

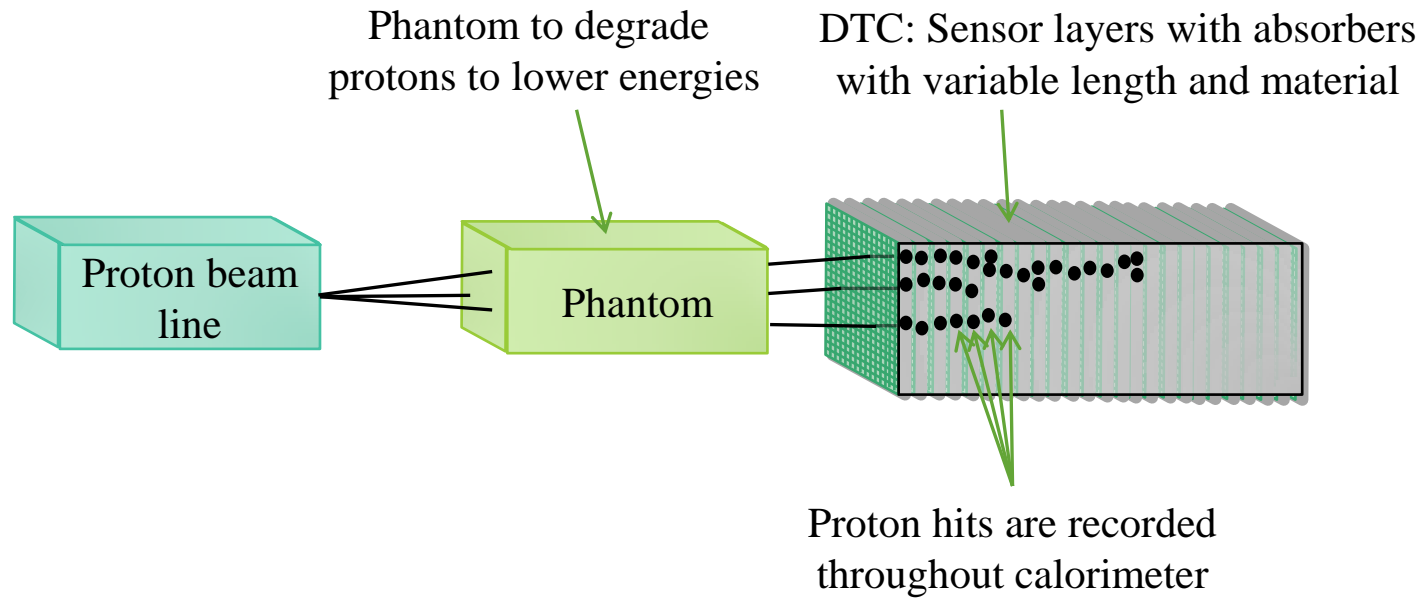
Optimizing the DTC

2017-03-03

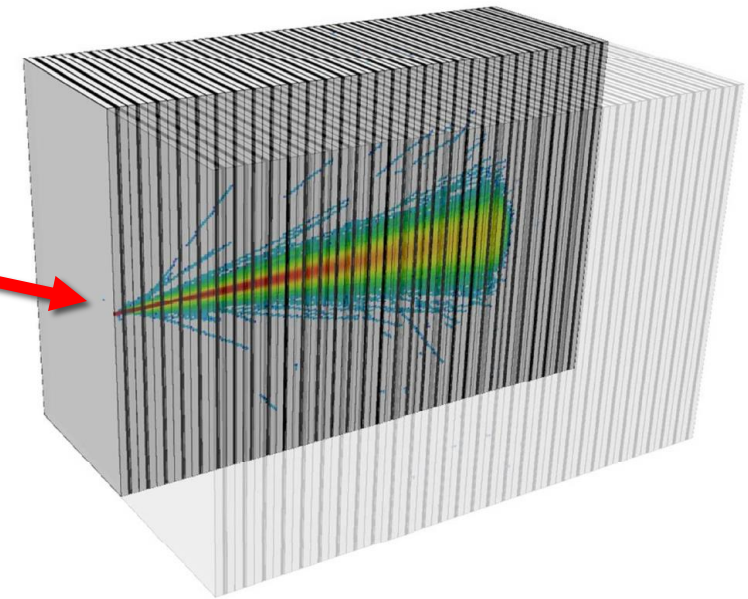
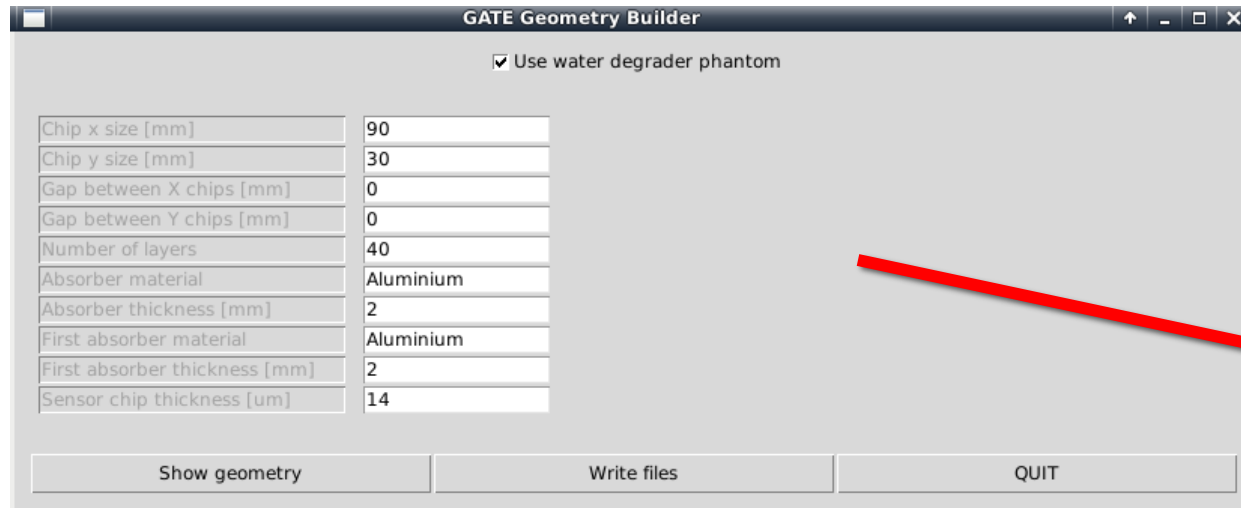
Goals

1. To find the best possible design for the range calorimeter part of DTC
2. Constrained by \$\$\$ and Range Straggling
 1. Need to contain a 250 MeV proton beam
 2. Linear resolution to be limited somewhat by range straggling $\sim 1\%$

Method



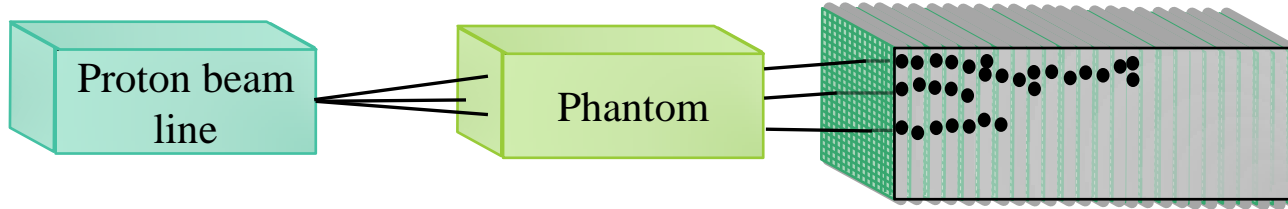
GATE setup



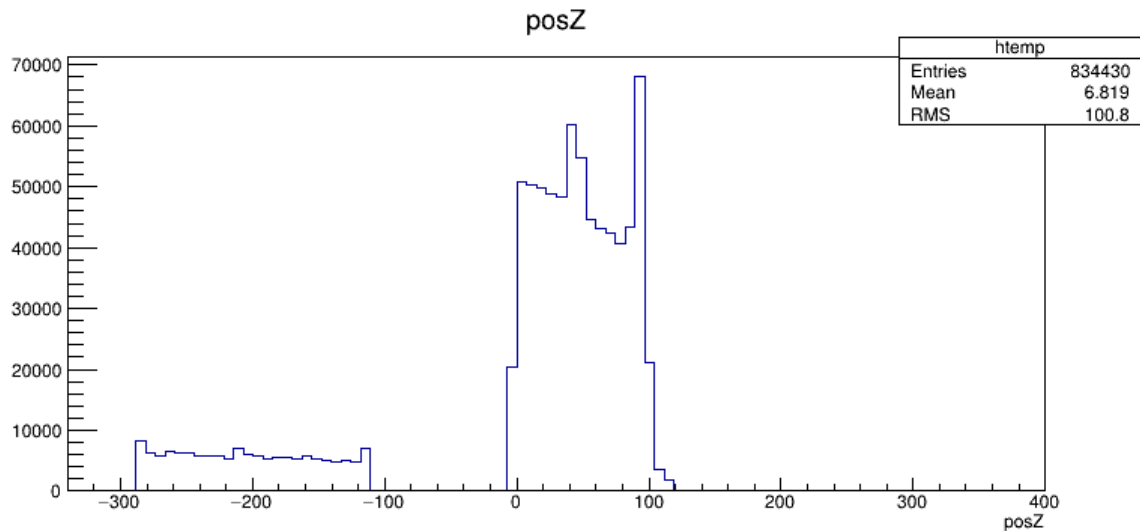
For example, with a 3 mm degrader, and simulating a 250 MeV beam passing through a phantom of 50, 55, 60, 65 and 70 mm water:

```
[gate/python] $ sh runDegrader.sh 3 50 5 70
```

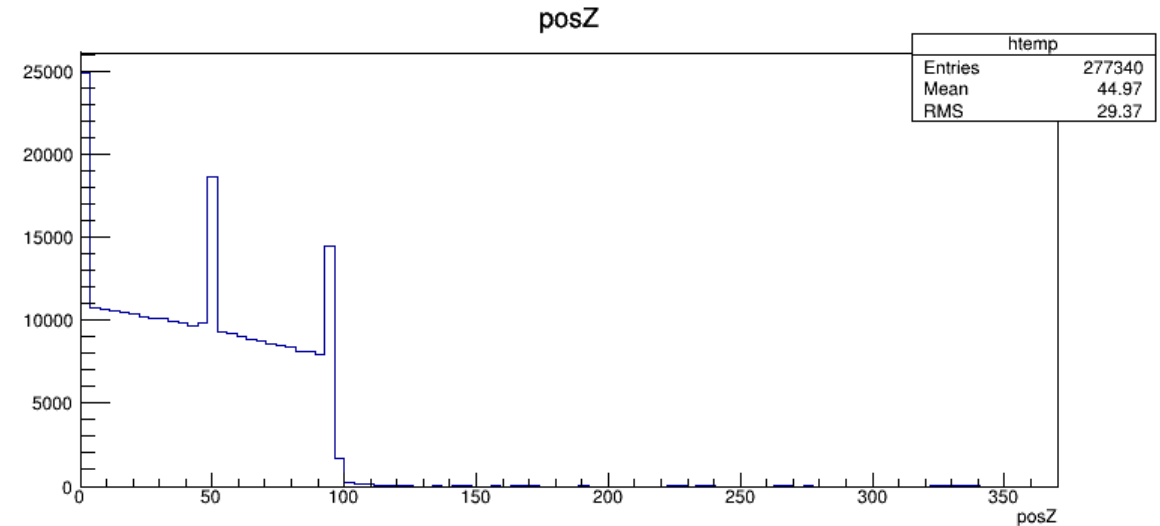
Monte Carlo simulations



Full simulation: Record everything: 5' primaries/energy



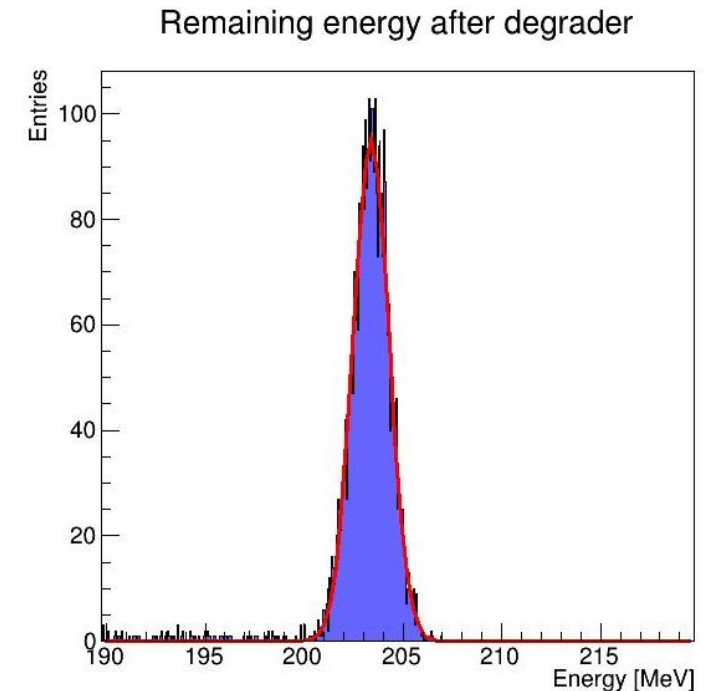
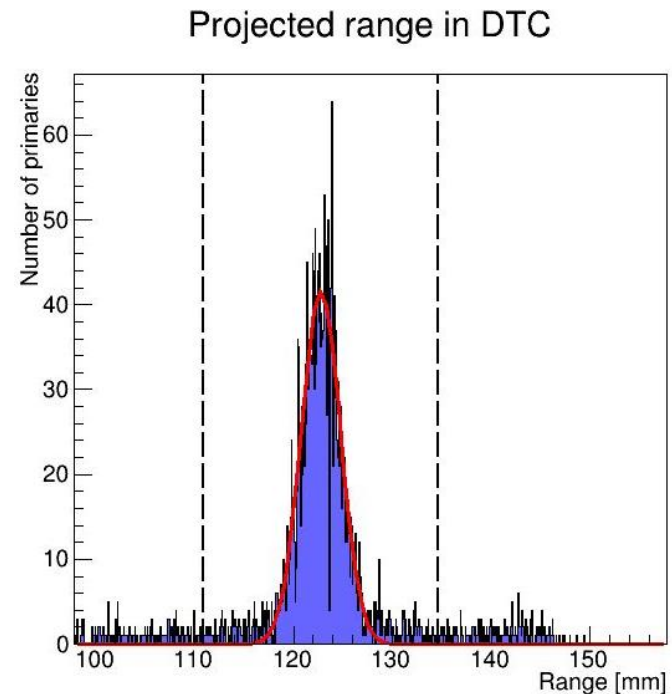
Chip simulation: Record only events in sensors:
15' primaries/energy



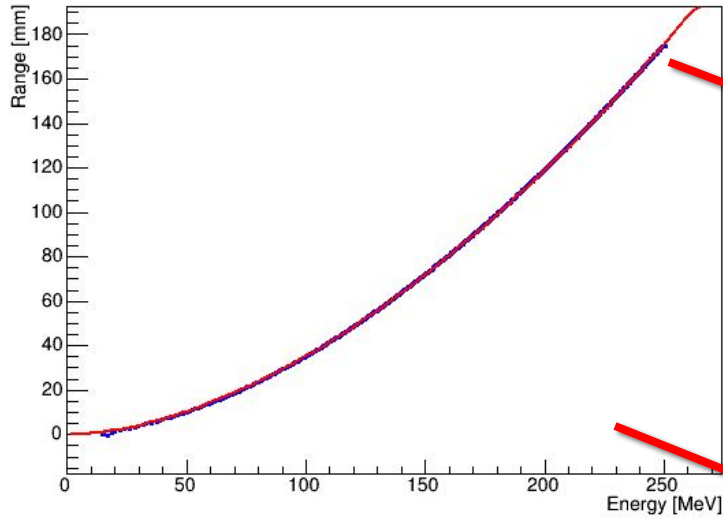
Full MC simulations

The «Gold standard» in this context

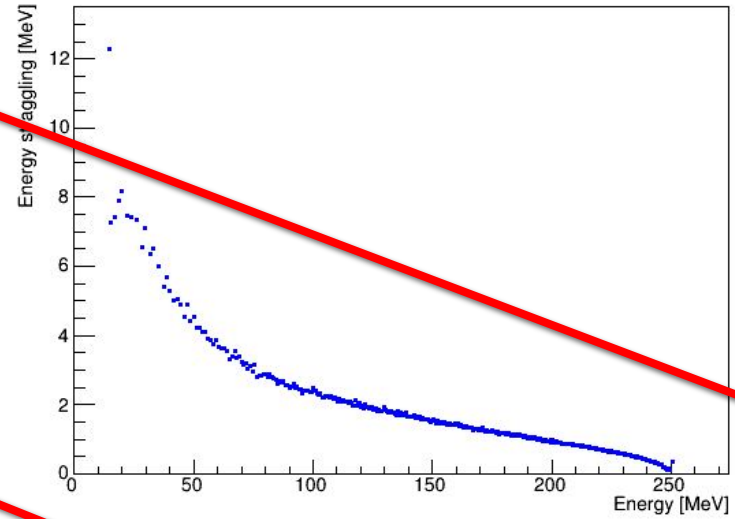
1. Range for each water phantom thickness and geometry configuration
2. Range straggling
3. Energy spread distal to DTC



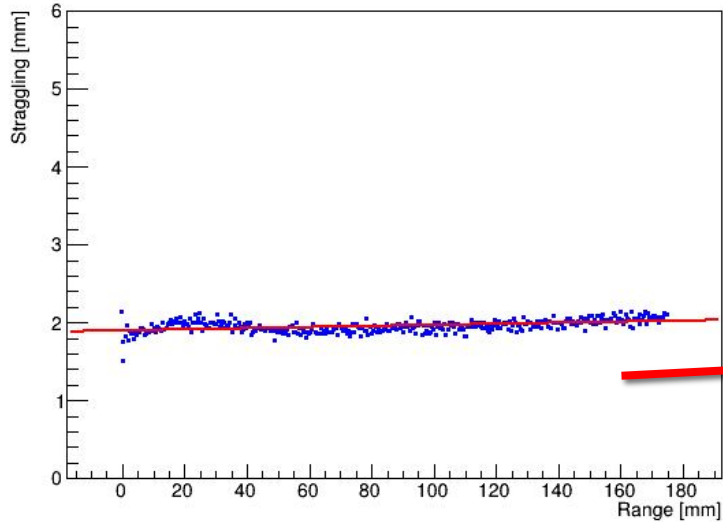
Ranges for material in 3 mm Al



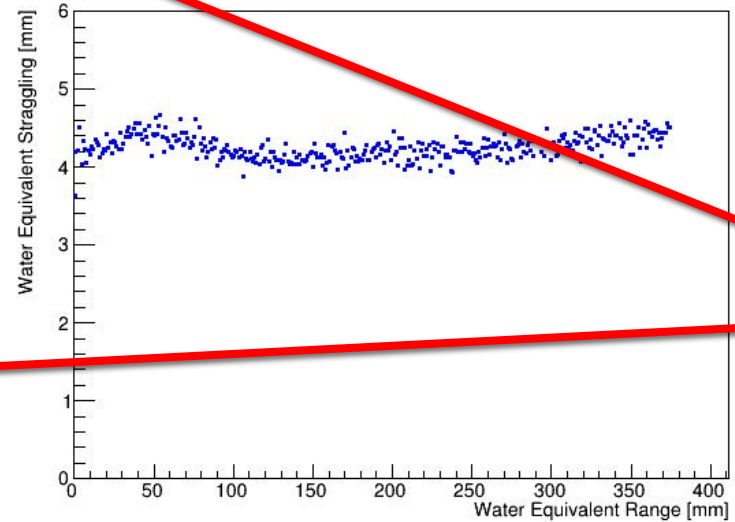
Energy straggling in DTC for 3 mm absorber



Straggling in Al for 3 mm absorber



WE straggling in Al for 3 mm absorber



1	#	ENERGY	RANGE
2	14.845	0.056788	
3	15.3795	0.104973	
4	17.2187	-0.215777	
5	18.6584	0.855986	
6	20.1045	1.2952	
7	22.0102	1.93137	
8	23.889	2.34508	
9	25.8993	2.82206	
10	28.2813	3.44492	

In addition, the following parameters should be extracted:

Bragg-Kleeman parameters: $R = 0.011626 E^{1.743151}$

Straggling = $1.8568 + 0.000856 R$

Configuring the chip MC simulations

```
81 void createSplines() {  
...  
107     else if (kAbsorbatorThickness = 3) {  
108         in.open("Data/Ranges/3mm_Al.csv");  
109     }  
...  
192     else if (kAbsorbatorThickness = 3) {  
193         alpha_aluminum = 0.011626;  
194         p_aluminum = 1.743151;  
195         stragglng_a = 1.8568;  
196         stragglng_b = 0.000856;  
197     }
```

Or in the corresponding material (alpha_pmma, alpha_carbon, etc.) and absorbatorthickness lines.

And in the file `DTCToolkit/Scripts/makePlots.C`, put the \alpha, p parameters.

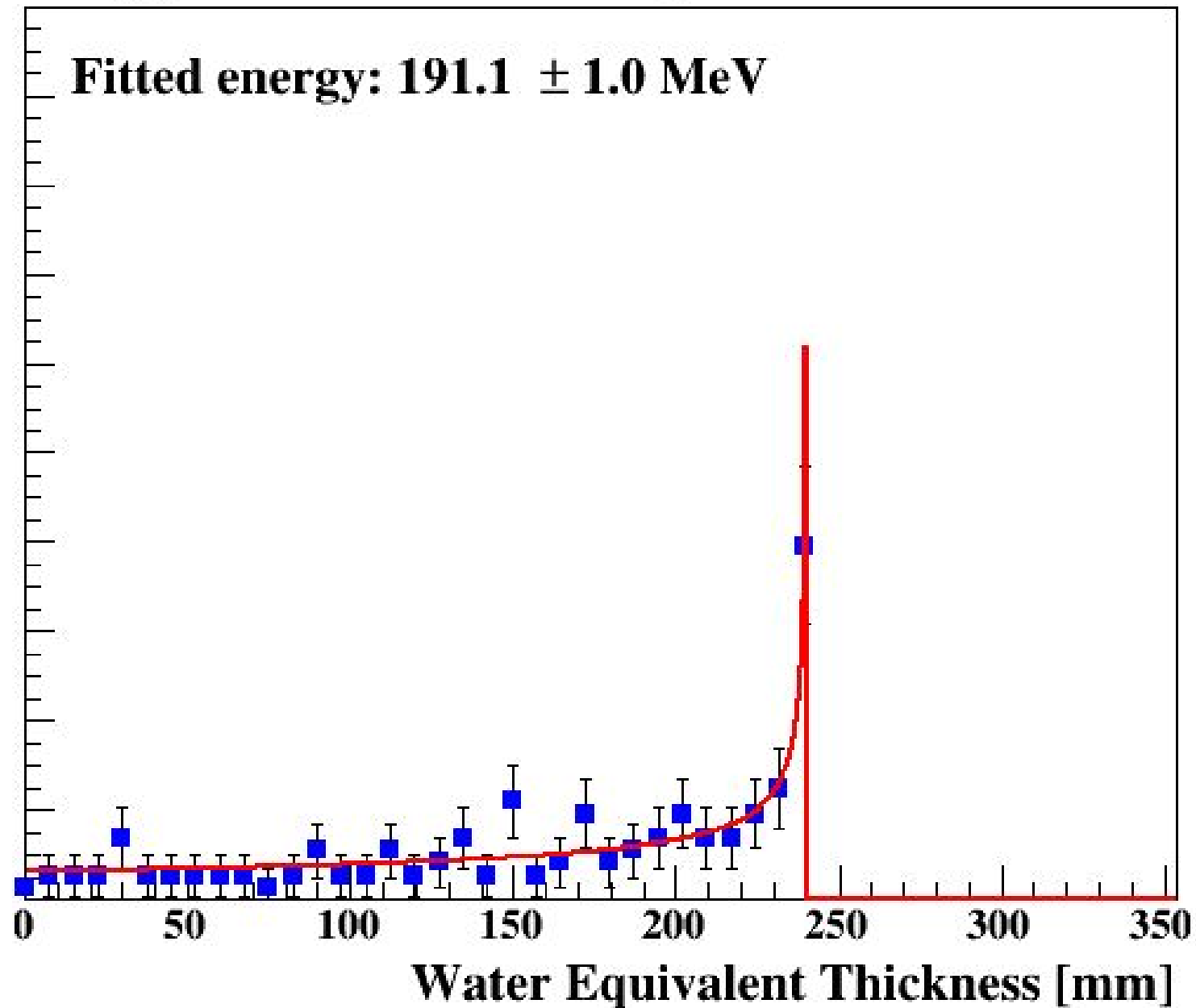
```
144     else if (absorberThickness == 3) {  
145         a_dtc = 0.011626;  
146         p_dtc = 1.743151;  
147     }
```


Chip MC simulations

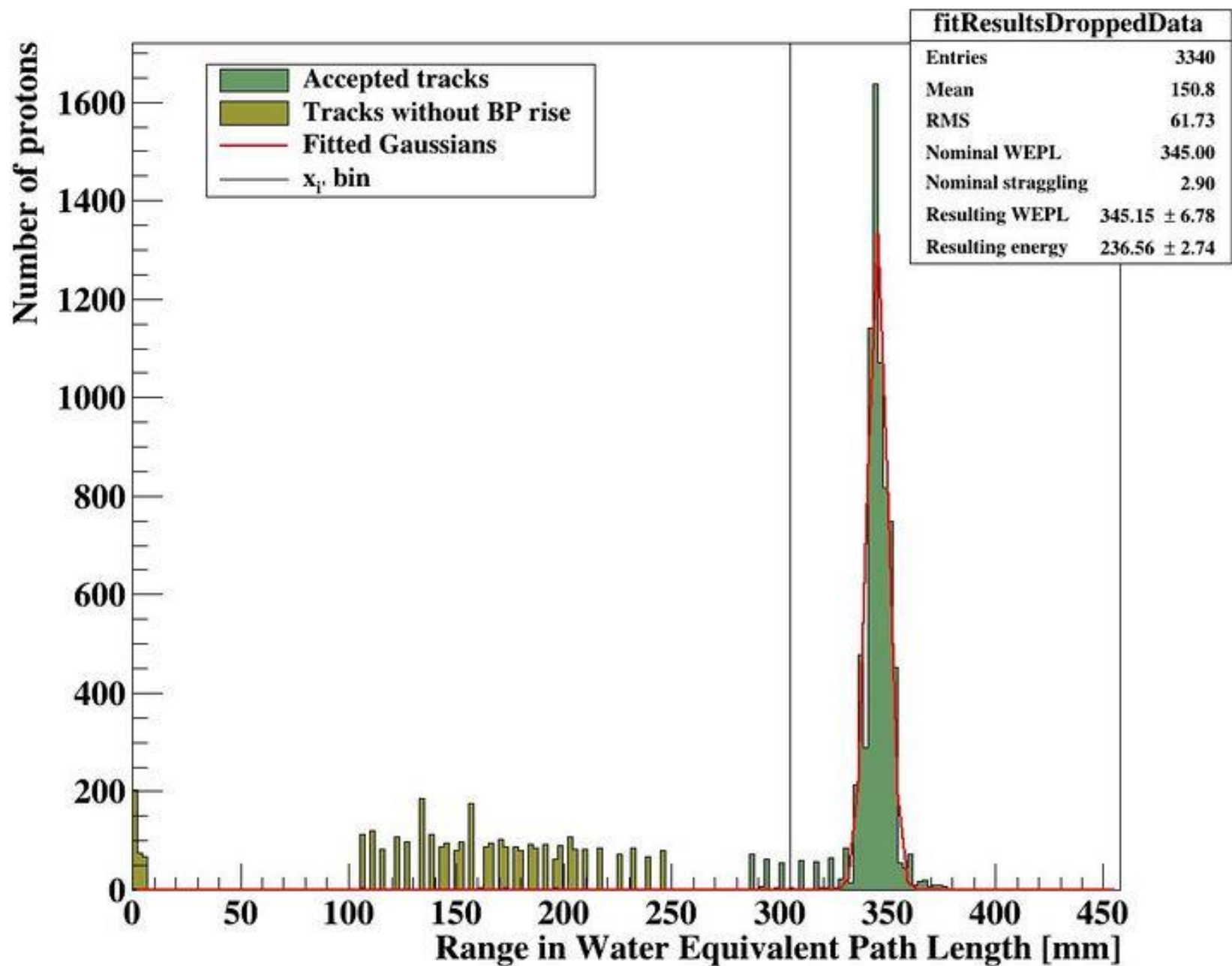
From the GATE data, we do:

1. Make Cluster objects containing (x,y,layer,edep,eventID) from the MC ROOT files **(simplified)**
2. Make Tracks using the eventID **(simplified)**
3. Find optimal range for each track using Bragg curve model fit with layer/edep
4. Find Mean and Width of the range distribution of all proton tracks
 1. Mean: DTC Accuracy
 2. Width: DTC Resolution
5. Repeat for all water degrader thicknesses (\simeq 400 energies per geometry)
6. Make plot of the variation of Accuracy and Resolution with water degrader thicknesses

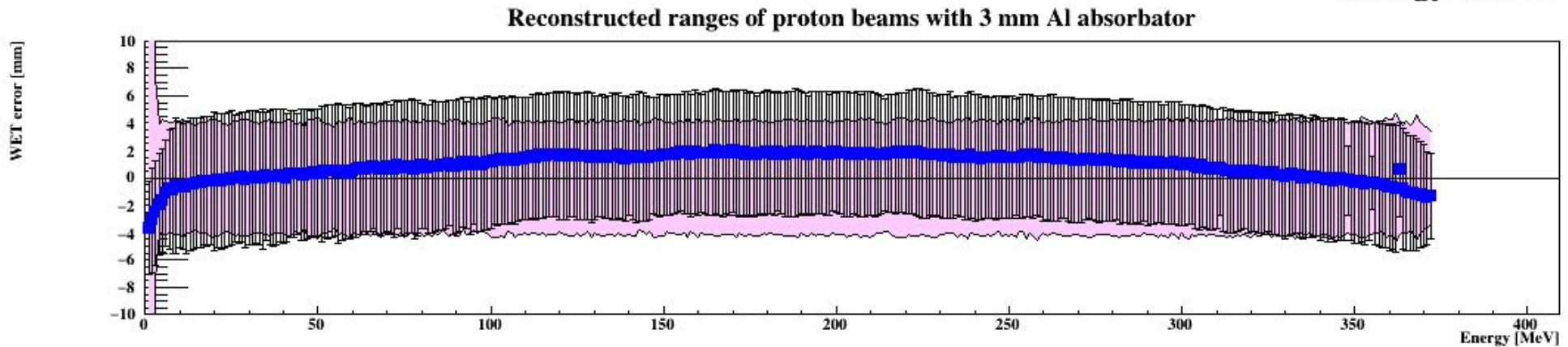
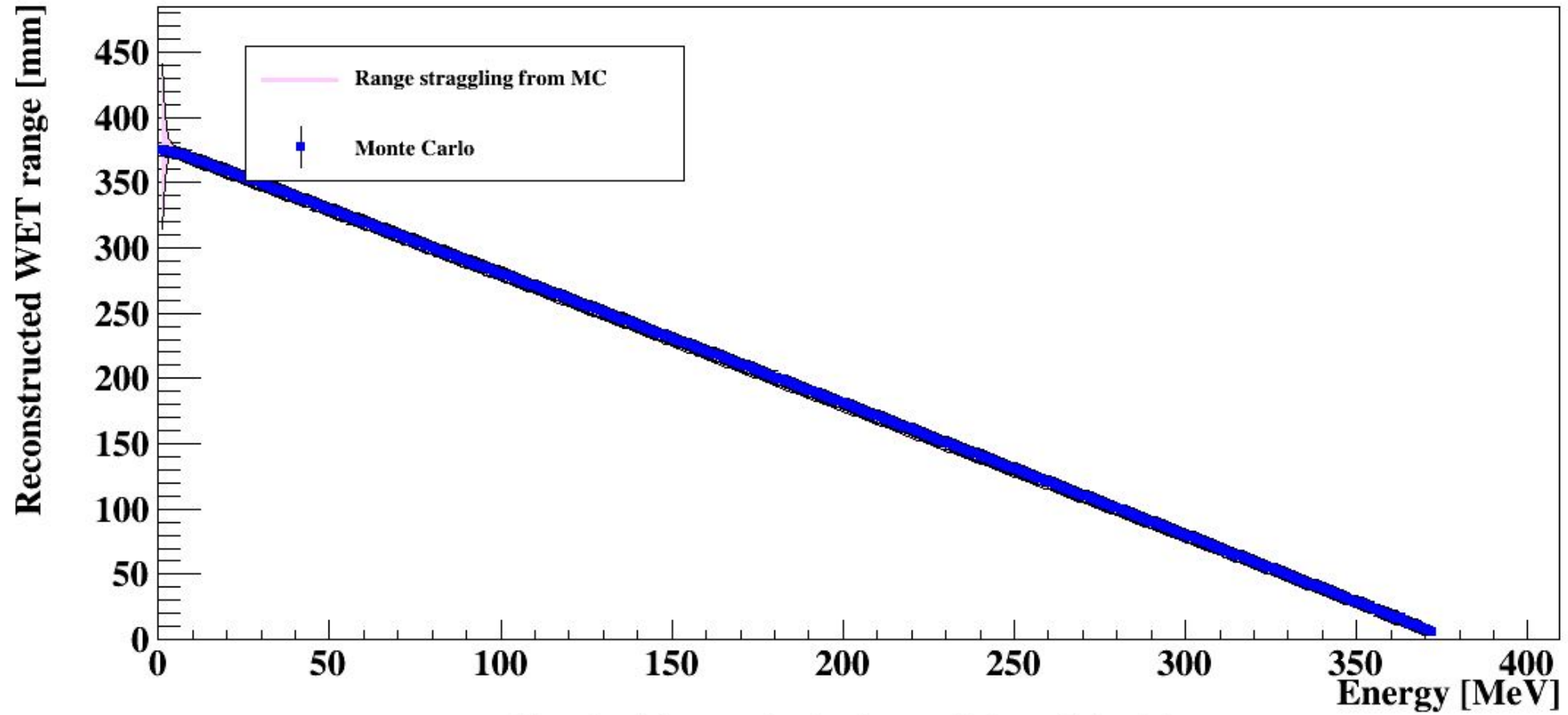
Bragg-Kleeman fit to exp. data at 192 MeV



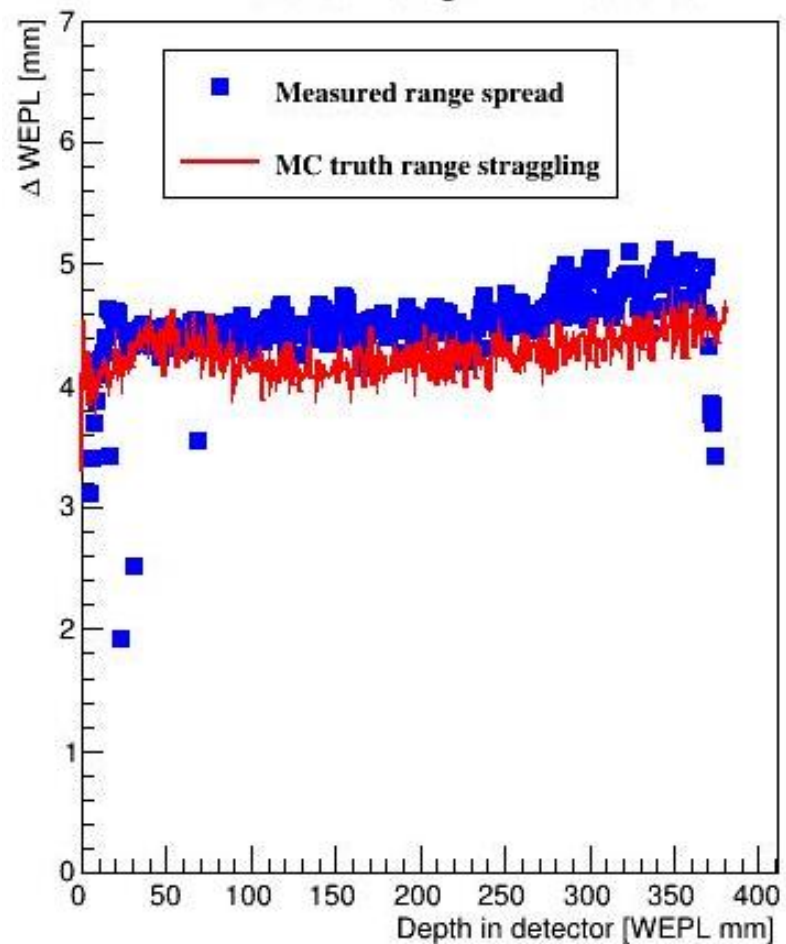
Fitted energy of a 250 MeV nominal beam on Aluminium DTC w/34.0 mm water degrader



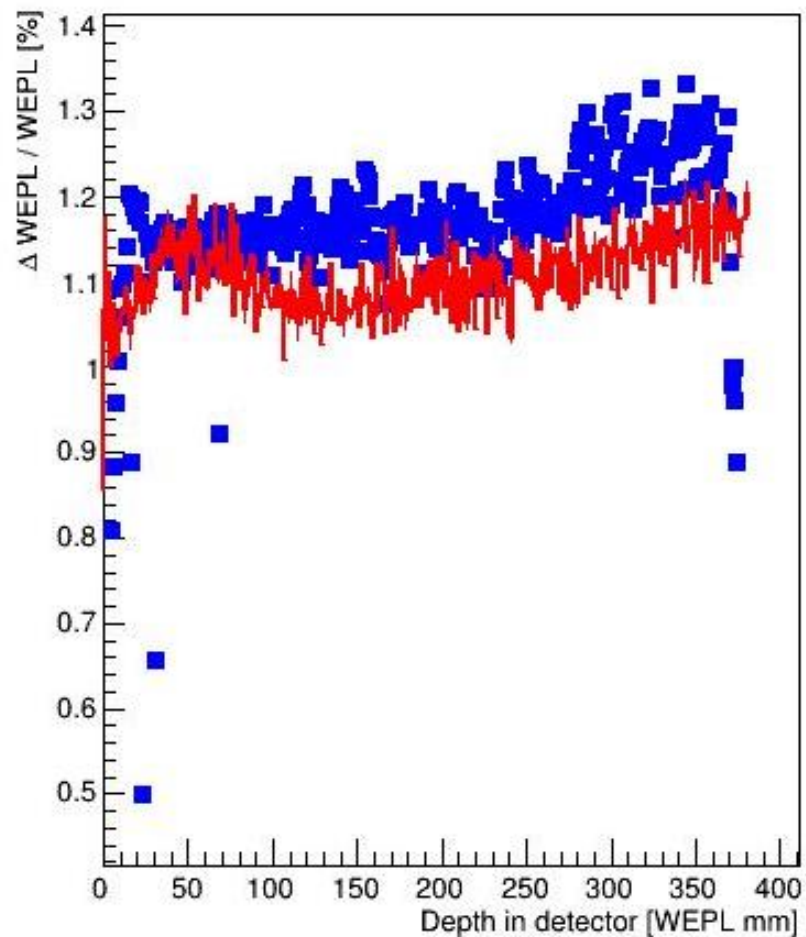
Reconstructed ranges of proton beams with 3 mm Al absorbator



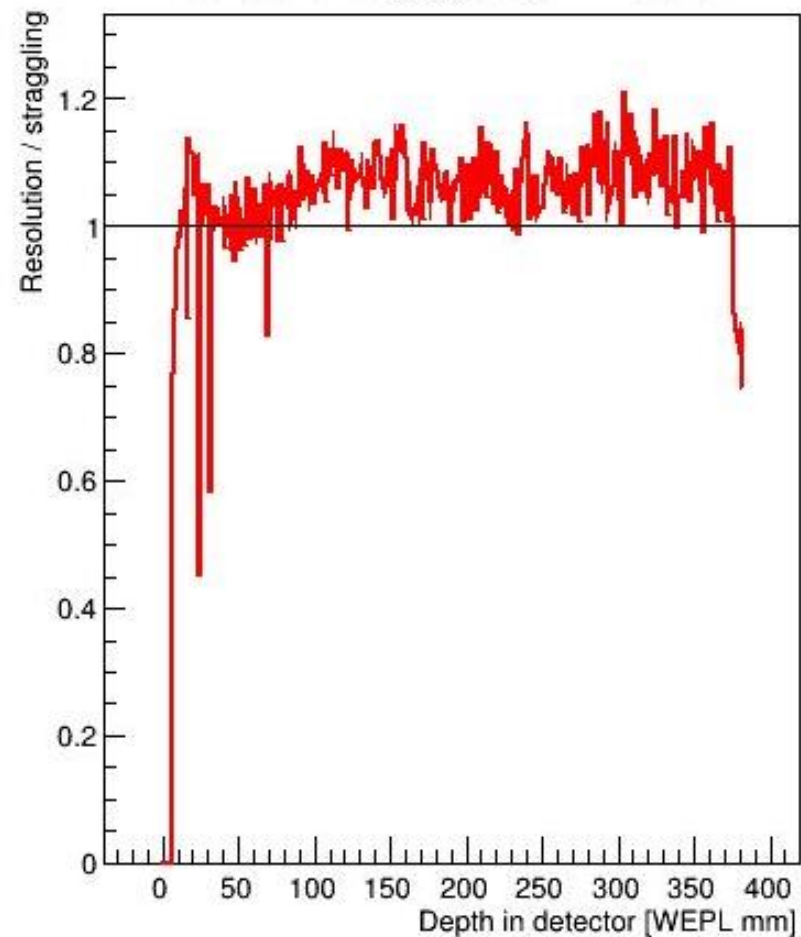
WEPL resolution using 3 mm Al absorber



WEPL resolution using 3 mm Al absorber



Ratio resolution / straggling using 3 mm Al absorber



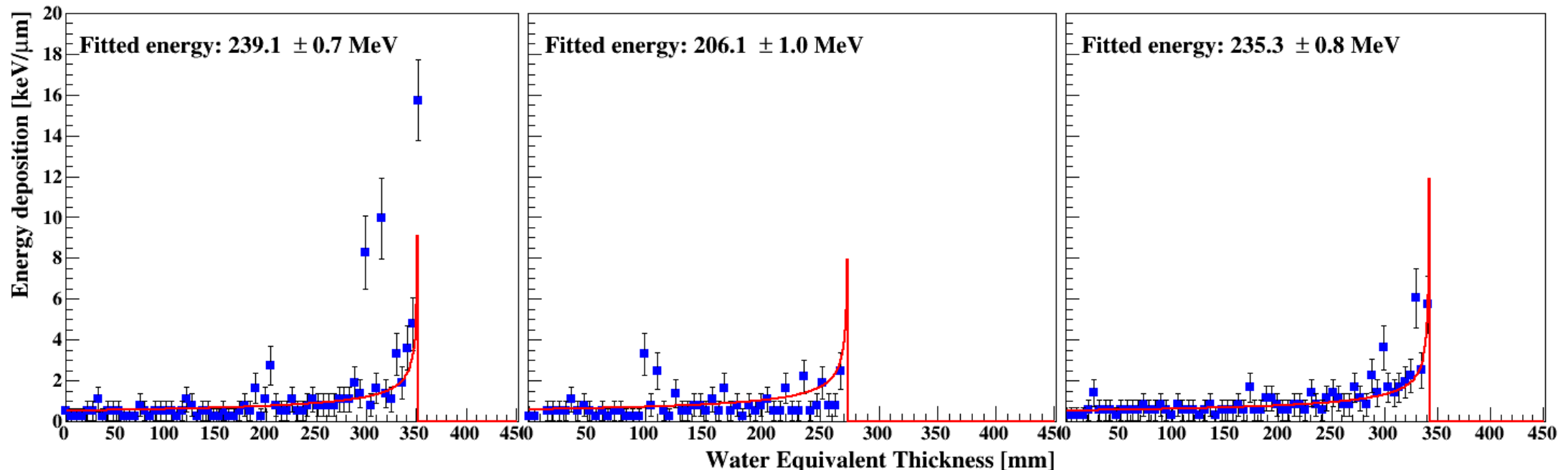
Points to discuss

1. Approaches to the analysis
2. Resolution determination
3. Features we **understand** and **don't understand**

Approaches to the analysis

1. How to properly fit the Bragg Peak?
 - Choice of limits (currently: 0 to last recorded hit + 1 layer)
 - Choice of bias (currently: Start of range fit is last recorded hit + 0.5 layer)
 - Choice of variables to fit (currently: Only range. More spread out if scale is included)

Bragg-Kleeman fit to exp. data at 237 MeV

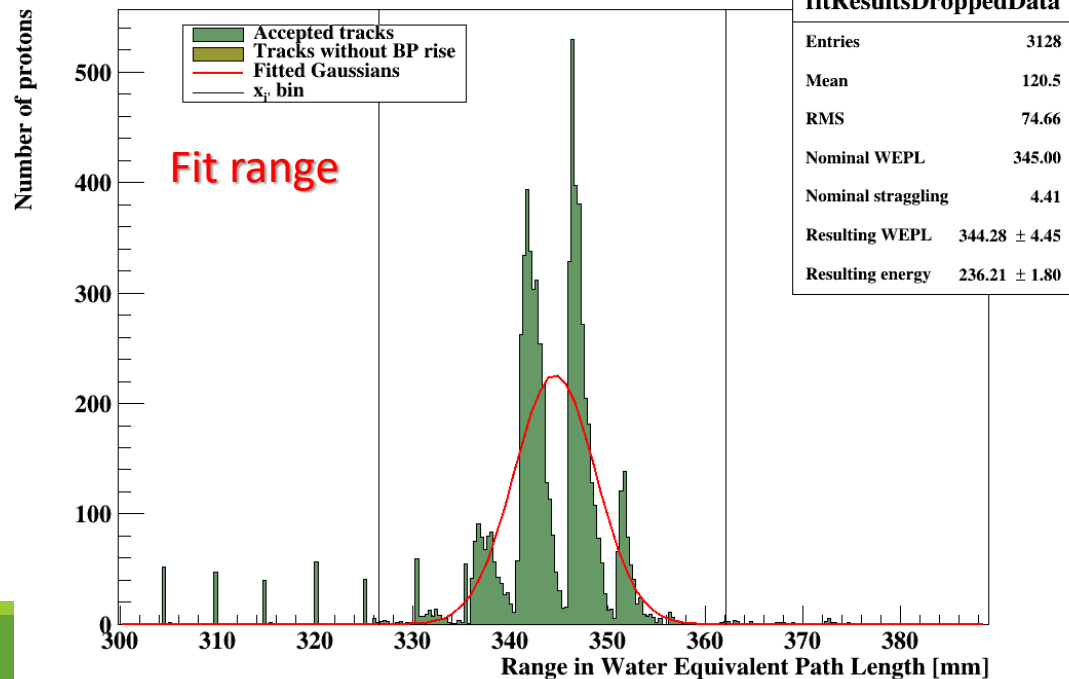


Approaches to the analysis

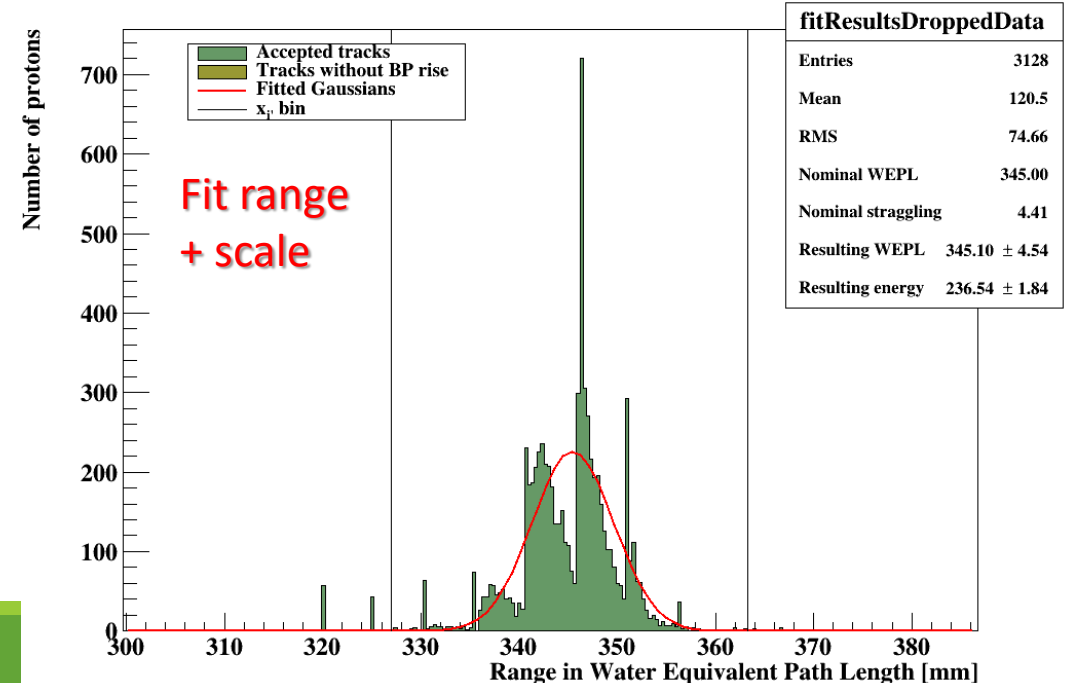
1. How to properly fit the Bragg Peak?

- Choice of limits (currently: 0 to last recorded hit + 1 layer)
- Choice of bias (currently: Start of range fit is last recorded hit + 0.5 layer)
- **Choice of variables to fit (currently: Only range. More spread out if scale is included)**

Fitted energy of a 250 MeV nominal beam on Aluminium DTC w/34.0 mm water degrader



Fitted energy of a 250 MeV nominal beam on Aluminium DTC w/34.0 mm water degrader

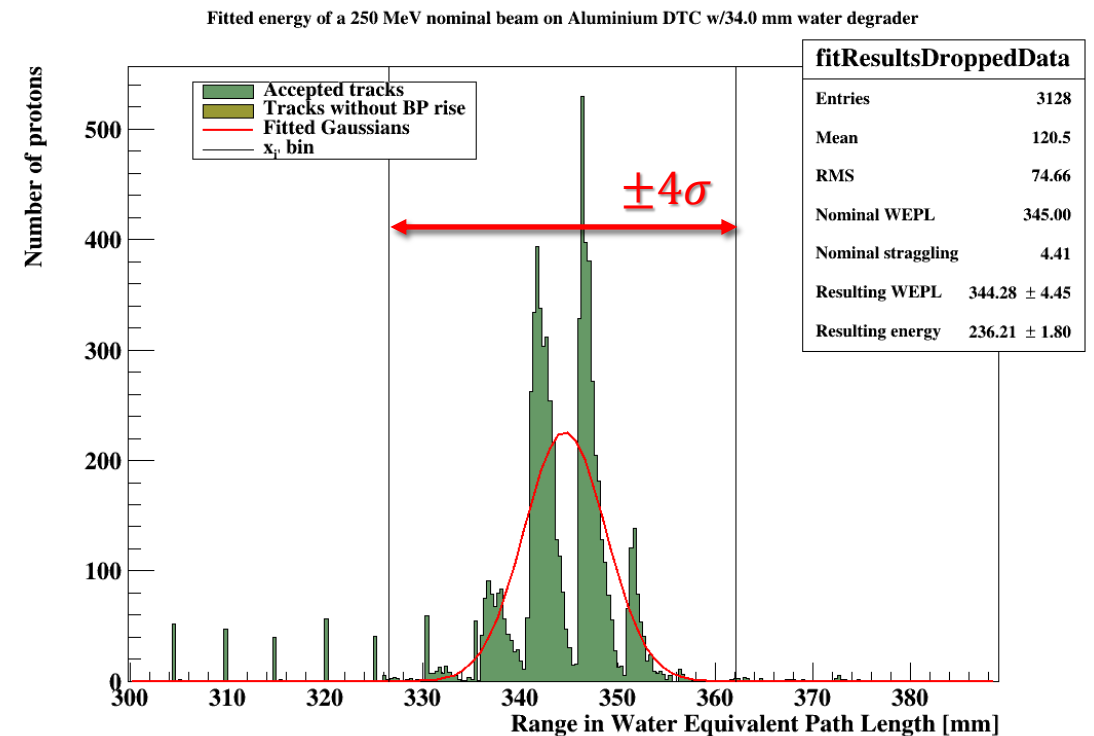


Approaches to the analysis

2. How to determine Mean and Width of range distribution?
 1. Gaussian fit parameters (NO) or empirical summed values (YES)

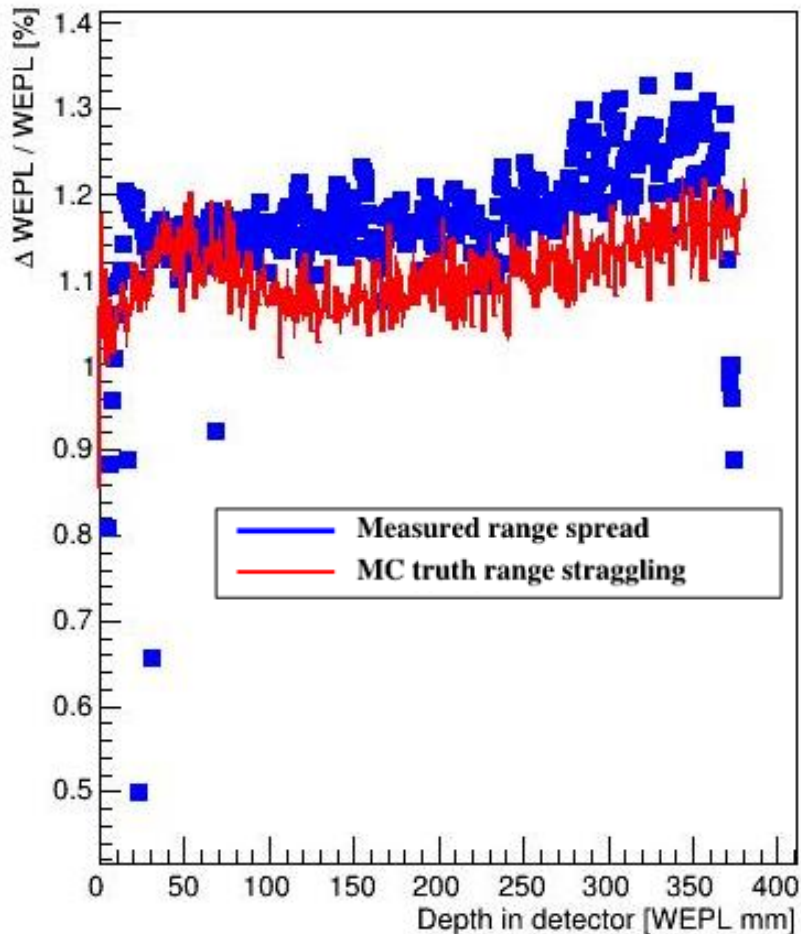
$$[\widehat{R}_0]_A = \sum_{-4\sigma}^{+4\sigma} \frac{R_i w_i}{w_i}$$
$$[\widehat{R}_0]_B = \text{Gauss}_\mu$$

2. Limits of empirical summed values?
(currently: $\pm 4\sigma$ from fitted Gaussian)
3. Q: How many layers in one «straggling» to best find the Range? 4-5? Better than FoCal's 1-2.



Resolution determination

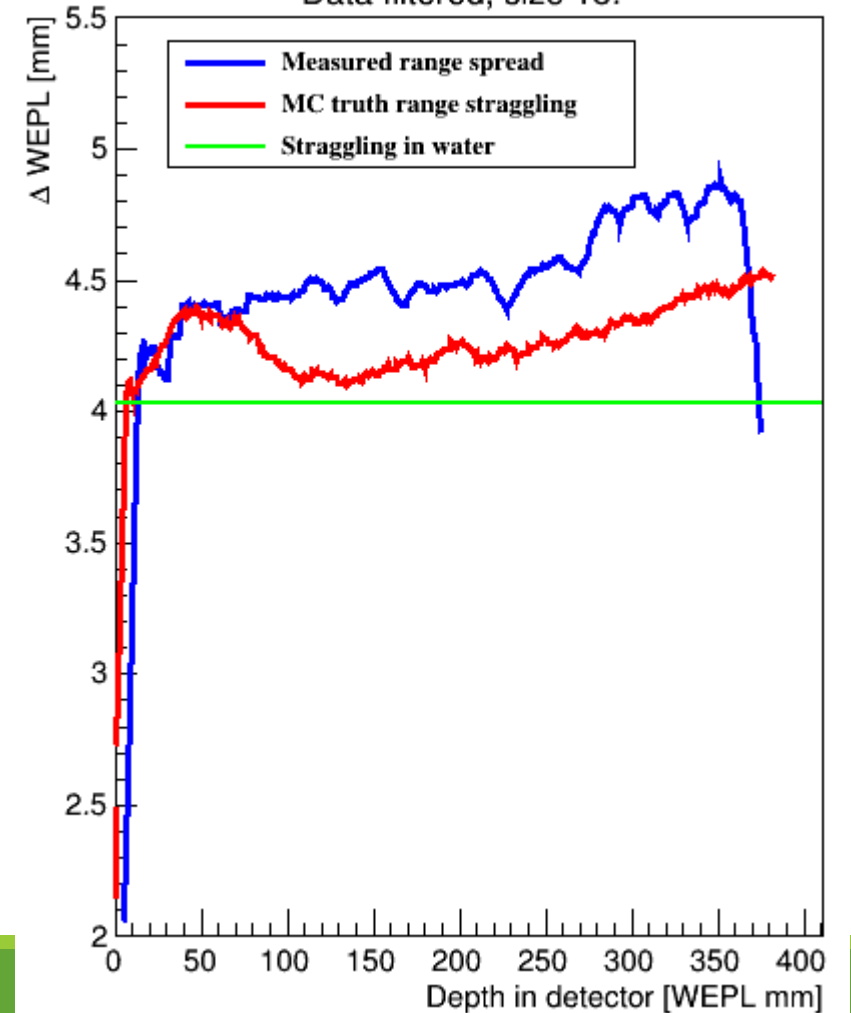
WEPL resolution using 3 mm Al absorber



TOO NOISY!!

```
for (Int_t i=0; i<arraySize; i++) {  
    value = 0;  
    n = 0;  
    for (Int_t j=i-filterSize/2; j<=i+filterSize/2; j++) {  
        if (j<0 || j>=arraySize) {  
            continue;  
        }  
        value += array[j];  
        n++;  
    }  
    value /= n;  
    tempArray[i] = value;  
}
```

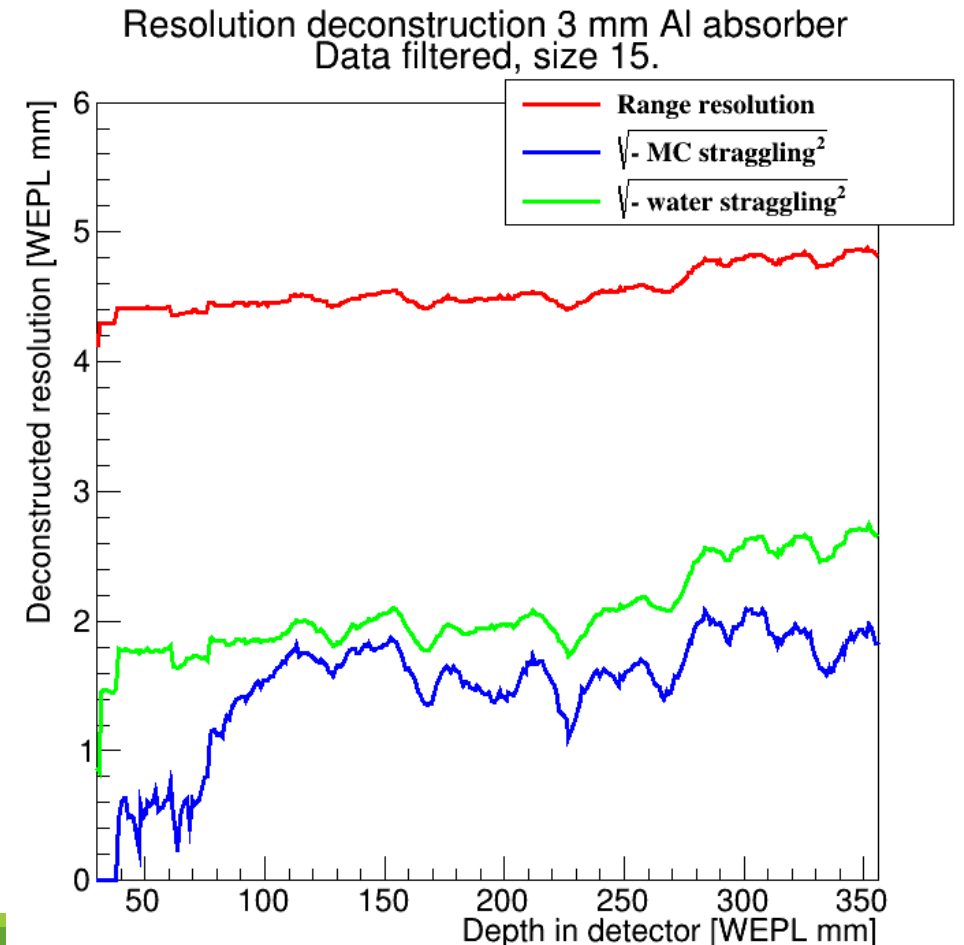
WEPL resolution using 3 mm Al absorber
Data filtered, size 15.



Resolution determination

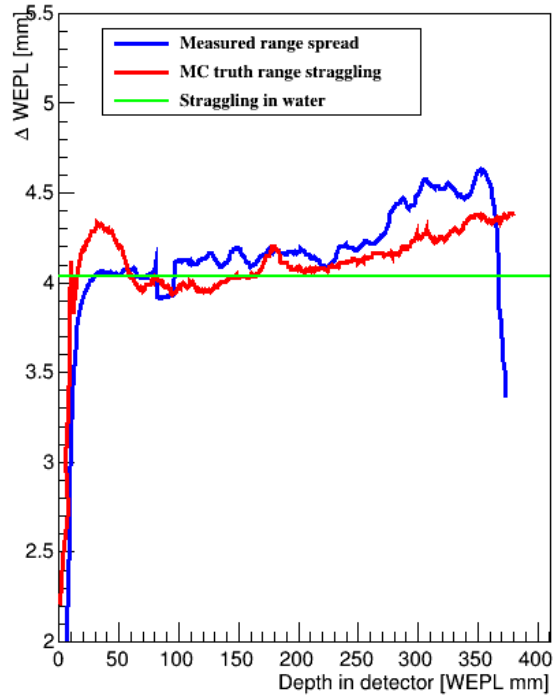
So... Lets do some calculations

1. Resolution can be added quadratically
2. We can remove the different (known) components by removing also in quadrature
3. Let's stay in the stable region of the detector (30 – 360 mm)

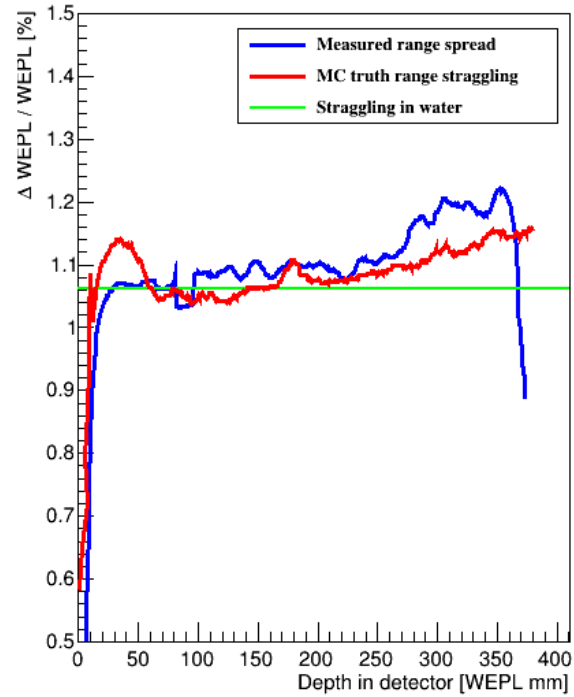


For the different geometries

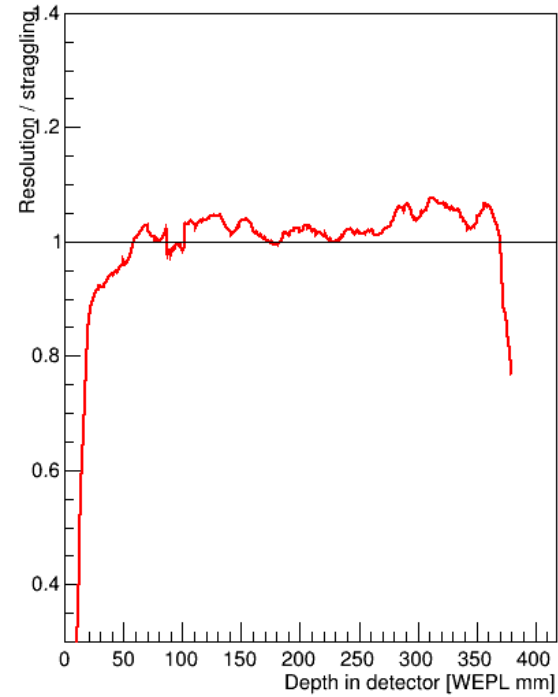
WEPL resolution using 2 mm Al absorber
Data filtered, size 15.



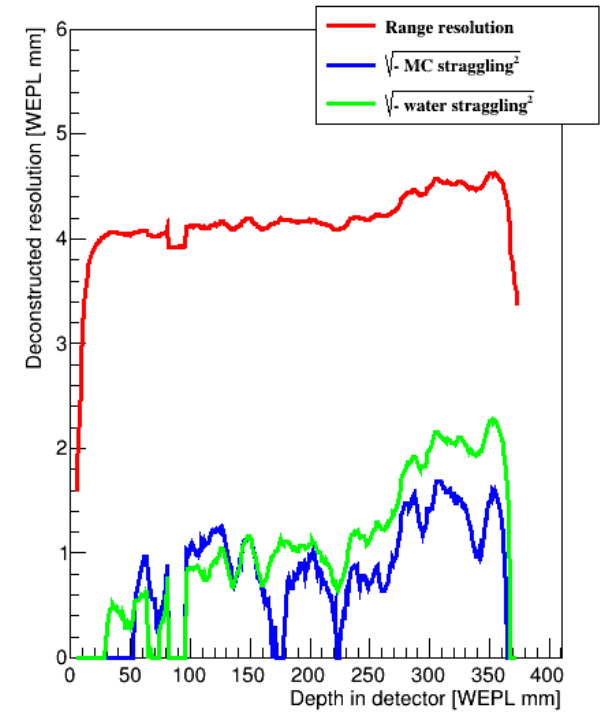
WEPL resolution using 2 mm Al absorber
Data filtered, size 15.



Resolution/straggling, 2 mm Al absorber
Data filtered, size 15.

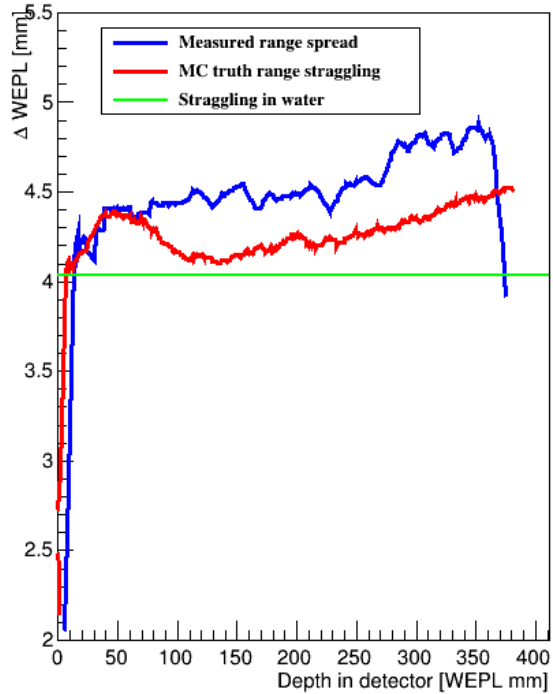


Resolution deconstruction 2 mm Al absorber
Data filtered, size 15.

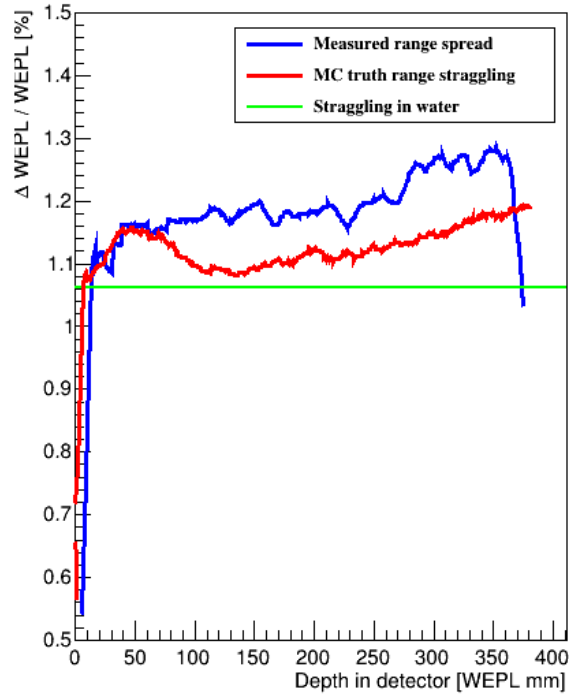


For the different geometries

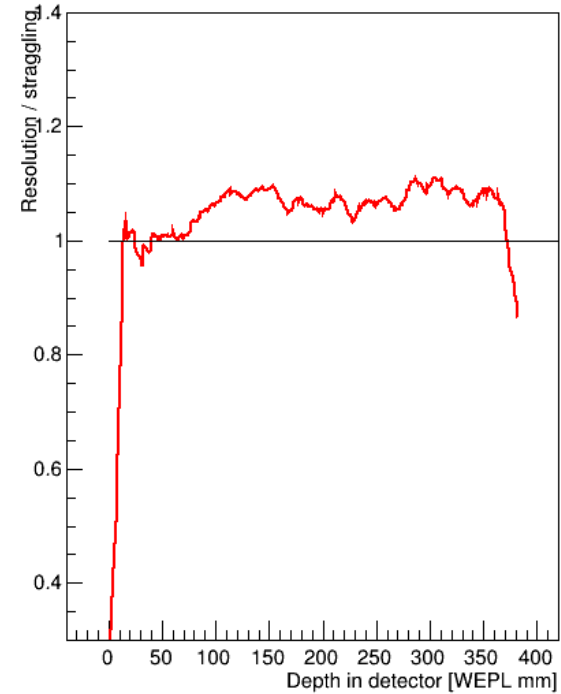
WEPL resolution using 3 mm Al absorber
Data filtered, size 15.



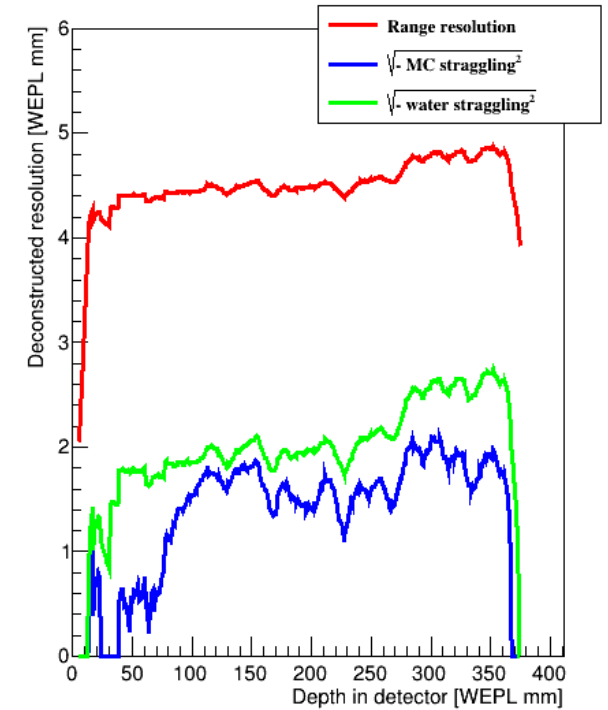
WEPL resolution using 3 mm Al absorber
Data filtered, size 15.



Resolution/straggling, 3 mm Al absorber
Data filtered, size 15.

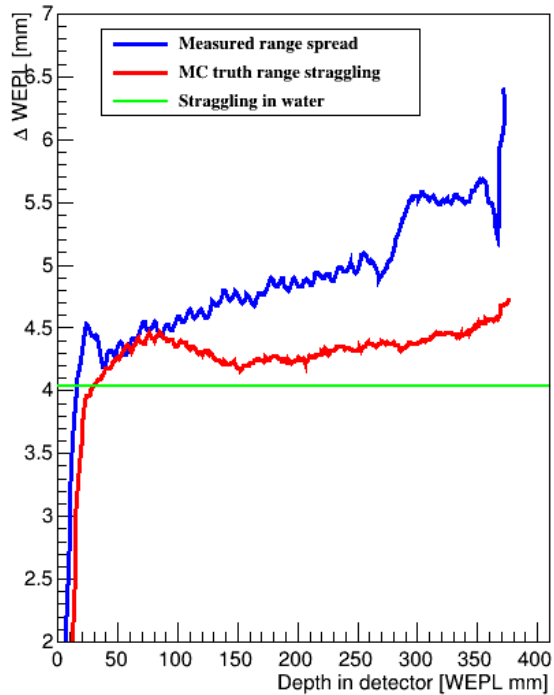


Resolution deconstruction 3 mm Al absorber
Data filtered, size 15.

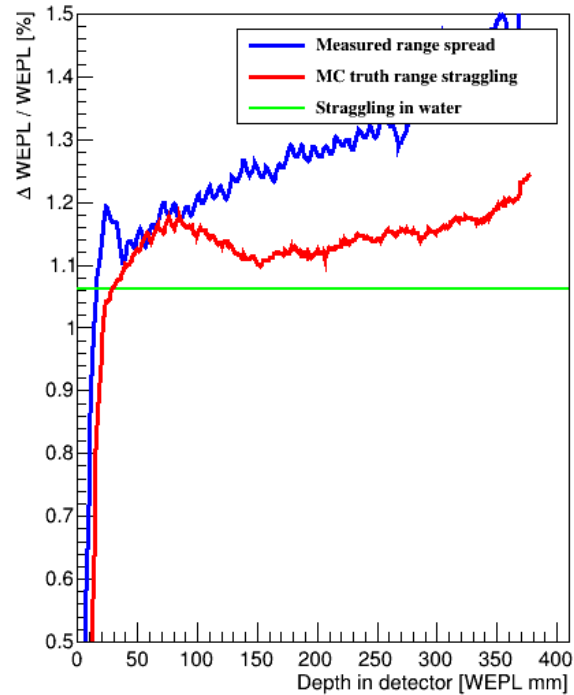


For the different geometries

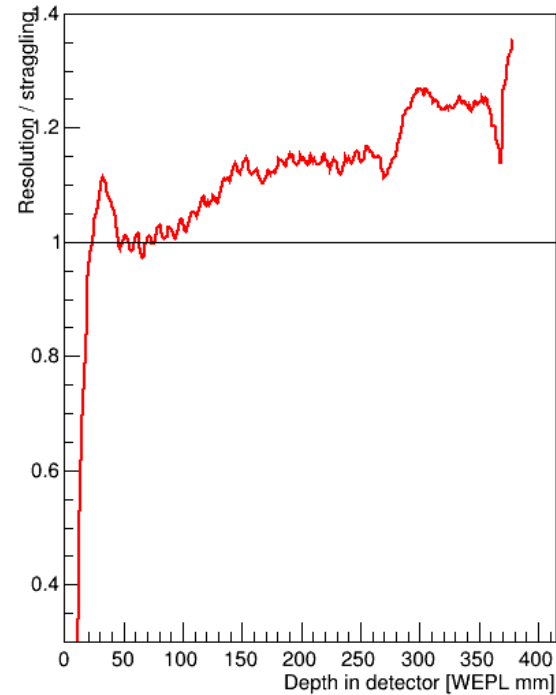
WEPL resolution using 4 mm Al absorber
Data filtered, size 15.



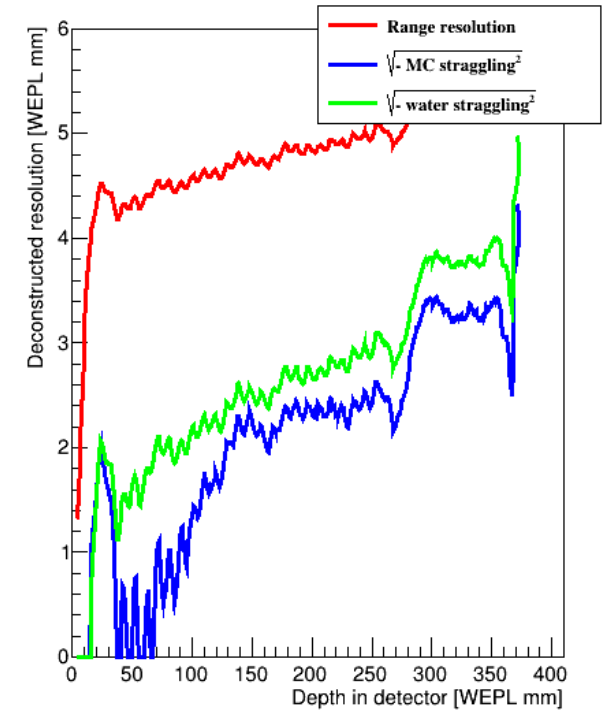
WEPL resolution using 4 mm Al absorber
Data filtered, size 15.



Resolution/straggling, 4 mm Al absorber
Data filtered, size 15.



Resolution deconstruction 4 mm Al absorber
Data filtered, size 15.



NOTE: This time the DTC had too few layers, so the deepest results are not OK

Resolution determination

So... Lets do some calculations

Setup	Average Resolution	Res. \ominus MC strag.	Res. \ominus H ₂ O strag.
2 mm	4.20 mm (1.11%)	1.03 mm	0.76 mm
3 mm	4.55 mm (1.19%)	1.37 mm	1.98 mm
4 mm	4.89 mm (1.29%)	2.10 mm	2.67 mm
Loma Linda (@ 200 MeV)	4.1 mm (2.19%)?		3.00 mm (1.16%)?
FOCAL @ 188 MeV	9.6 mm (4%) !!		

Remember...

For the discretization uncertainty to be sub-dominant to range straggling, we would require $\Delta/\sqrt{12} < 3$ mm (Table 4).

The WEPL factor is approx. 2.18

At 2 mm: $4.4 / \sqrt{12} = 1.27$ ✓

At 3 mm: $6.54/\sqrt{12} = 1.89$ ✓

At 4 mm: $8.72/\sqrt{12} = 2.51$ ✓

At 5 mm: $10.9/\sqrt{12} = 3.14$ ✗

$$\underline{3 * \sqrt{12} / 2.18 = 4.76}$$

Interpreting the Loma Linda results...

IV. FIVE-STAGE DETECTOR PERFORMANCE

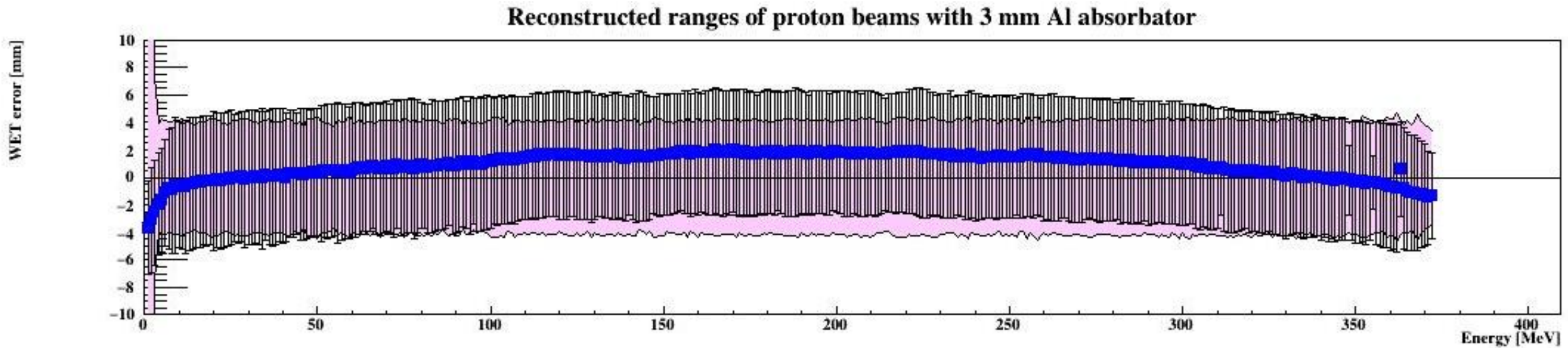
When we look at just the signals in the first stage from 200 MeV protons passing through a 2 cm square in the center of the 5-stage detector (with no phantom in place), we find a Gaussian distribution with $\sigma = 3.0\%$. Since Geant-4 simulations [18] predict $\sigma = 2.8\%$ just from variations in the energy deposition, by subtracting the two numbers in quadrature we estimate that the detector resolution is around 1%, more than adequate for this application.

in 6.35 mm steps. From those data we derive the calibration constants needed for scanner operation. We have verified that this procedure yields a WEPL resolution that is approximately 3 mm rms for all values of WEPL, only slightly higher than the predicted lower limit of 2.8 mm that arises from range straggling of 200 MeV protons.

200 MeV -> 259 mm
3 mm is 1.16 %

BUT... Is that after subtracting the expected straggling..? Do they do that twice??

Features we don't understand



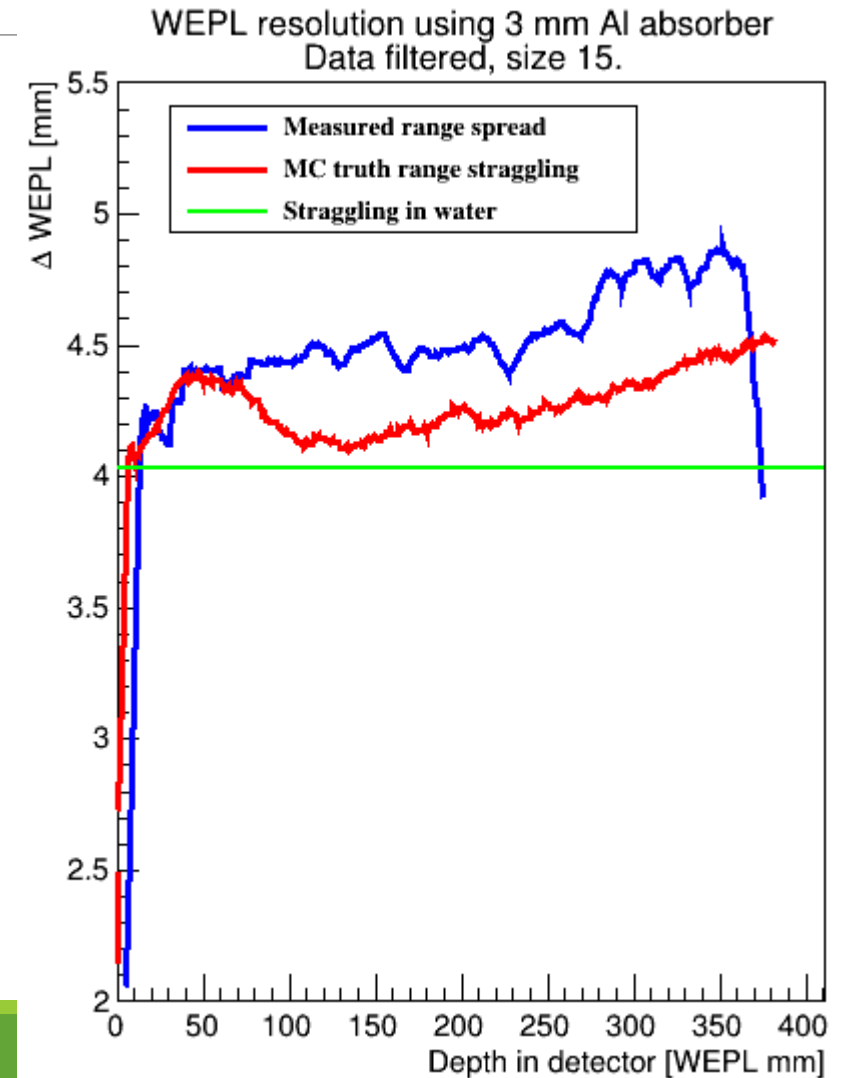
Why do we have these «nice» systematic errors in range determination?

- Looks like Bragg-Kleeman-type error of range-energy, but I can't think of any place I use the equation... (I use LUTs to calculate this)

Features we don't understand

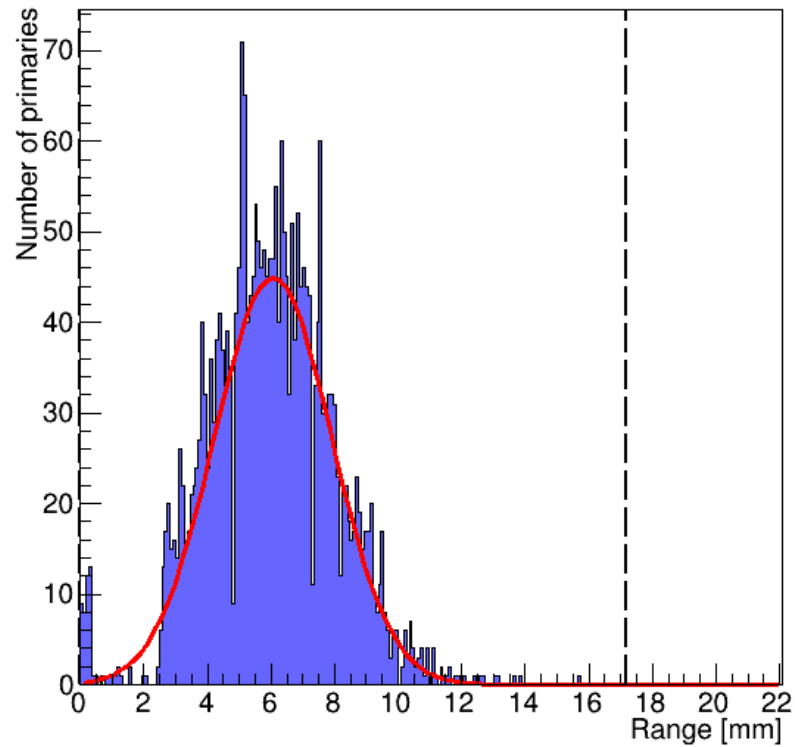
What about the «bump» in the MC truth range straggling here? It's in the raw data as well

1. Edge effects from the proton beam hitting the DTC just before the Bragg Peak?
2. It's not mis-fits (I've studied the range distribution data)
3. We expect the MC truth range straggling to start at the water straggling levels (since beam passes through 99% water at that point), and increase when the fraction of the beam traversing the DTC increases.
4. I'll try anchoring the water phantom just proximal to the first layer, and not to its midpoint

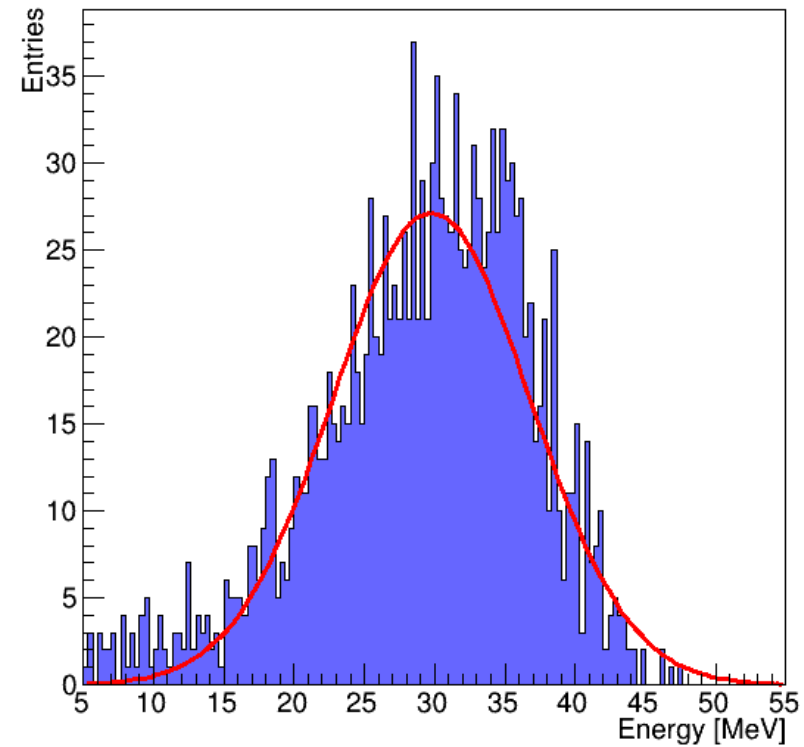


Animation of the straggling raw data

Projected range in DTC



Remaining energy after degrader



Bump at 5-30 mm with peak at 18 mm

NOTE: It's the Gaussian σ that is used, not a histogram sum

Going forward

1. We need to fully understand the artifacts before going through with a full scan of all geometries and materials, otherwise it's wasted effort
2. All code is located at github: <http://github.com/HelgeEgil/focal>
3. A full user's guide for its usage is located on the wiki: [https://wiki.uib.no/pct/index.php/Software for design optimization](https://wiki.uib.no/pct/index.php/Software_for_design_optimization)

Known bugs

1. It's still not possible to use geometries with floating point values for absorber thicknesses
2. The DTC geometry creator does NOT propagate the first layer material. We want to use Air (to increase the Dynamic Range at low energies), but it defaults to Aluminium...
 - Solution: Edit the material in *Module.mac* after using *makeGeometryDTC.py*.