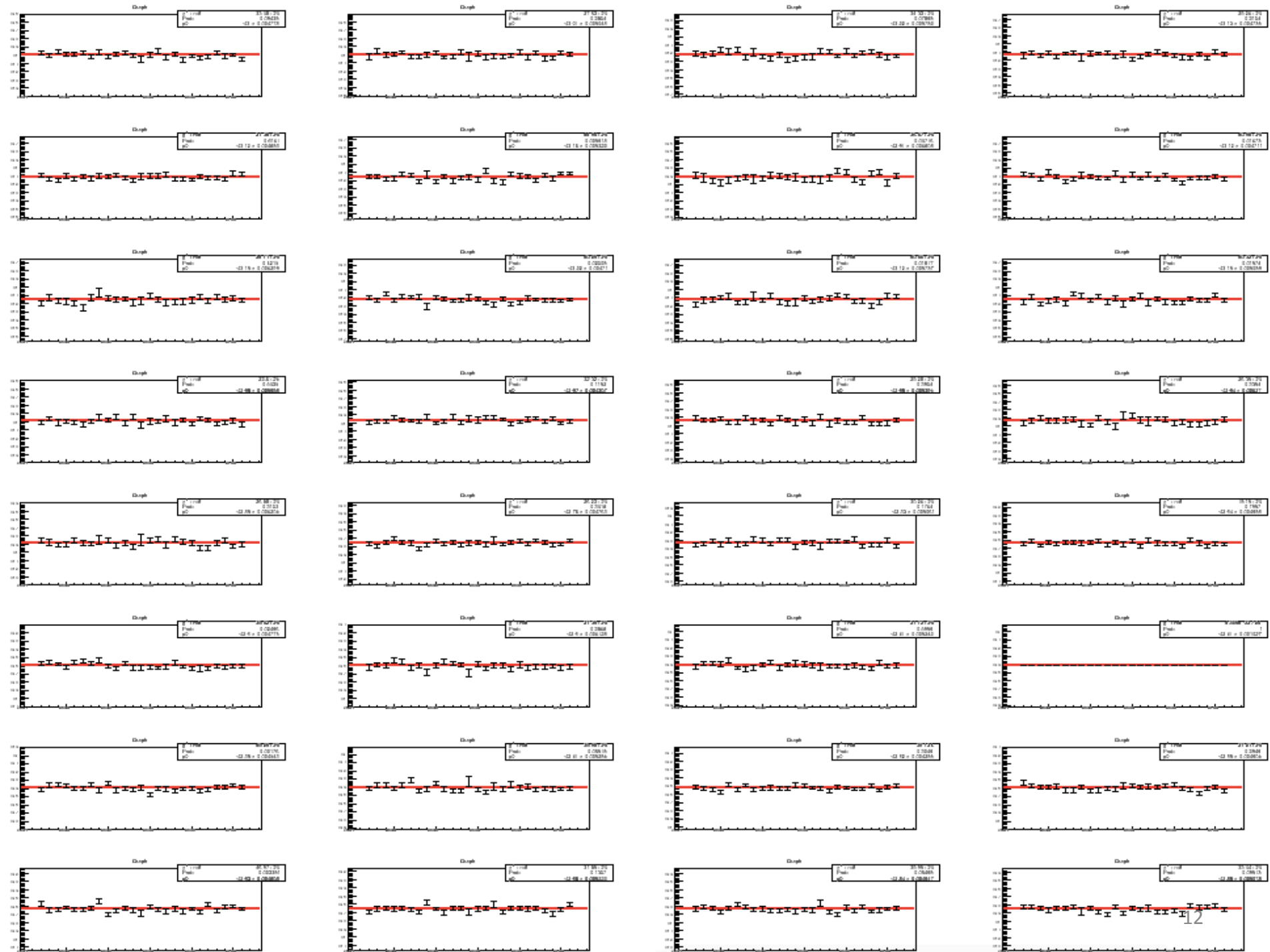
1.1mm case



Raw Distributions

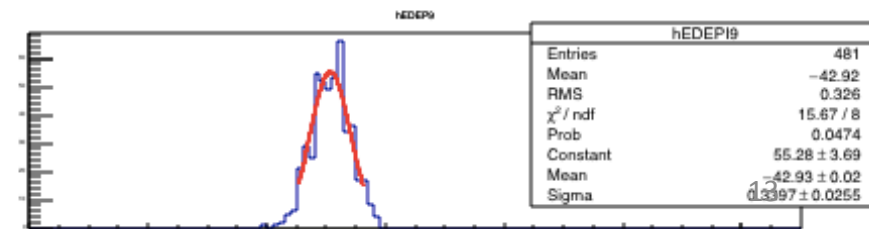
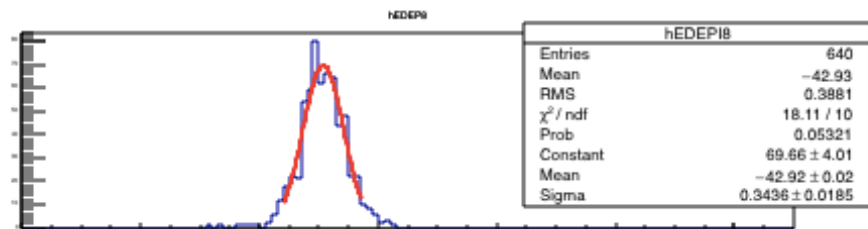
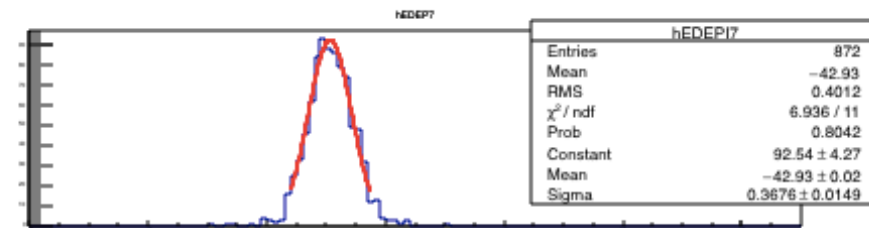
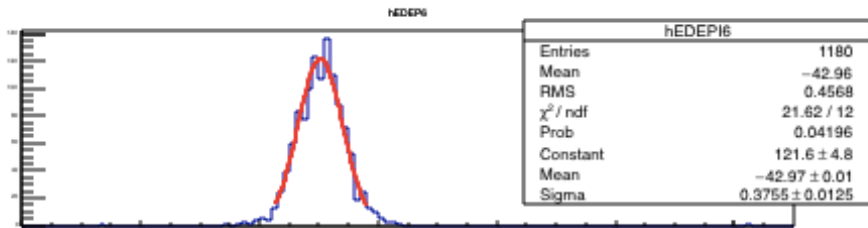
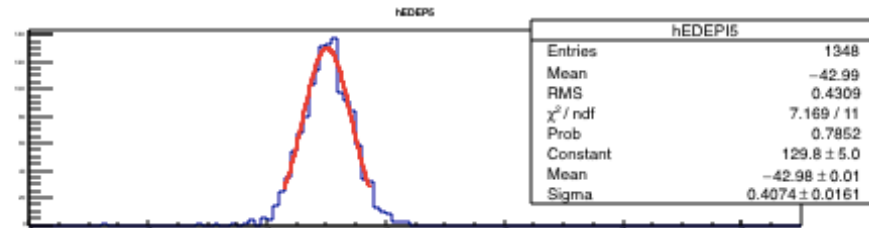
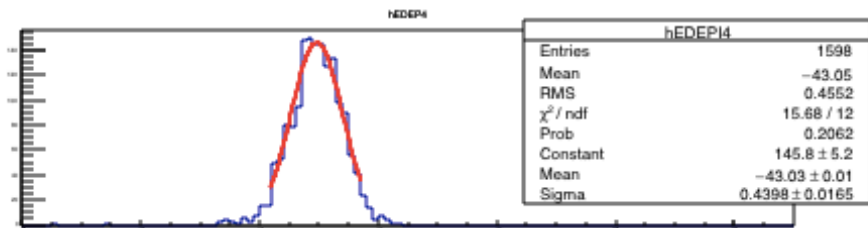
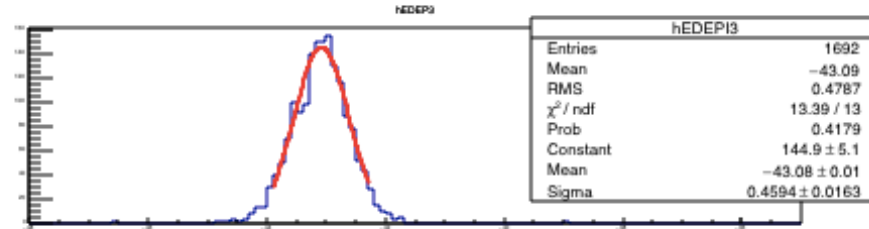
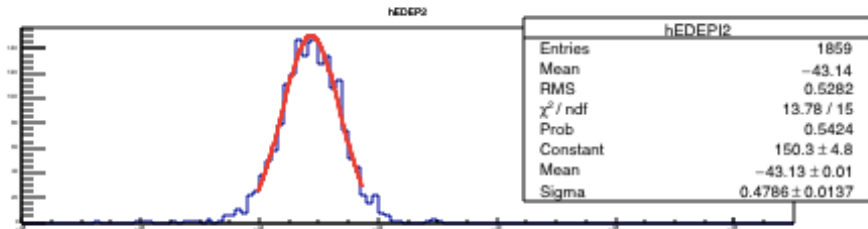
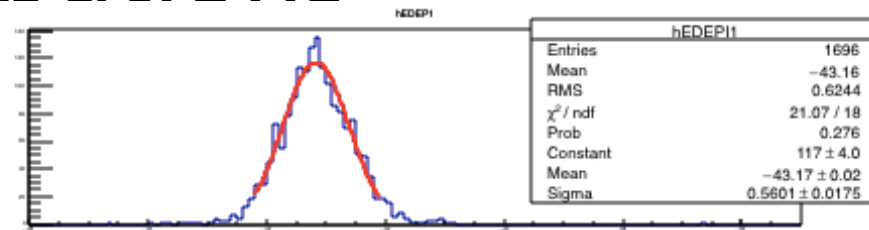
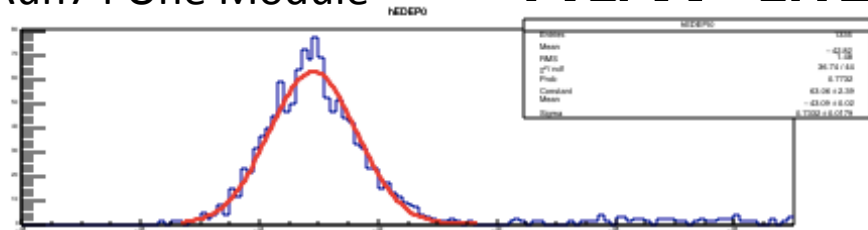
Pedestal vs Pulse Stability





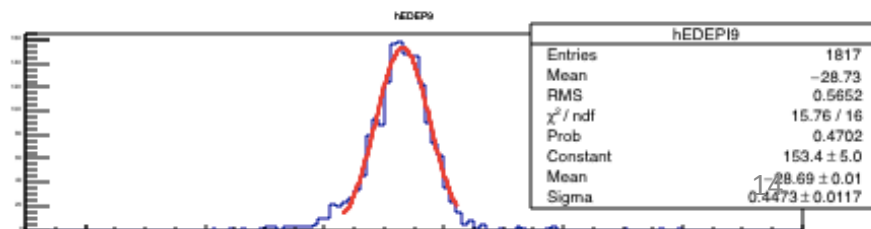
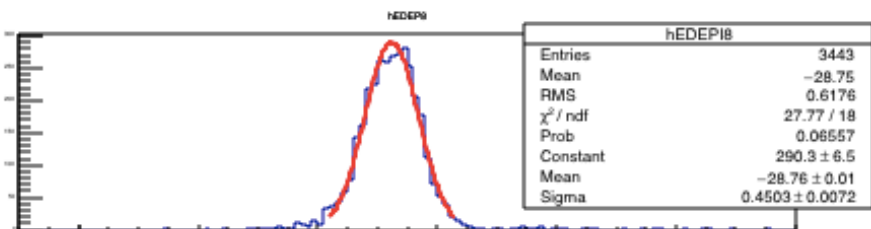
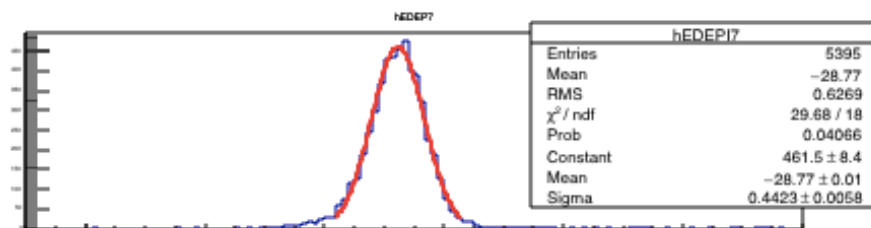
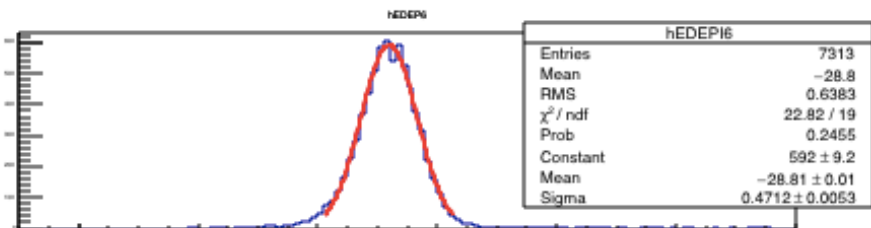
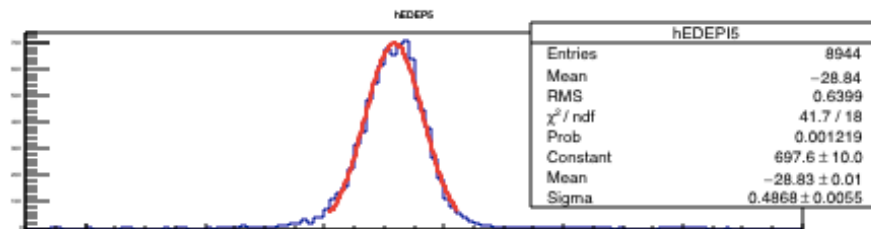
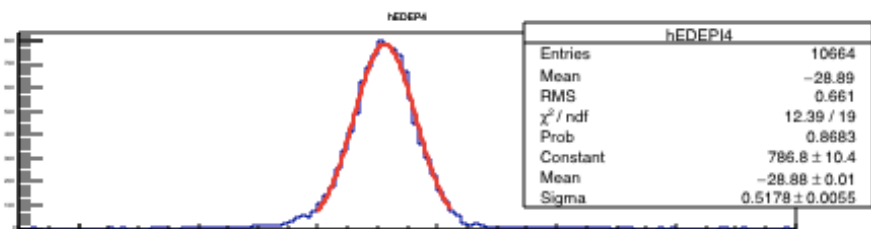
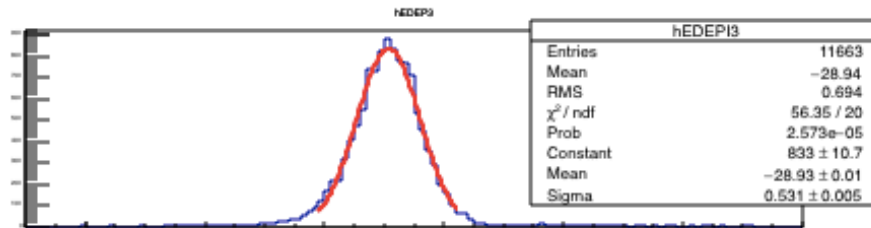
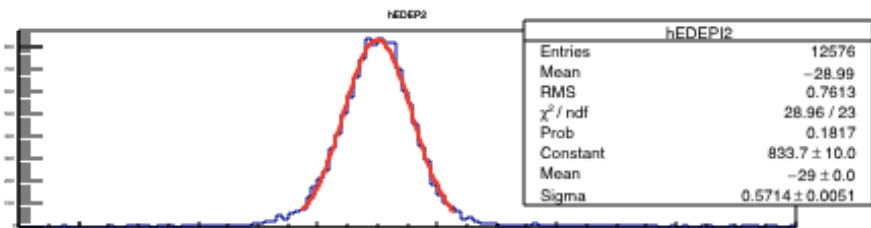
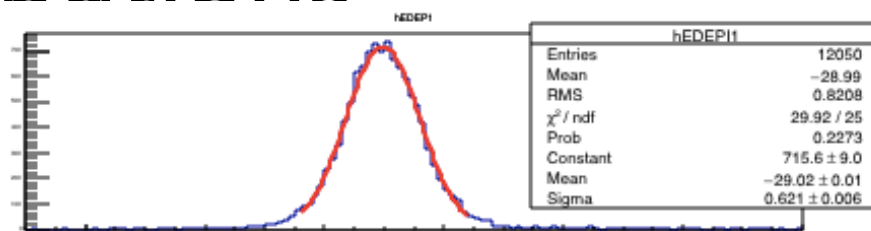
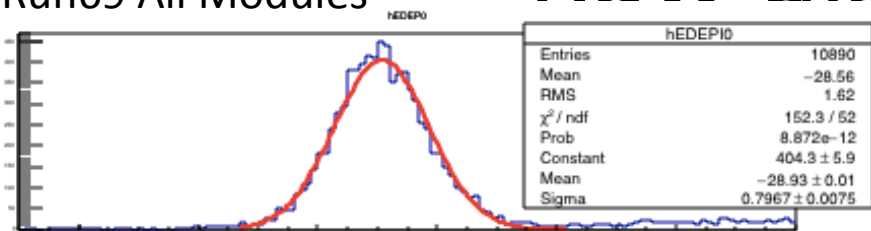
Raw distributions

Run74 One Module



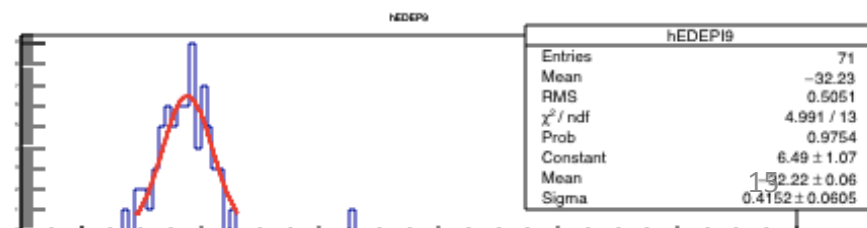
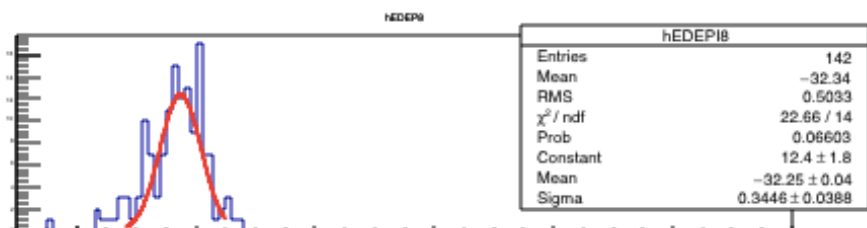
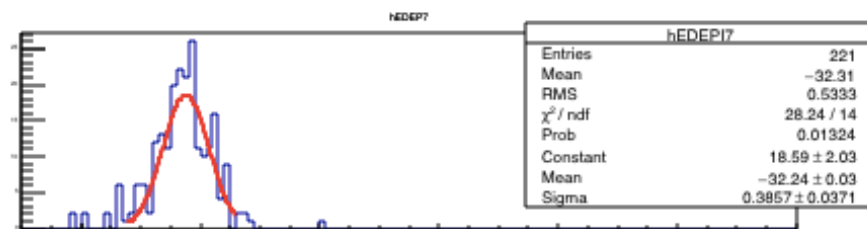
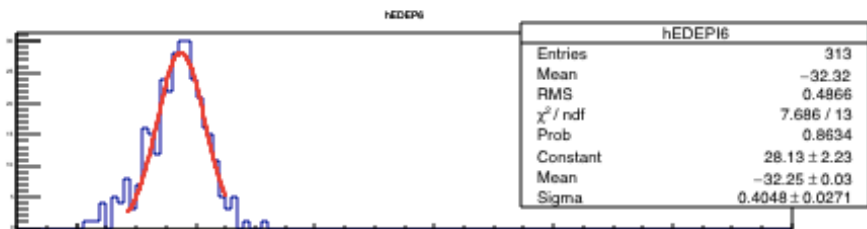
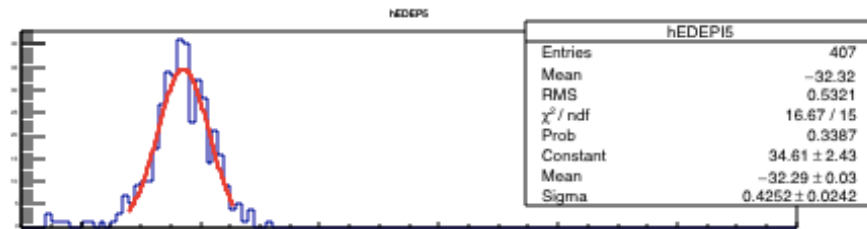
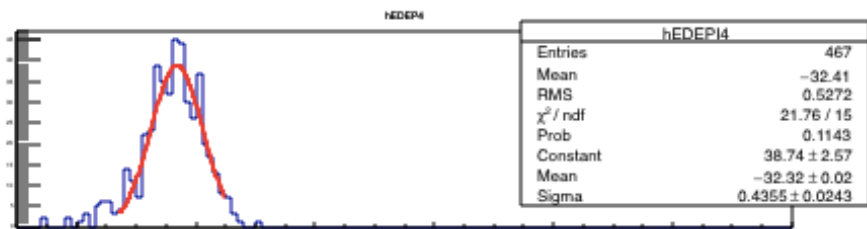
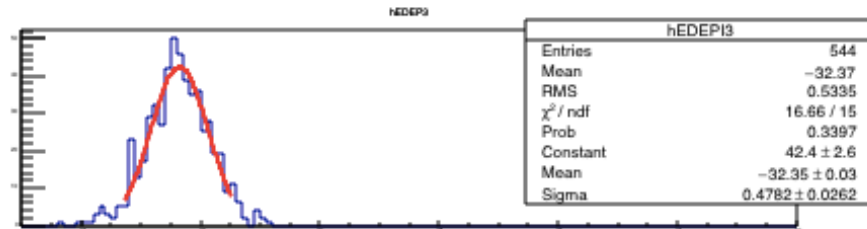
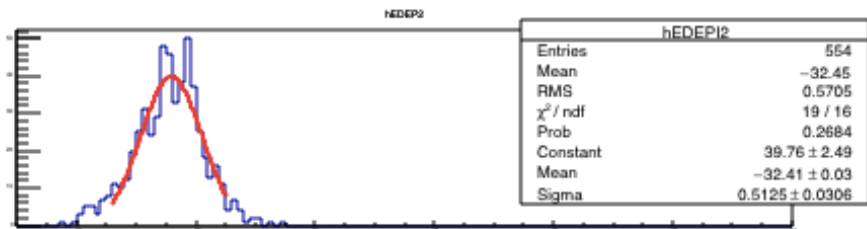
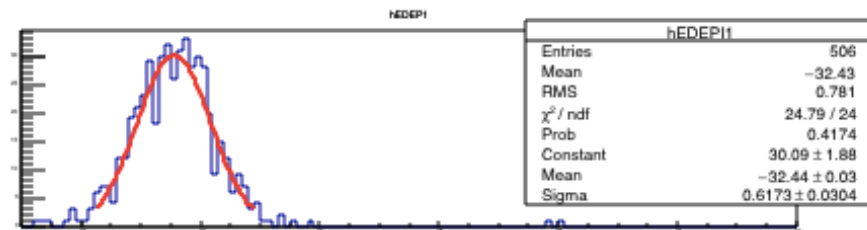
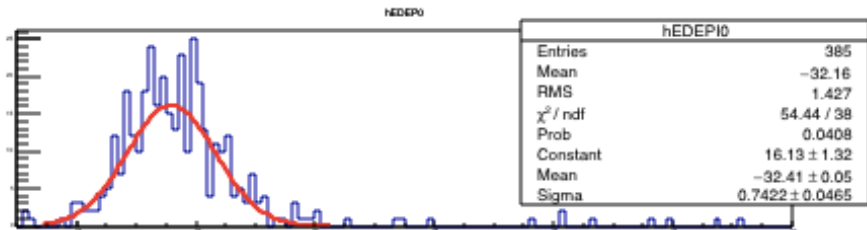
Raw distributions

Run69 All Modules



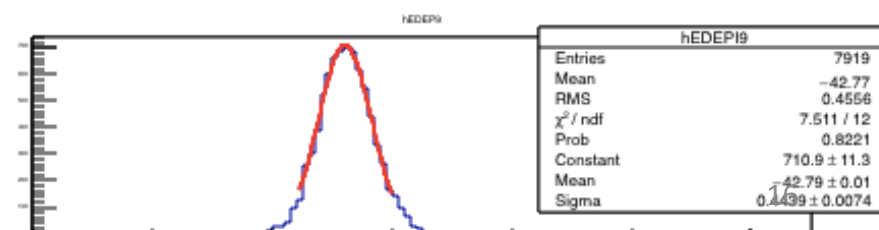
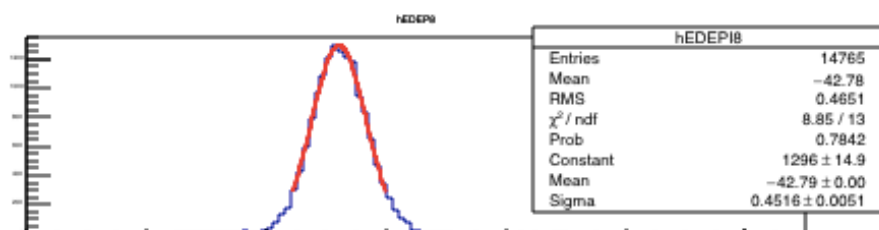
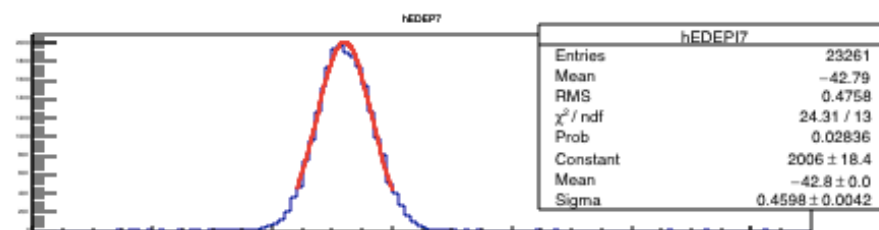
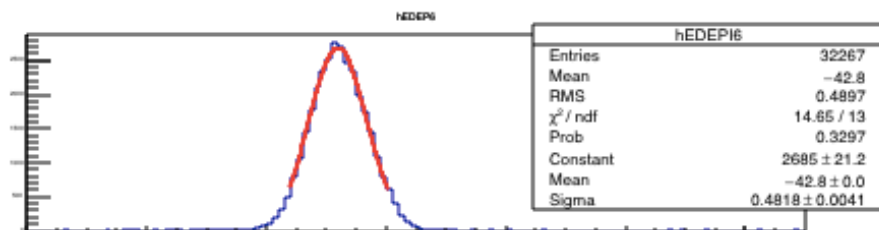
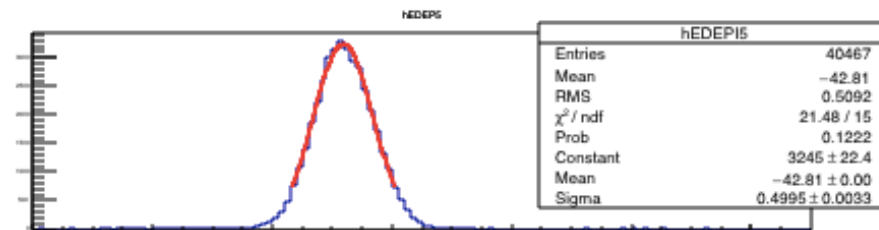
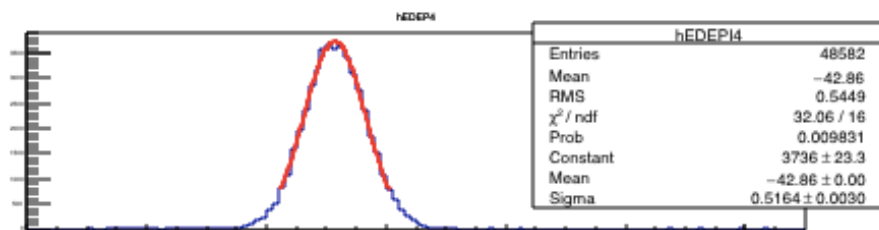
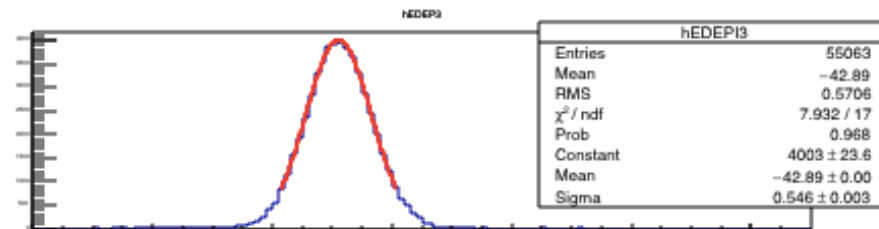
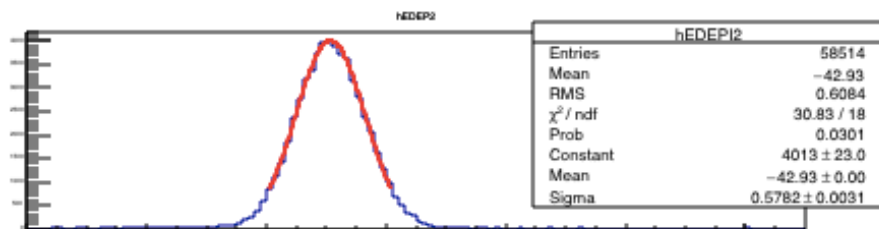
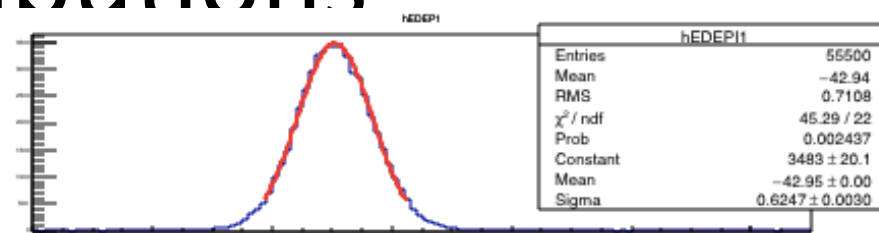
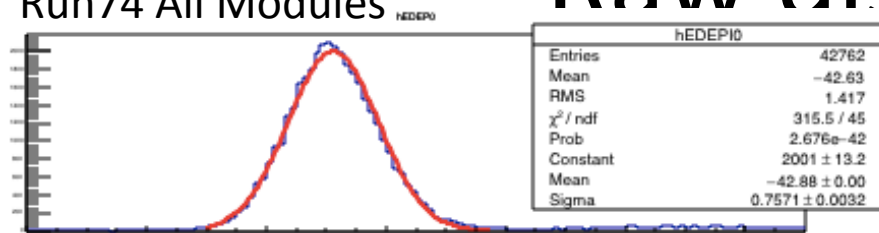
Raw distributions

Run69 One Module



Raw distributions

Run74 All Modules



Energy Deposit

Black : Original MC

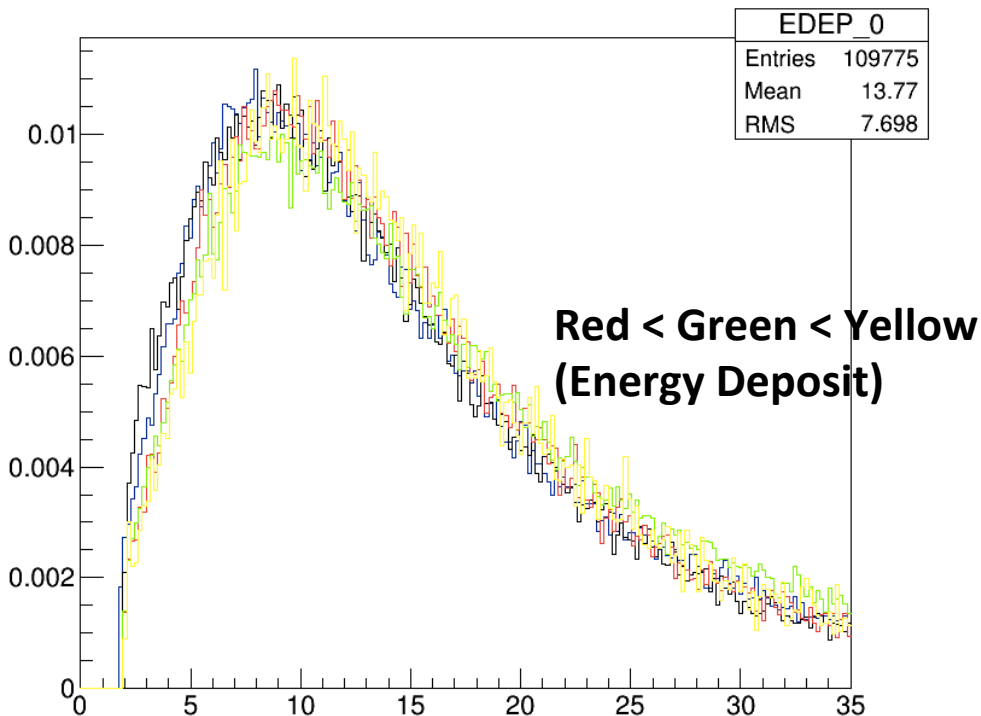
Blue : 0.9 Scaled Energy

Red : MC with 0.1mm reduction

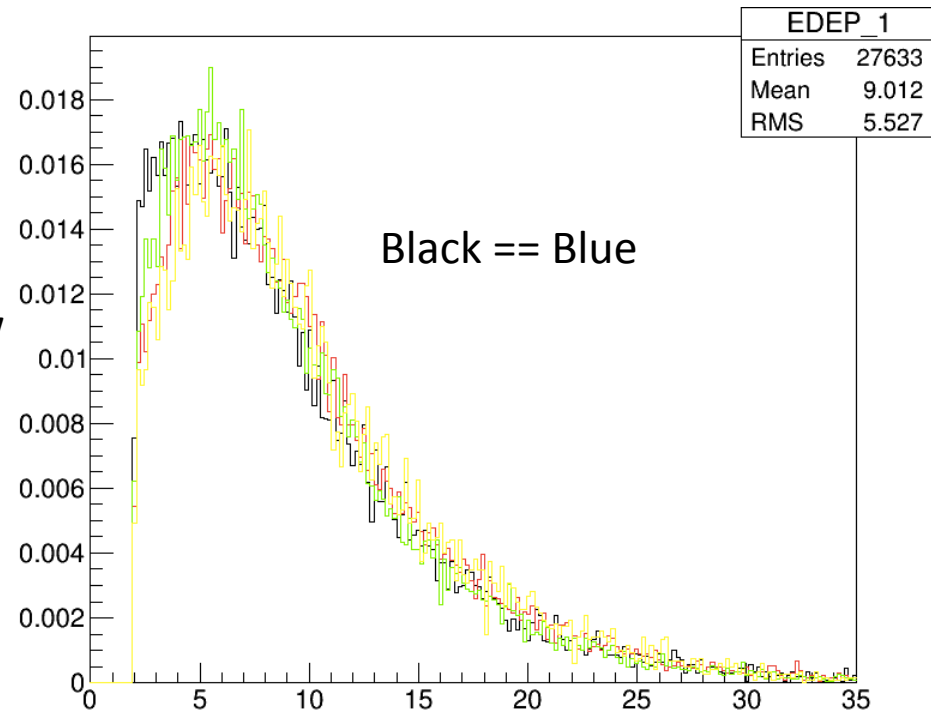
Green : Original Data

Yellow : MC with 0.2mm reduction

IMB



OMB

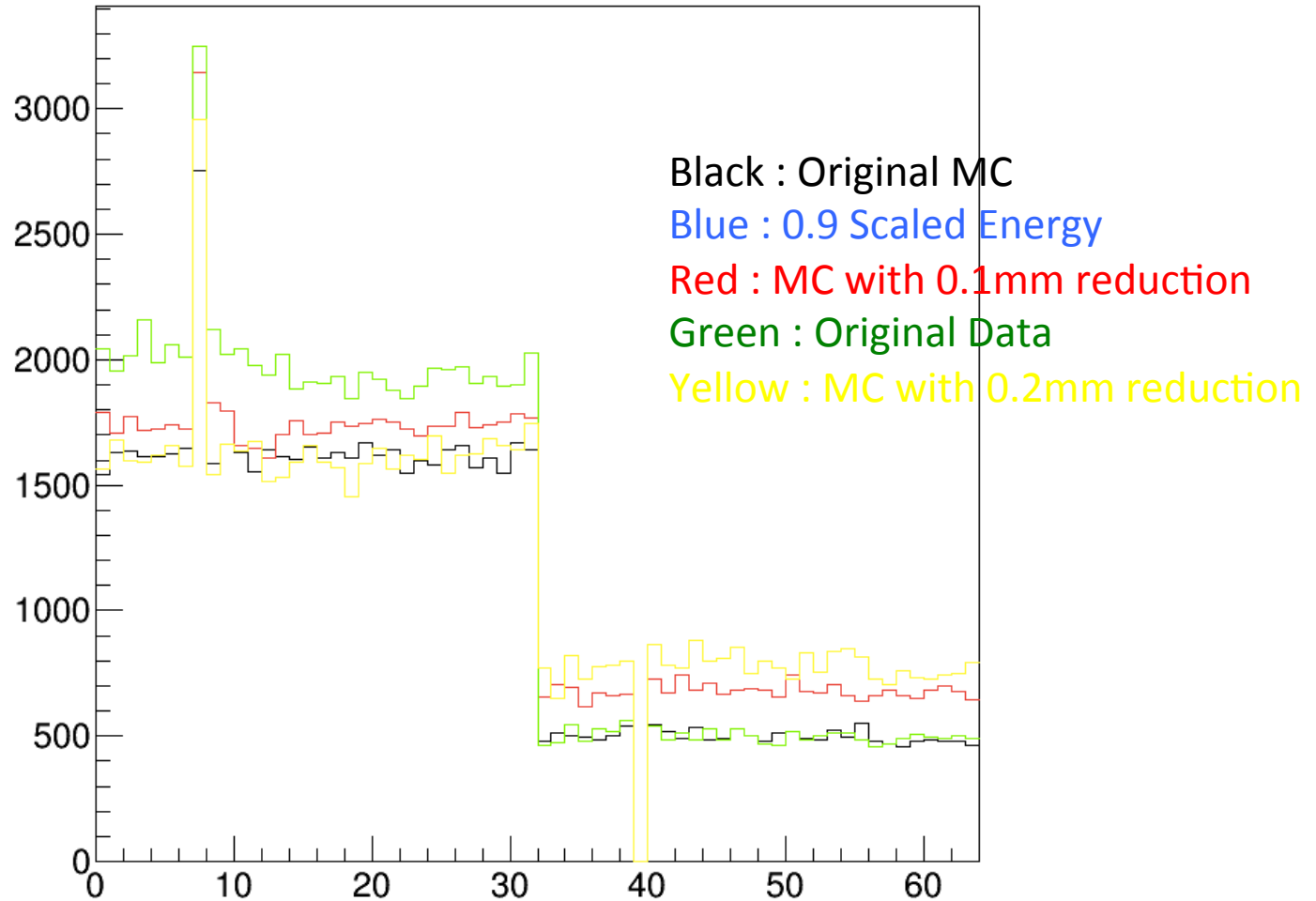


P.O.T. to KL

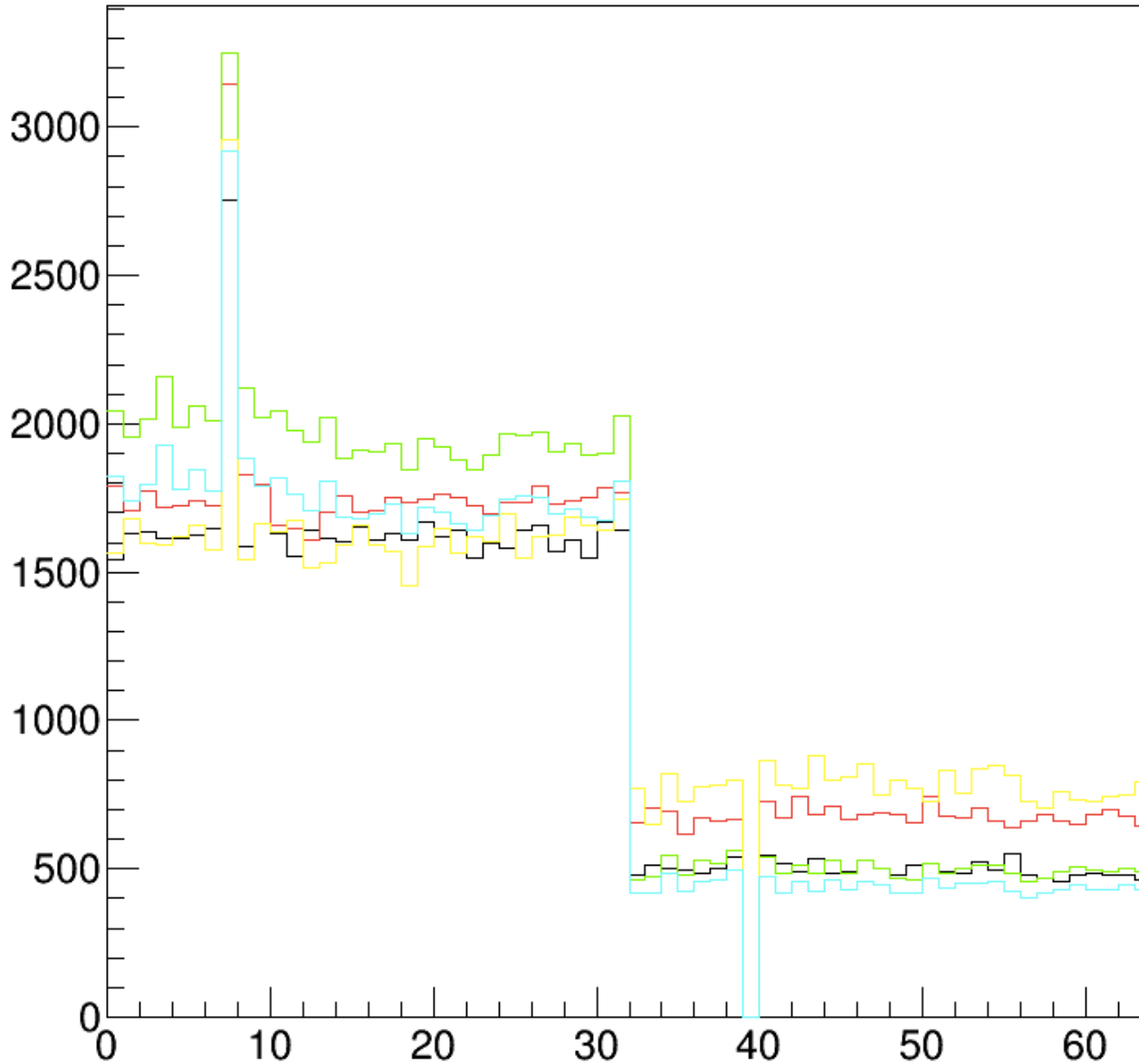
- P.O.T. in Run65 : $1.286e19$
- #KLs = $8.146e9$ in Min. bias
- #KLs in MC = $1e9$
 - KL3pi0 only, Total KL = $1e9 / 0.192 = 5.12e9$

After P.O.T. Normalization

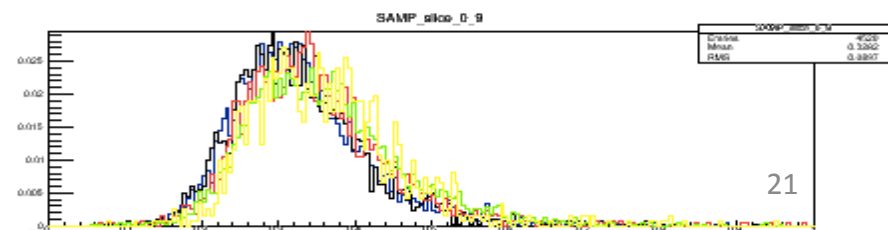
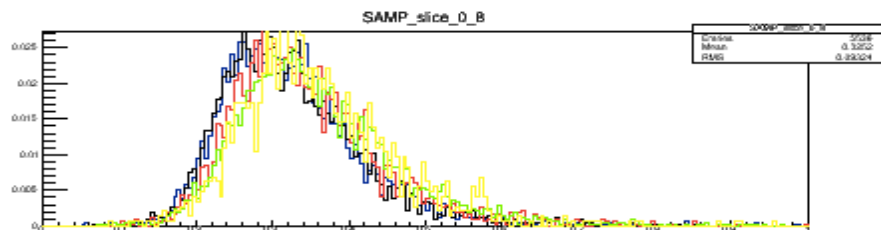
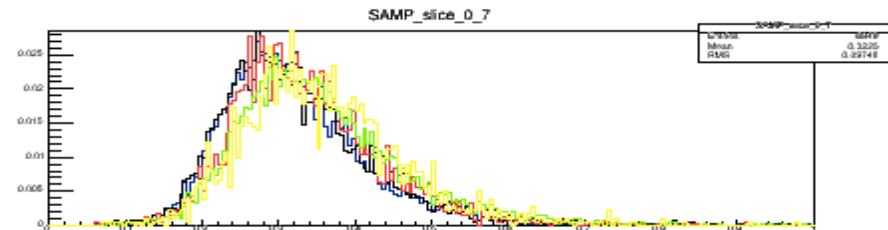
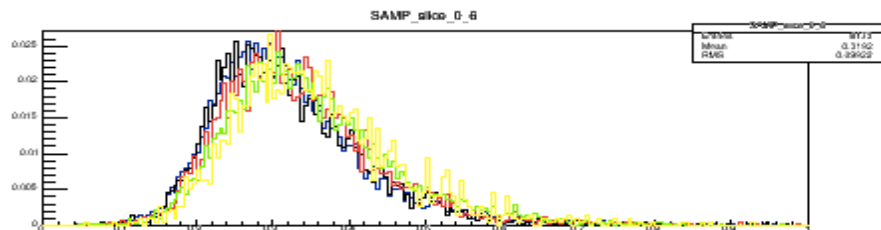
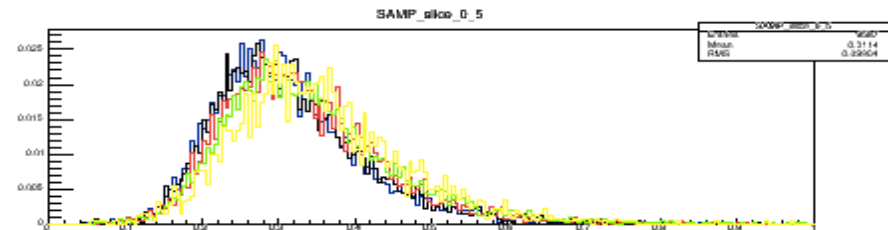
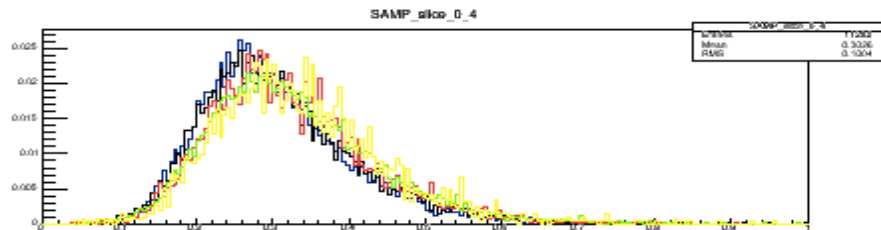
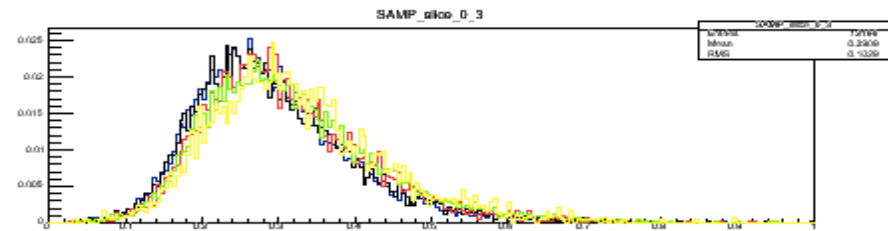
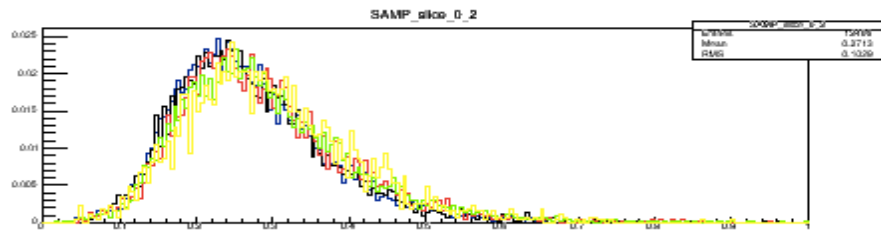
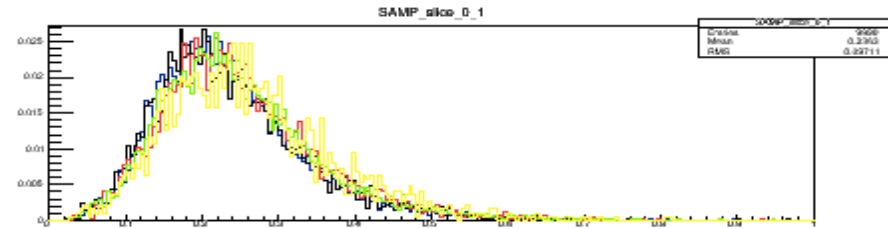
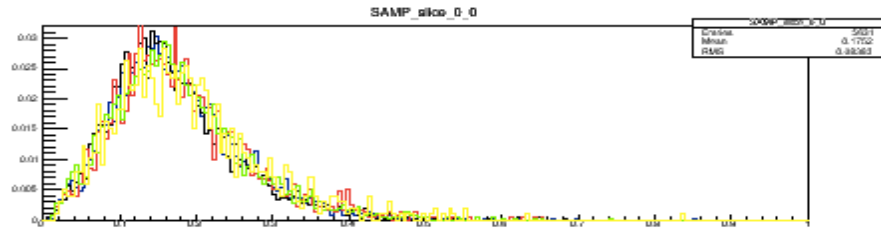
With Hit Requirement (#Hits == 1)



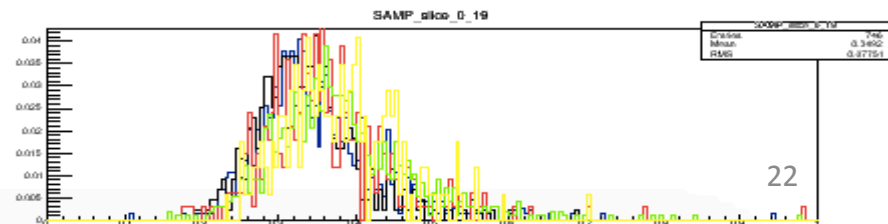
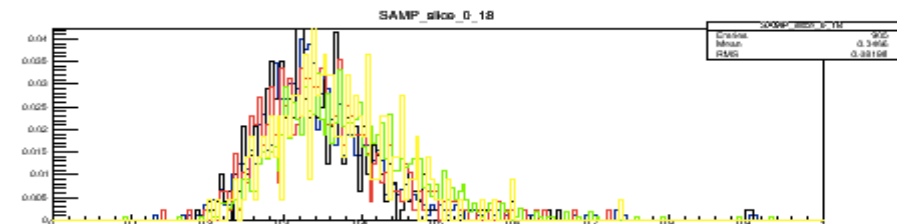
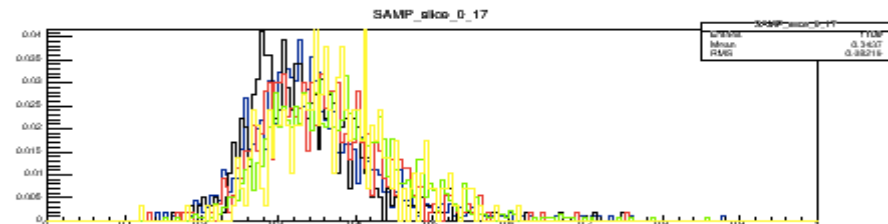
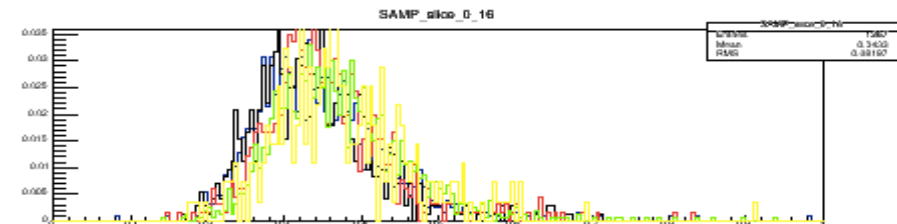
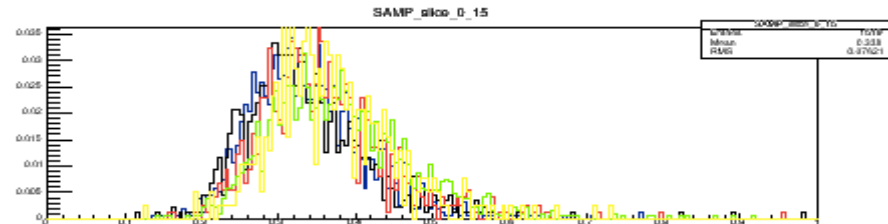
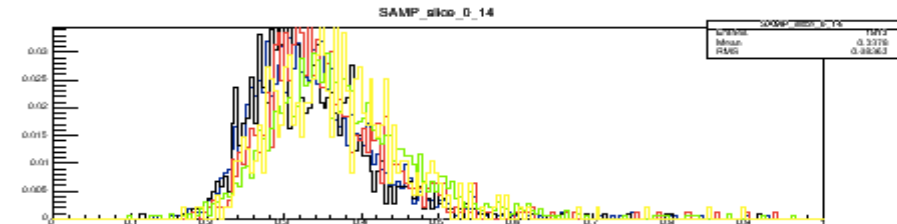
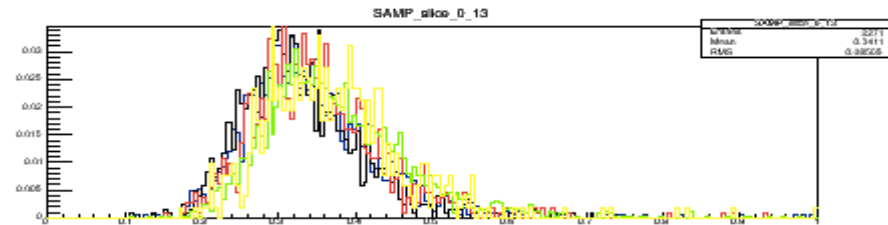
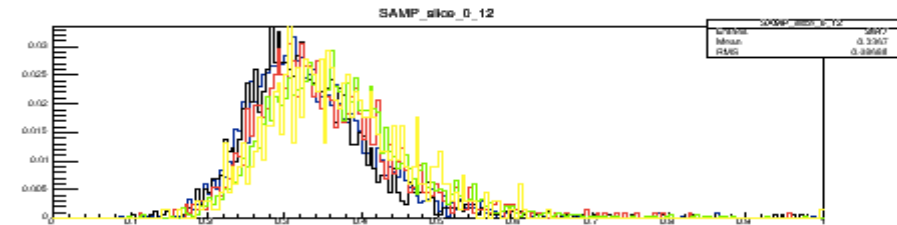
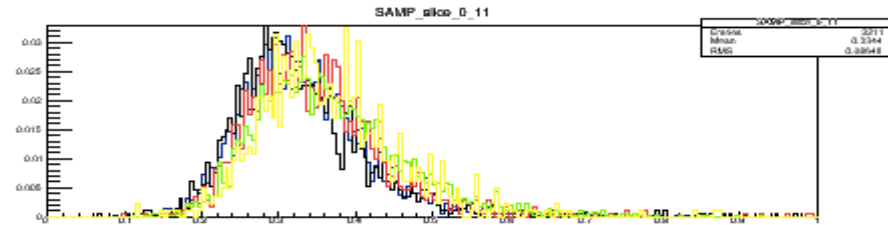
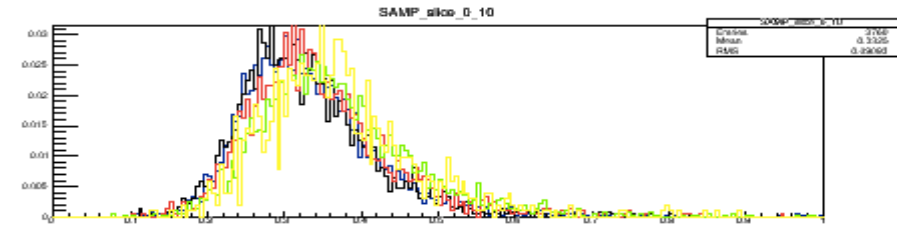
Without Hit Req



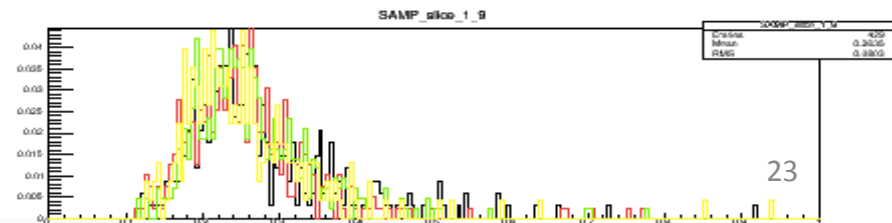
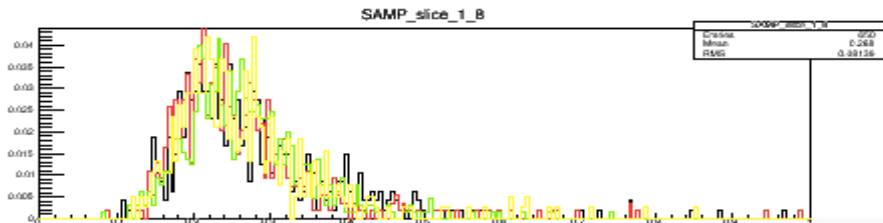
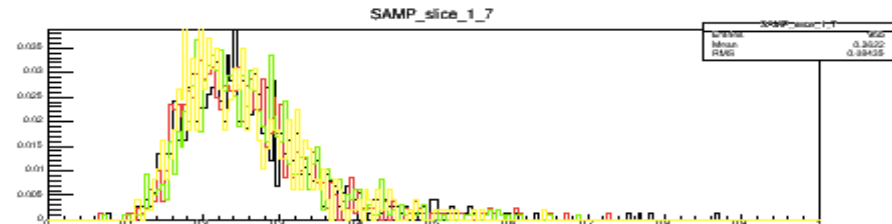
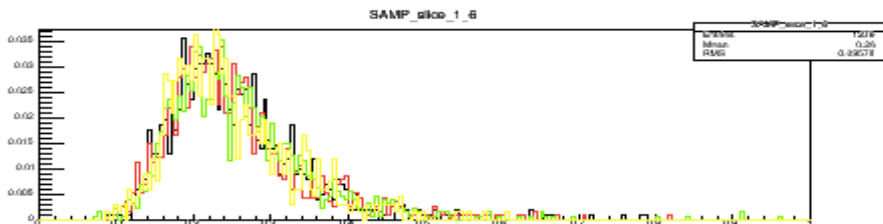
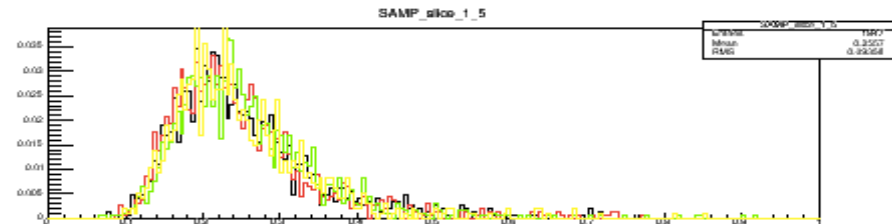
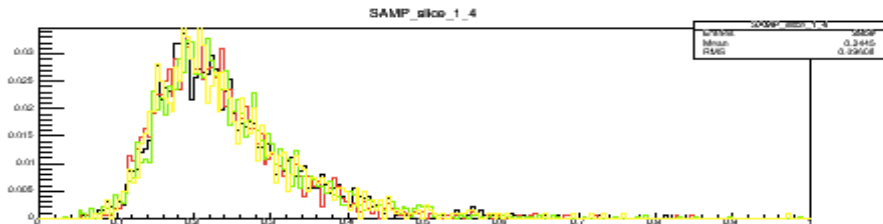
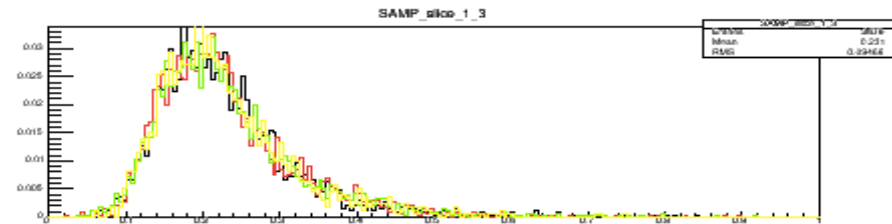
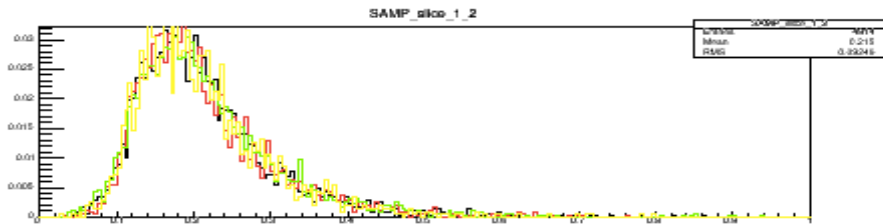
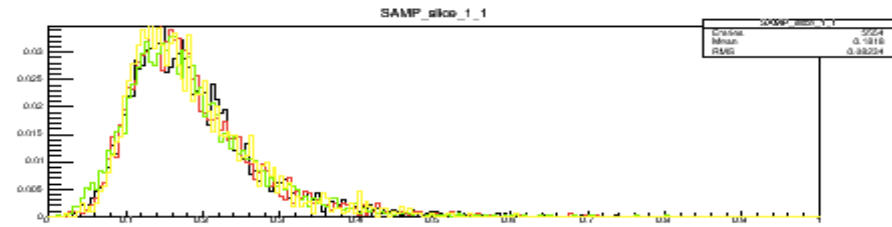
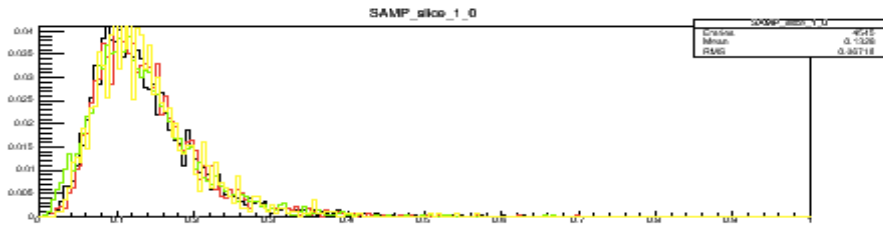
Sampling Fraction (IMB)



Sampling Fraction (IMB)



Sampling Fraction (OMB)



Material	Z	A	$\langle Z/A \rangle$	Nucl.coll. length λ_T {g cm ⁻² }	Nucl.inter. length λ_I {g cm ⁻² }	Rad.len. X_0 {g cm ⁻² }	$dE/dx _{\min}$ { MeV g ⁻¹ cm ² }	Density {g cm ⁻³ } {gℓ ⁻¹ }	Melting point (K)	Boiling point (K)	Refract. index @ Na D
H ₂	1	1.008(7)	0.99212	42.8	52.0	63.04	(4.103)	0.071(0.084)	13.81	20.28	1.11[132.]
D ₂	1	2.01410177803(8)	0.49650	51.3	71.8	125.97	(2.053)	0.169(0.168)	18.7	23.65	1.11[138.]
He	2	4.002602(2)	0.49967	51.8	71.0	94.32	(1.937)	0.125(0.166)		4.220	1.02[35.0]
Li	3	6.94(2)	0.43221	52.2	71.3	82.78	1.639	0.534	453.6	1615.	
Be	4	9.0121831(5)	0.44384	55.3	77.8	65.19	1.595	1.848	1560.	2744.	
C diamond	6	12.0107(8)	0.49955	59.2	85.8	42.70	1.725	3.520			2.42
C graphite	6	12.0107(8)	0.49955	59.2	85.8	42.70	1.742	2.210			
N ₂	7	14.007(2)	0.49976	61.1	89.7	37.99	(1.825)	0.807(1.165)	63.15	77.29	1.20[298.]
O ₂	8	15.999(3)	0.50002	61.3	90.2	34.24	(1.801)	1.141(1.332)	54.36	90.20	1.22[271.]
F ₂	9	18.998403163(6)	0.47372	65.0	97.4	32.93	(1.676)	1.507(1.580)	53.53	85.03	[195.]
Ne	10	20.1797(6)	0.49555	65.7	99.0	28.93	(1.724)	1.204(0.839)	24.56	27.07	1.09[67.1]
Al	13	26.9815385(7)	0.48181	69.7	107.2	24.01	1.615	2.699	933.5	2792.	
Si	14	28.0855(3)	0.49848	70.2	108.4	21.82	1.664	2.329	1687.	3538.	3.95
Cl ₂	17	35.453(2)	0.47951	73.8	115.7	19.28	(1.630)	1.574(2.980)	171.6	239.1	[773.]
Ar	18	39.948(1)	0.45059	75.7	119.7	19.55	(1.519)	1.396(1.662)	83.81	87.26	1.23[281.]
Ti	22	47.867(1)	0.45961	78.8	126.2	16.16	1.477	4.540	1941.	3560.	
Fe	26	55.845(2)	0.46557	81.7	132.1	13.84	1.451	7.874	1811.	3134.	
Cu	29	63.546(3)	0.45636	84.2	137.3	12.86	1.403	8.960	1358.	2835.	
Ge	32	72.630(1)	0.44053	86.9	143.0	12.25	1.370	5.323	1211.	3106.	
Sn	50	118.710(7)	0.42119	98.2	166.7	8.82	1.263	7.310	505.1	2875.	
Xe	54	131.293(6)	0.41129	100.8	172.1	8.48	(1.255)	2.953(5.483)	161.4	165.1	1.39[701.]
W	74	183.84(1)	0.40252	110.4	191.9	6.76	1.145	19.300	3695.	5828.	
Pt	78	195.084(9)	0.39983	112.2	195.7	6.54	1.128	21.450	2042.	4098.	
Au	79	196.966569(5)	0.40108	112.5	196.3	6.46	1.134	19.320	1337.	3129.	
Pb	82	207.2(1)	0.39575	114.1	199.6	6.37	1.122	11.350	600.6	2022.	
U	92	[238.02891(3)]	0.38651	118.6	209.0	6.00	1.081	18.950	1408.	4404.	
Polystyrene ((C ₆ H ₅ CHCH ₂) _n)			0.53768	57.5	81.7	43.79	1.936	1.06			1.59