

Equalizing the response of the FOOT Calorimeter as a function of the ion energy and charge



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on behalf of the FOOT collaboration

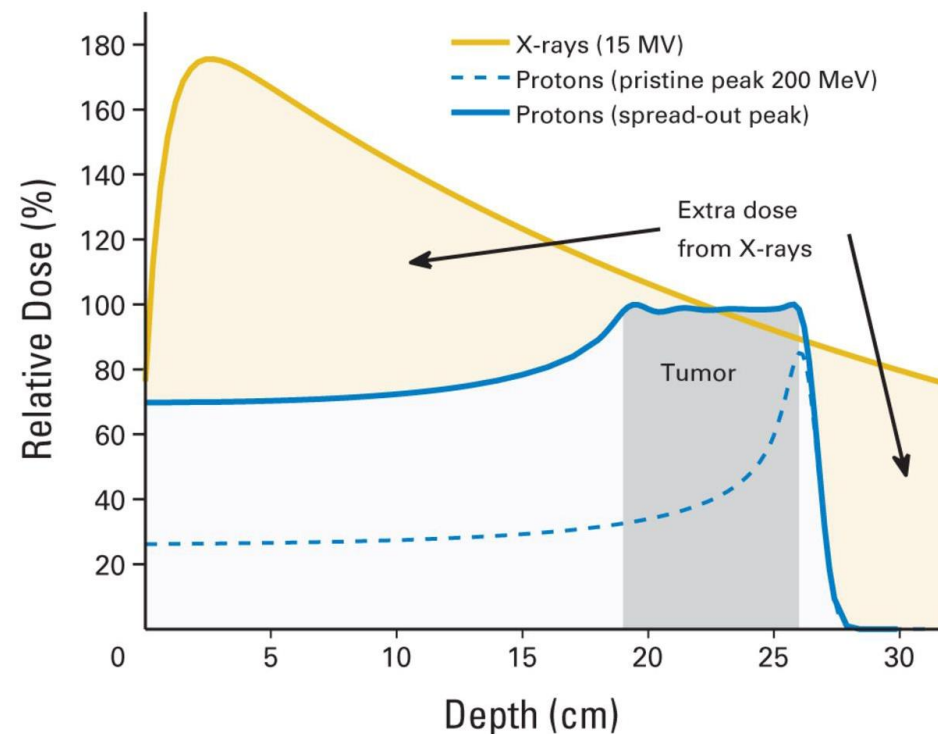
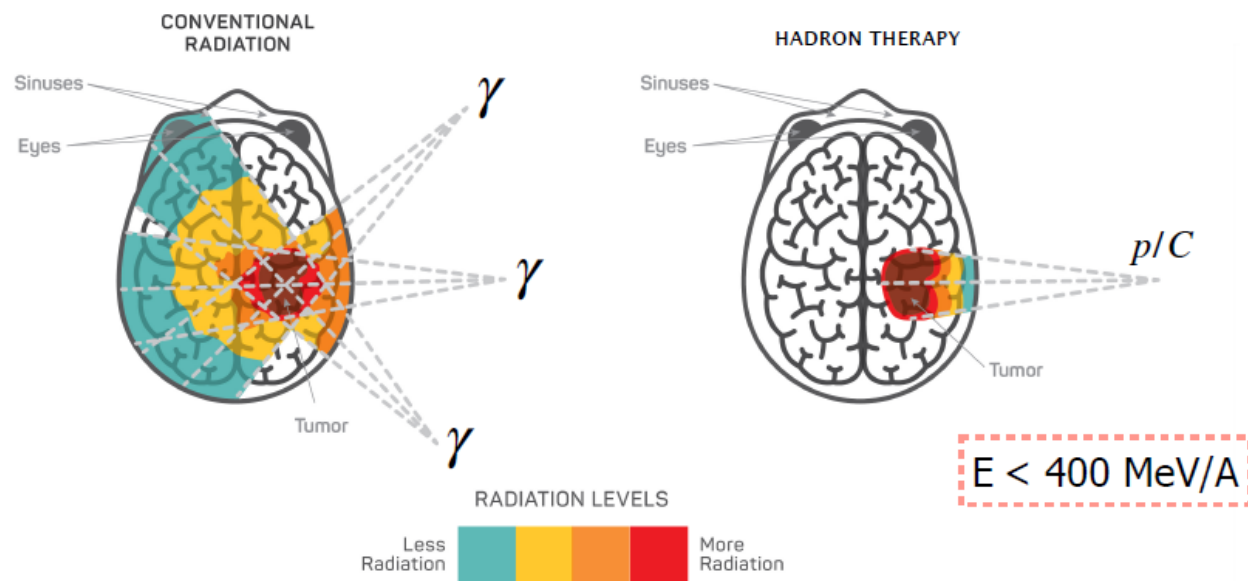
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Particle Therapy

- ✓ Particle therapy is performed with p or C ions
- ✓ The Bragg peak is very effective in minimizing the dose delivered to healthy tissues



Secondary fragments are created in the interaction between treatment beams and tissues



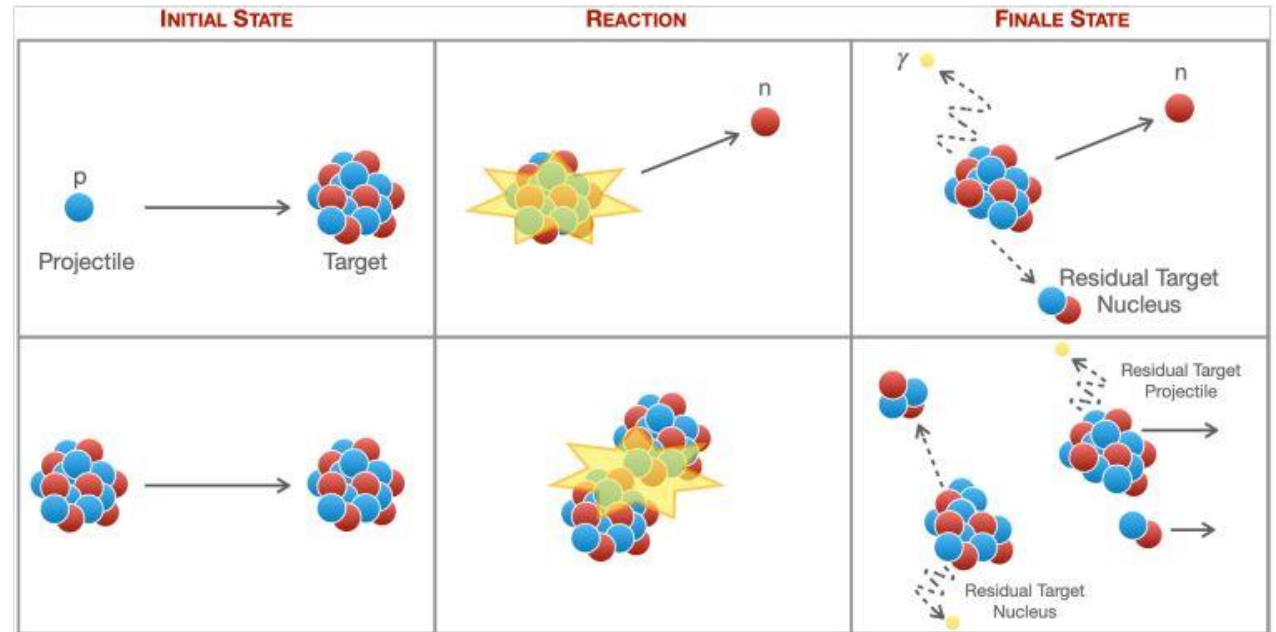
Nuclear Fragmentation

Target Fragmentation:

- ✓ Target fragments are produced with low energies (short range, hundreds of μm)
- ✓ Difficult to detect

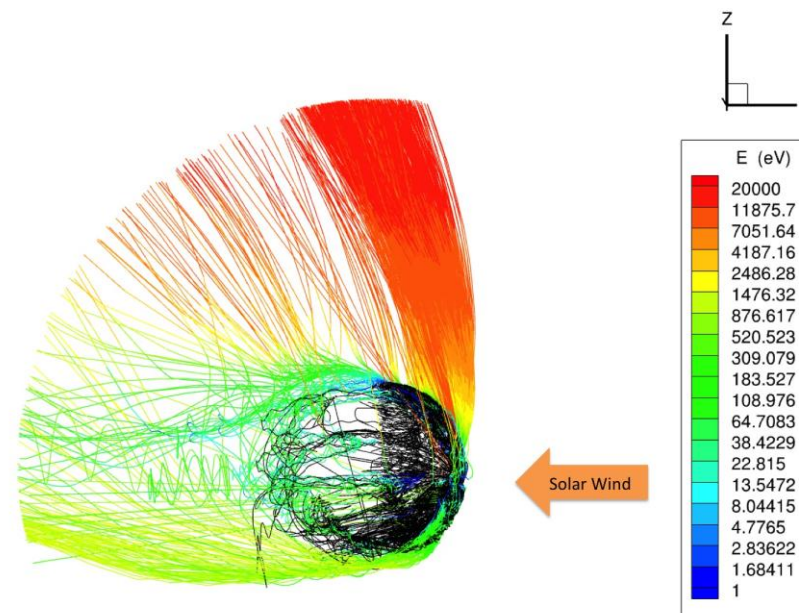
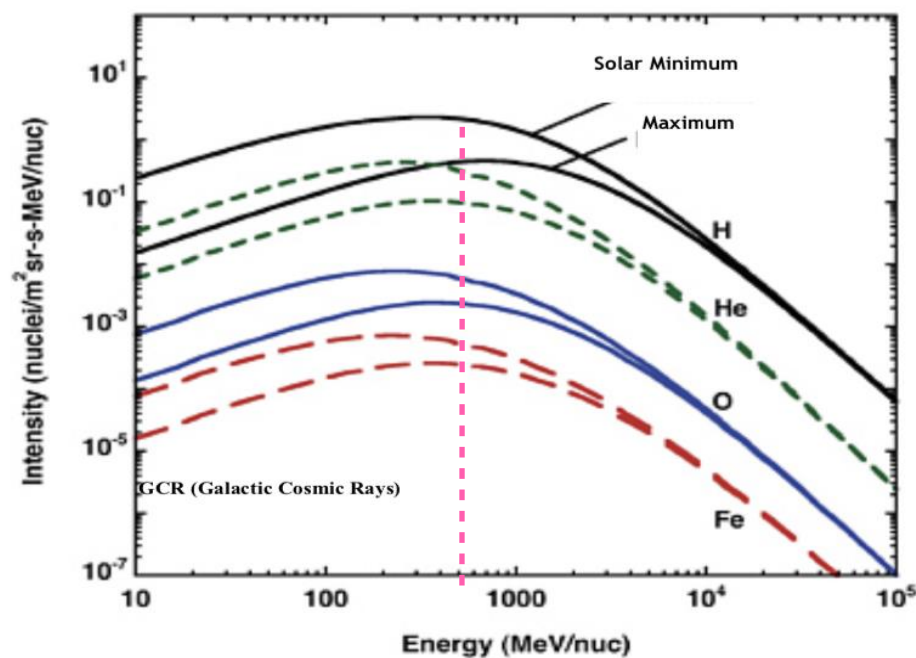
Beam Fragmentation:

- ✓ Projectile fragments (longer range)
- ✓ Non-zero dose beyond the Bragg peak to address
- ✓ Not present in protontherapy



Spacecraft Shielding

- ✓ Charged particles in space: Solar Particles Events (SPEs), Galactic Cosmic Rays (GCRs), geomagnetically trapped particles
- ✓ Interaction with walls/shielding of spacecraft produce secondary fragments



FragmentatiOn Of Target: FOOT

- ✓ Aim: measurement of beam and target fragmentation differential cross sections with 5% accuracy
- ✓ Goal: charge and mass identification at 2-3% and 5% accuracy, respectively
- ✓ Inverse kinematics approaches
- ✓ ^4He , ^{12}C , ^{16}O beams of 200-400 MeV/u on ^{12}C , C_2H_4 and $\text{C}_5\text{O}_2\text{H}_8$ targets (Hadrontherapy)
- ✓ ^{12}C , ^{16}O beams of 800 MeV/u on ^{12}C , C_2H_4 and $\text{C}_5\text{O}_2\text{H}_8$ targets (Radioprotection)

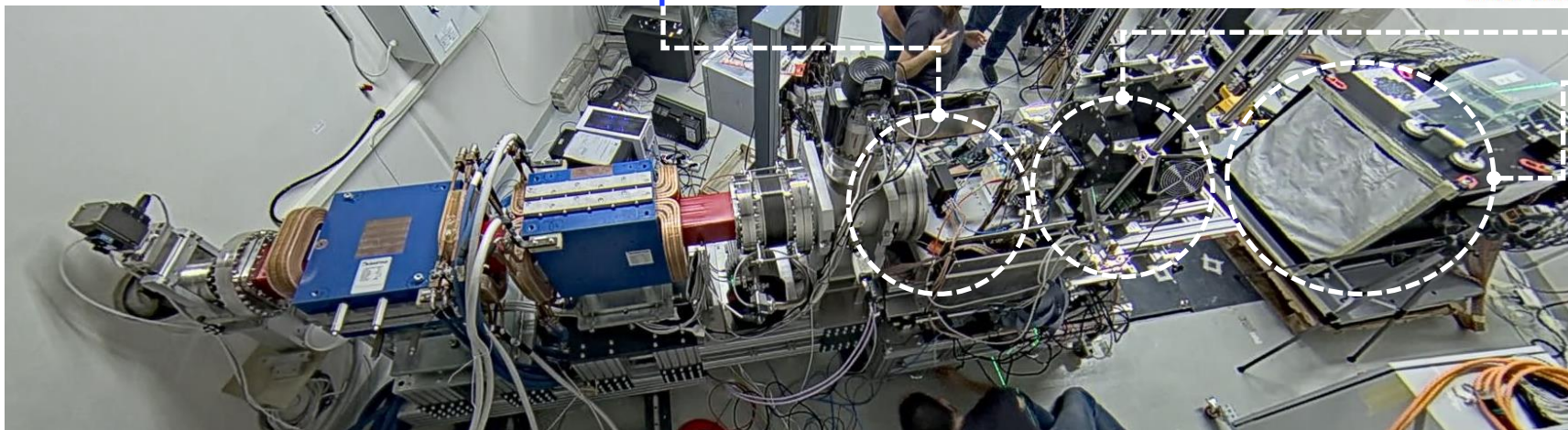
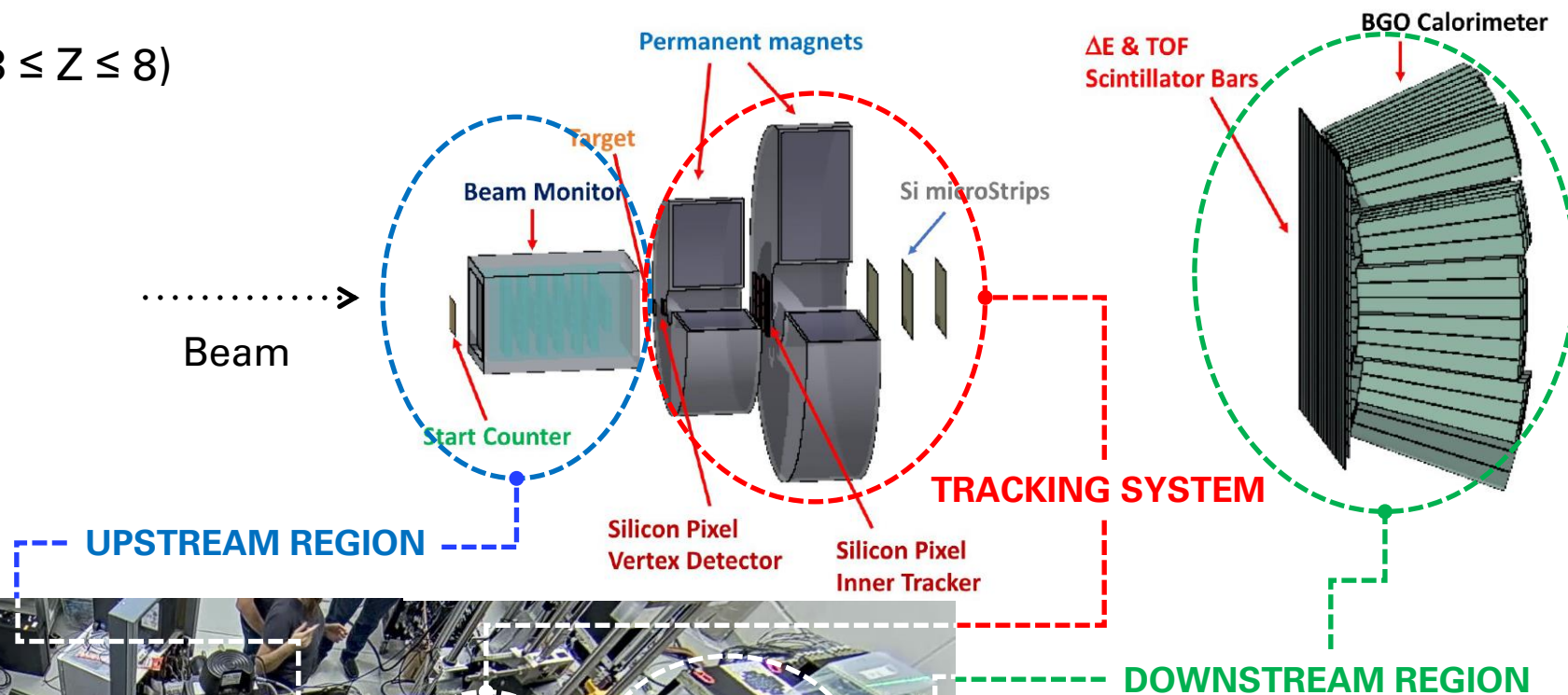
FOOT is a collaboration with about 100 members coming from

- INFN: 10 sections
- 3 laboratories: CNAO, GSI, IPHC
- 15 universities: France, Italy, Japan, Germany



Electronic Setup

- ✓ Designed for fragments ($3 \leq Z \leq 8$)
- ✓ Angular acceptance $< 10^\circ$



Mass Reconstruction

TOF (β) – TRACKER (p)

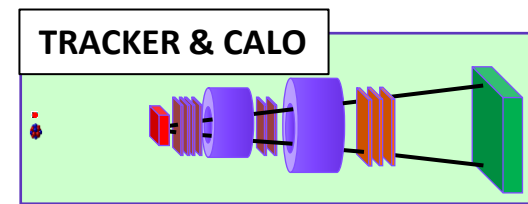
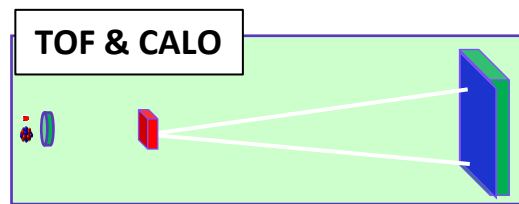
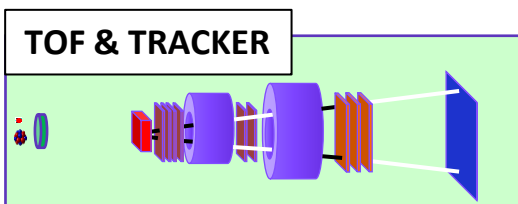
$$A_1 = \frac{m}{u} = \frac{p\sqrt{1-\beta^2}}{u\beta}$$

TOF (β) – CALO (E_{kin})

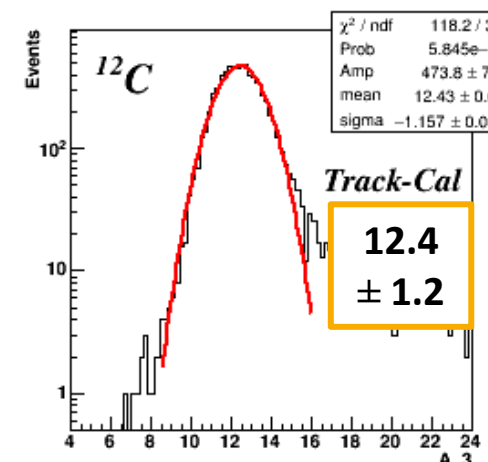
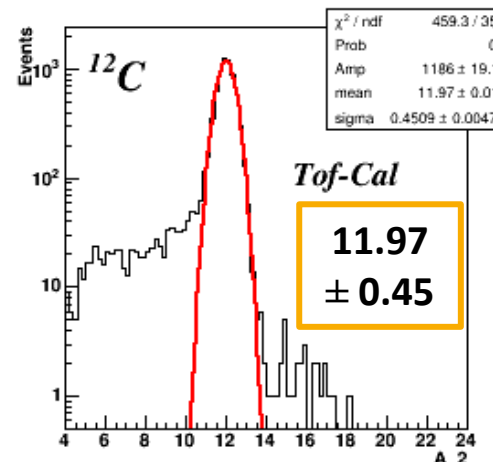
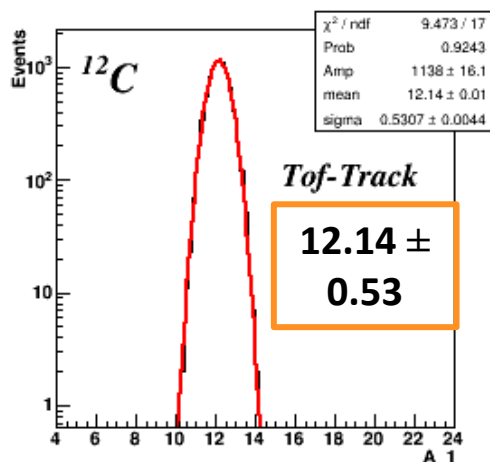
$$A_2 = \frac{m}{u} = \frac{E_{kin}}{u(\gamma-1)}$$

TRACKER (p) – CALO (E_{kin})

$$A_3 = \frac{m}{u} = \frac{p^2 - E_{kin}^2}{2E_{kin}}$$



Fluka simulation ^{16}O (200 MeV/u) + C_2H_4

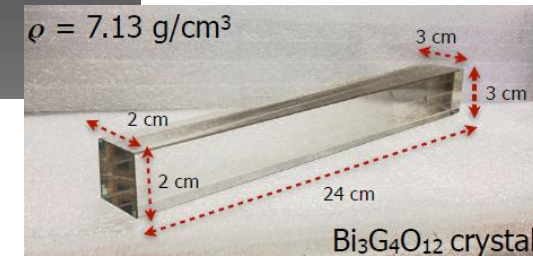
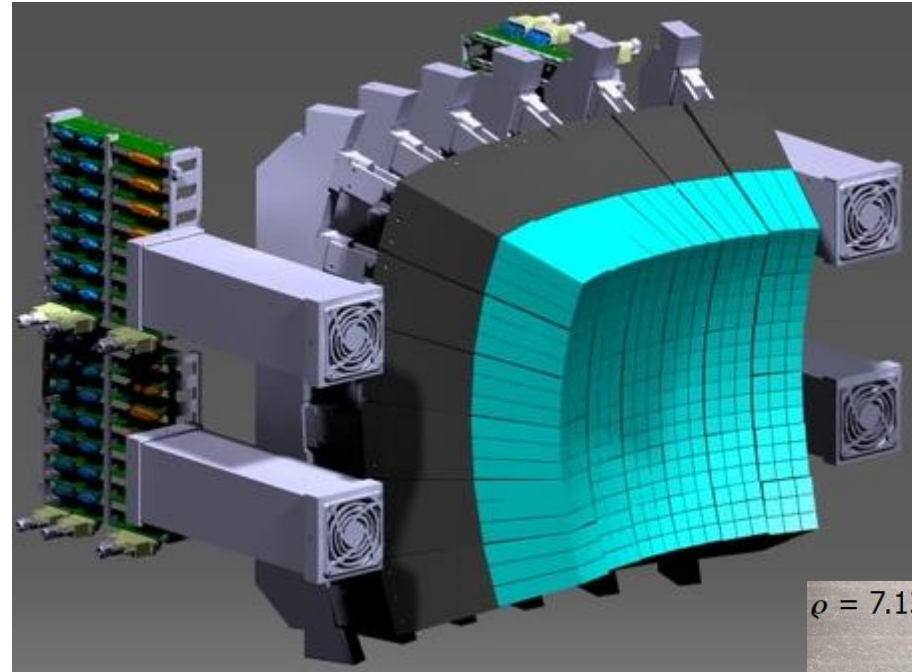


Courtesy of G. Battistoni



Calorimeter

- ✓ 320 BGO crystals grouped in modules (9 crystals for each module)
- ✓ Crystals dimension: 2x2 cm² (front) 3x3 cm² (back) 24 cm (length)
- ✓ SiPM based readout
- ✓ 36/36 modules fully assembled
- ✓ Measurement of the kinetic energy



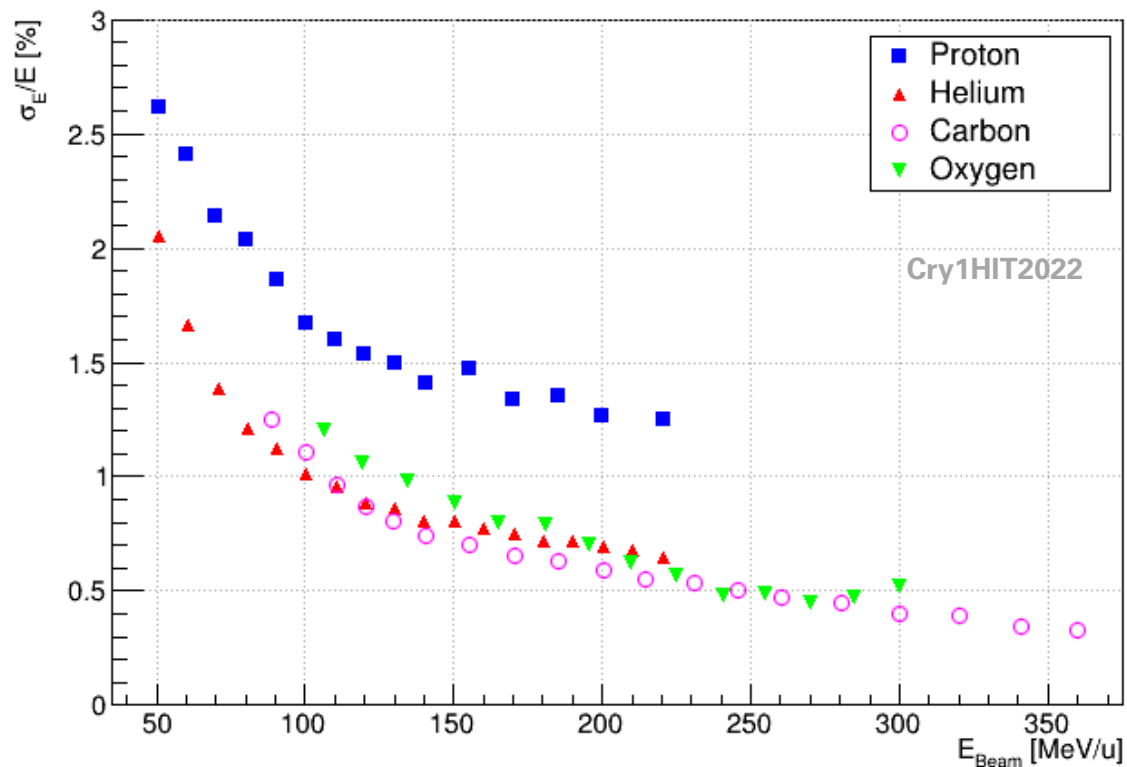
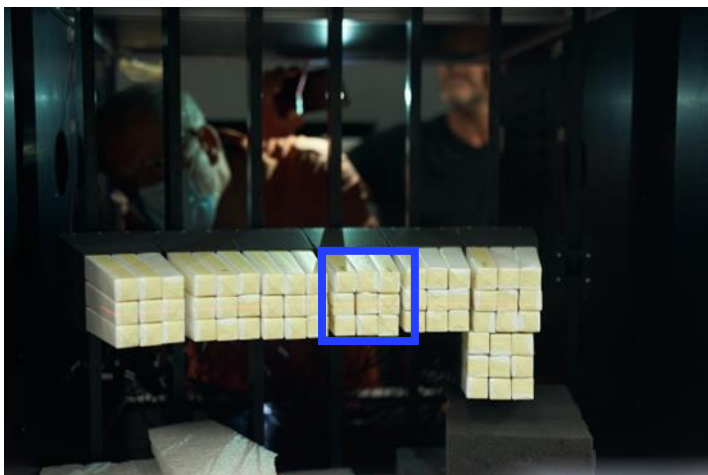
Multiple data acquisition campaigns for calibration and equalisation at:

- ✓ Heidelberg Ion Therapy Center ([HIT](#))
- ✓ Centro Nazionale di Adroterapia Oncologica ([CNAO](#))



HIT – First Energy Calibration Study

- ✓ Simple setup with just one **fully assembled module**
- ✓ Beam focused on central crystal
- ✓ No other detector between Calorimeter and beam nozzle (autotrigger)
- ✓ Energy scan from 50 to 400 MeV/u with **Proton**, **Helium**, **Carbon** and **Oxygen** ions



Calibrated crystal resolution for different ions



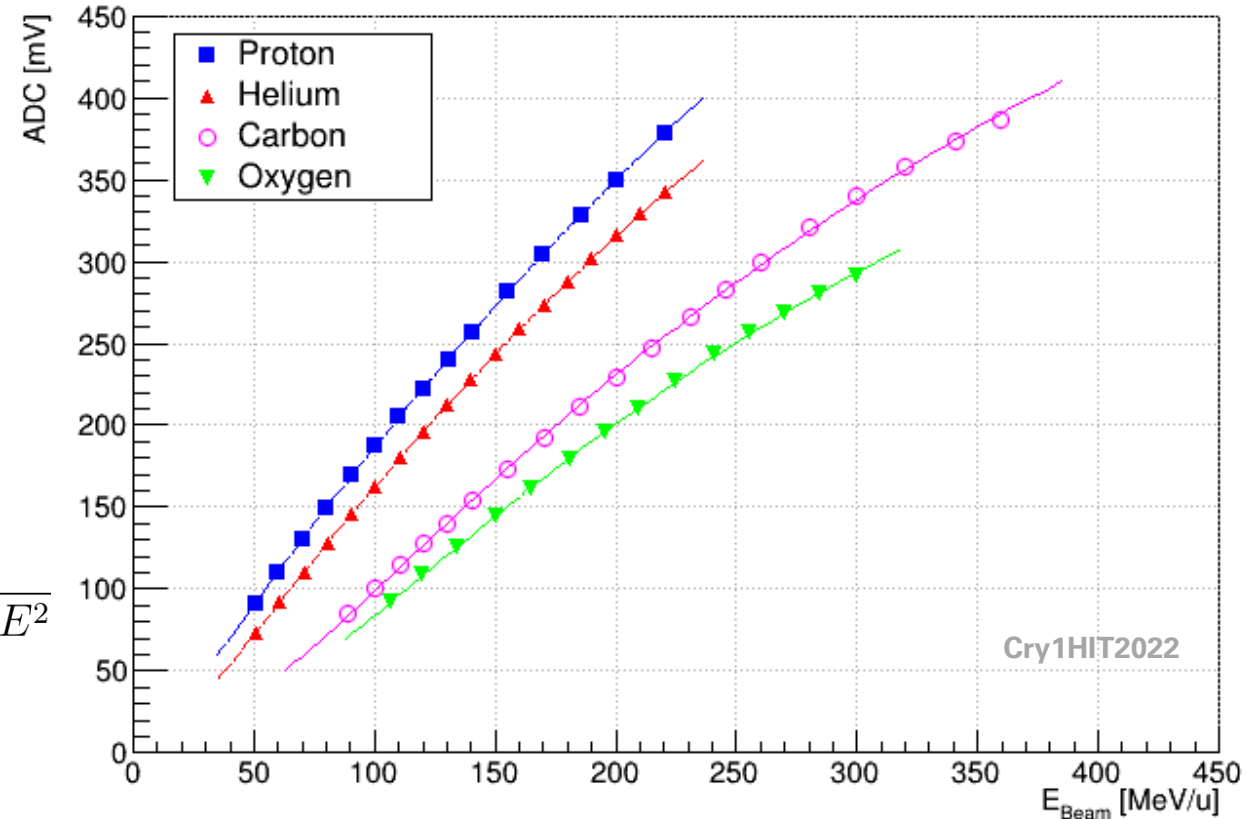
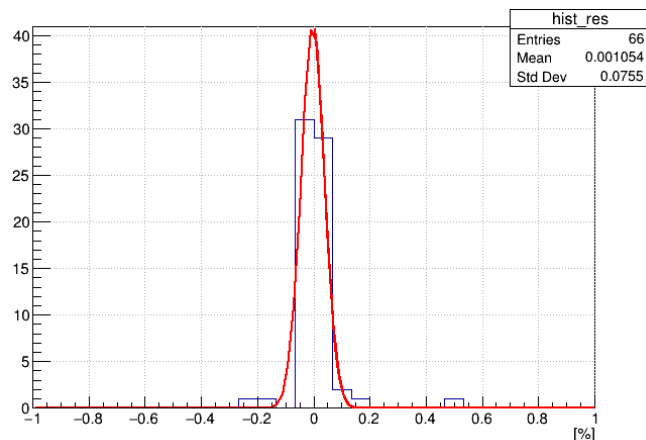
Energy Calibration Curves

- ✓ The calorimeter linearity is affected by the **Birks effect**

$$\frac{dS}{dx} = \frac{A \cdot dE/dx}{1 + k \cdot B \cdot dE/dx}$$

- ✓ There is clear dependence on **Z**
- ✓ The chosen fit function is derived from Birks formula, we call it **modified Birks function (MBF)**

$$ADC(E) = \frac{P_0 E^2}{1 + P_1 E + P_2 E^2}$$



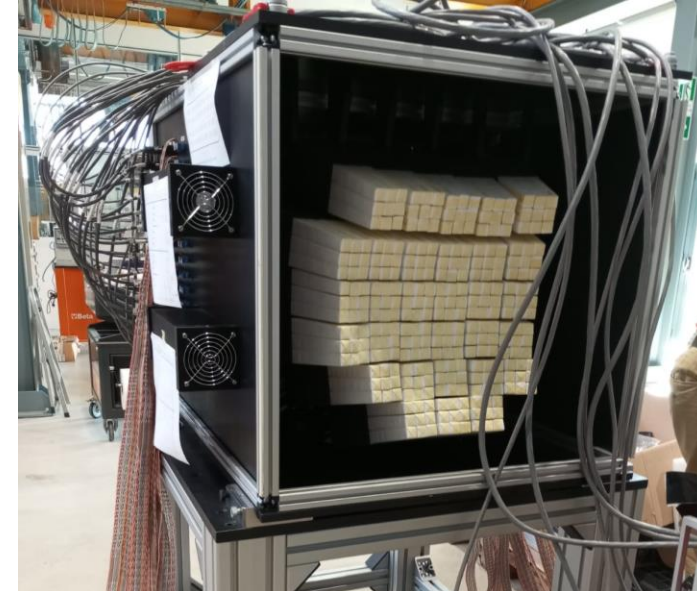
- ✓ Good fitting of experimental data

$$\frac{|E_{fit} - E_{ADC}|}{E_{fit}} < 1\%$$



CNAO – Calibration Validation

- ✓ Beam in “Screensaver Mode”: sweeping a quarter of the modules at a time during each run greatly reducing data acquisition time
- ✓ Setup with 31 fully assembled modules
- ✓ Scan at four different energies 115, 190*, 200, 250 MeV/u with Carbon ions



Crystal intercalibration is performed using the Modified Birks function

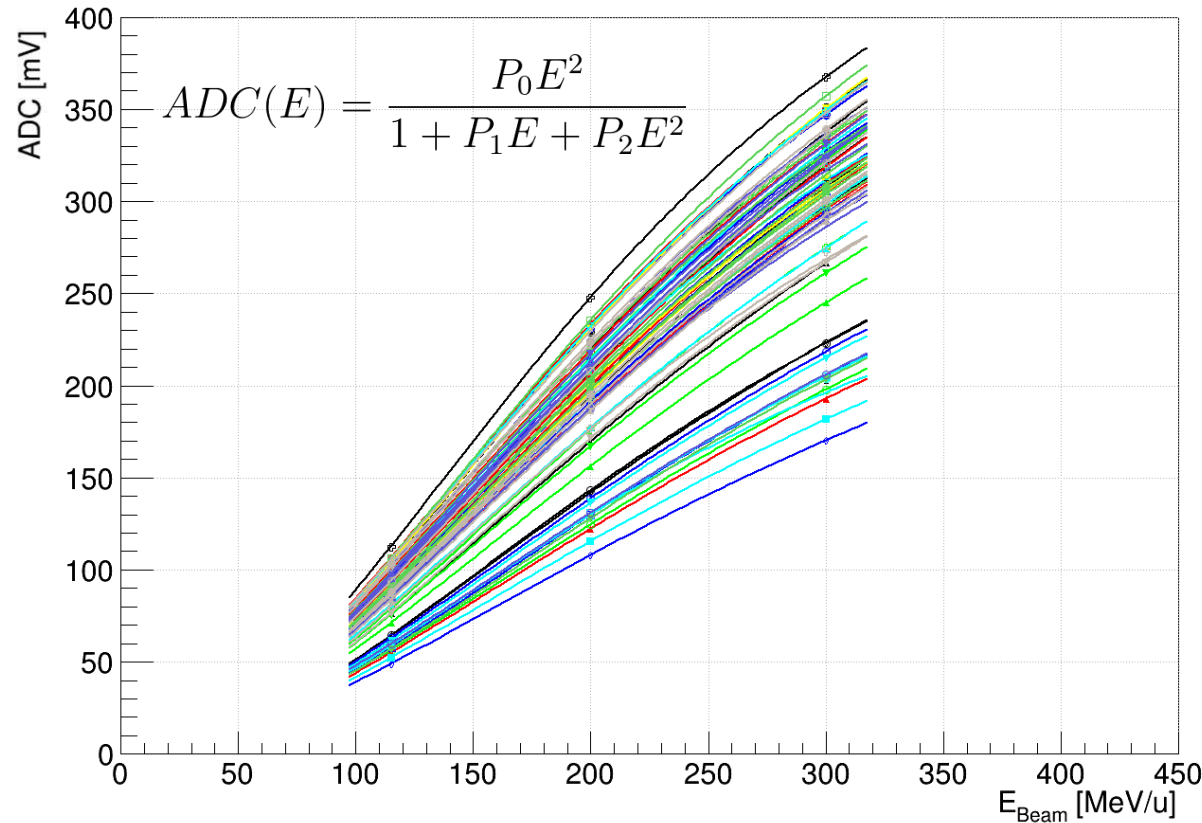
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		3	4	5	12	13	14	21	22	23										
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1		27	28	29	36	37	38	45	46	47	51	55	56	63	64	65	2			
		30	31	32	39	40	41	48	49	50	57	58	59	66	67	68				
		53	54	55	62	63	64	71	72	73	79	83	84	91	92	93				
72	73	74	81	82	83	90	91	92	99	100	101	108	109	110	117	118	119	126	127	128
75	76	77	84	85	86	93	94	95	102	103	104	111	112	113	120	121	122	129	130	131
78	79	80	87	88	89	96	97	98	105	106	107	114	115	116	123	124	125	132	133	134
135	136	137	144	145	146	153	154	155	162	163	164	171	172	173	180	181	182	189	190	191
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4			306	307	308	315	316	317	324	325	326	332	333	334				3		
			309	310	311	318	319	320	327	328	329									
			312	313	314	321	322	323	330	331	332									

* foreseen at 150 MeV, issue with beam delivery



Intercalibration Strategy

How to test whether screensaver run achieve the performance needed?



Four Carbon energy screensaver points

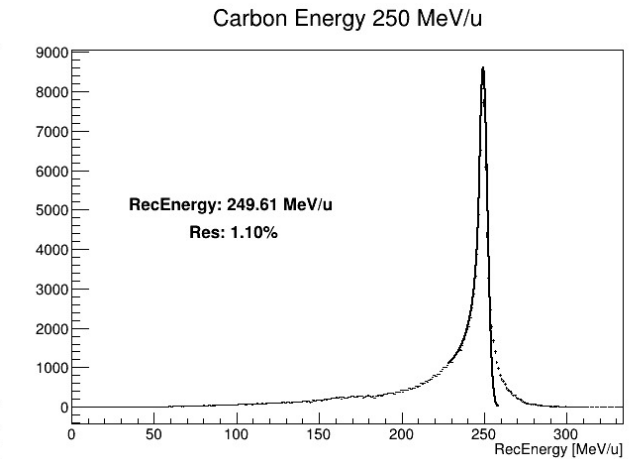
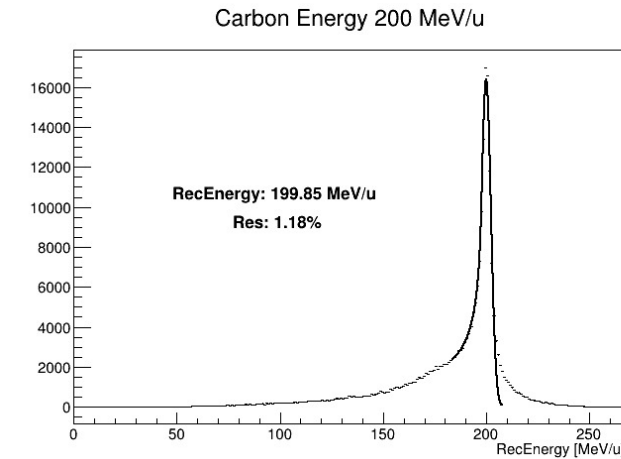
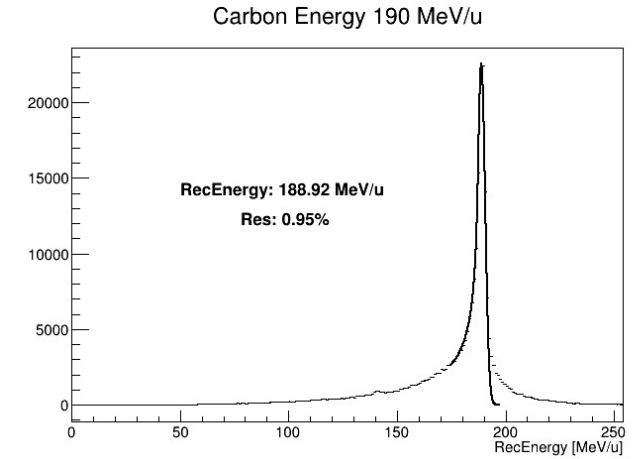
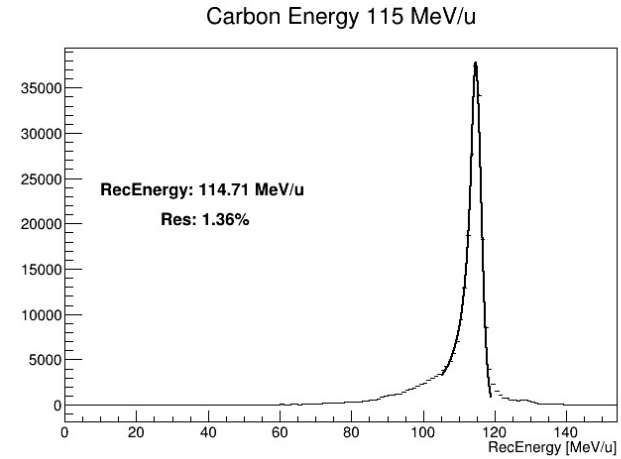
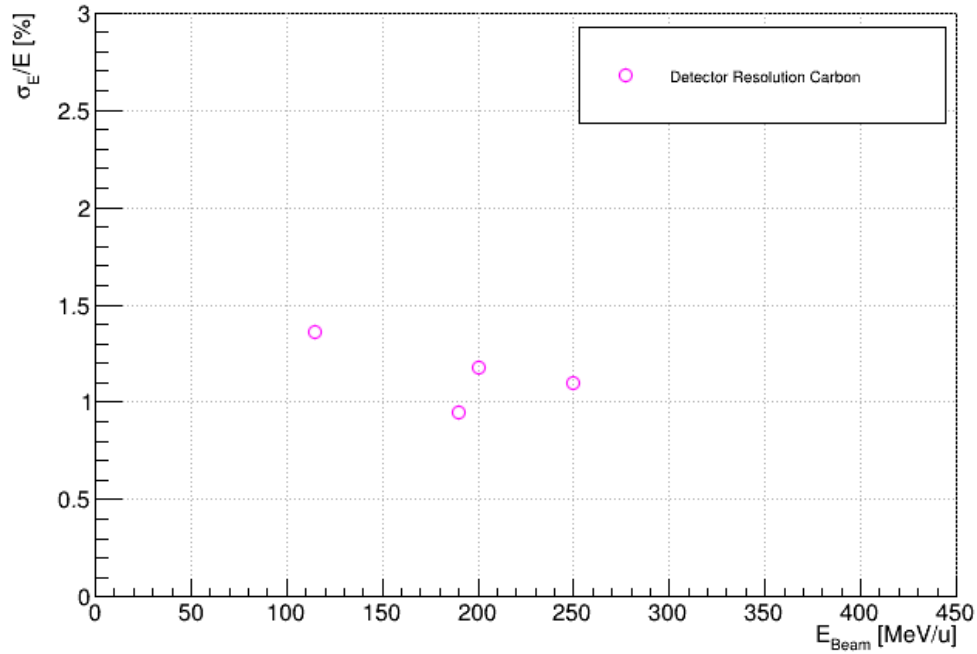
For each crystal we obtain p_0 , p_1 and p_2

Energy calculated as inverse modified Birks

$$E(ADC) = \frac{-P_1 ADC - \sqrt{(P_1 ADC)^2 - 4ADC(P_2 ADC - P_0)}}{2(P_2 ADC - P_0)}$$



Calorimeter Resolution



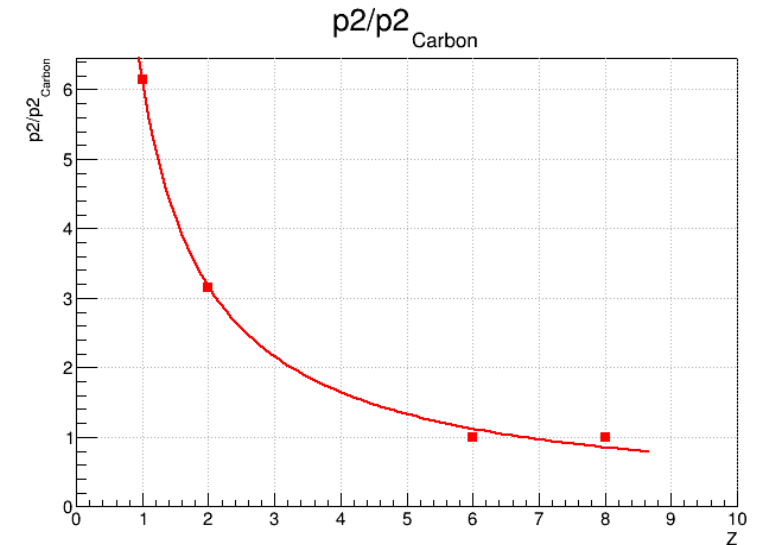
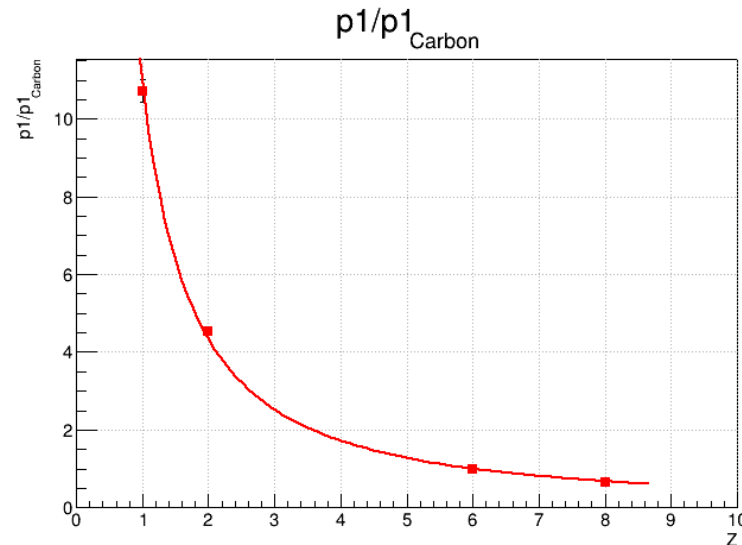
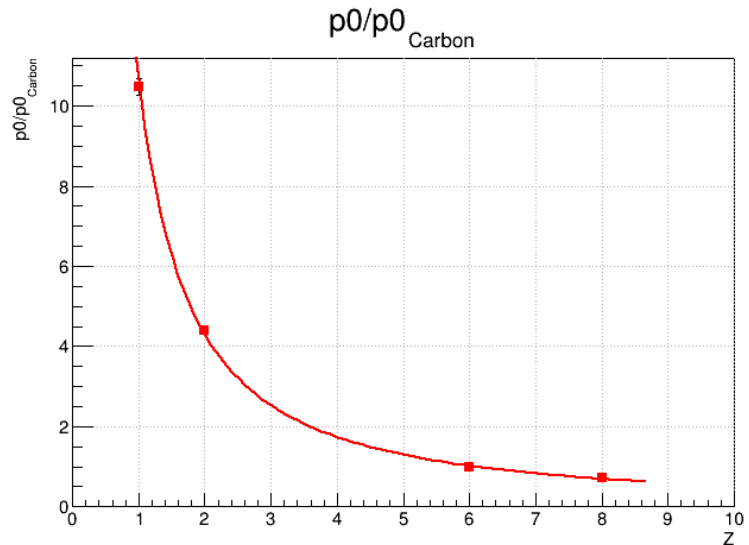
After equalization the integrated energy resolution is $\sim 1\%$



Atomic Number Dependence

Is it possible to address the Z dependence problem?

$$ADC(E) = \frac{P_0 E^2}{1 + P_1 E + P_2 E^2}$$



Ratio of energy calibration curve parameters to carbon curve parameters follow a power law



By measuring Z we can convert the ADC response to Energy for all the possible fragments

$$\frac{P_x}{P_{x_{carbon}}} = a_0 Z^{a_1}$$

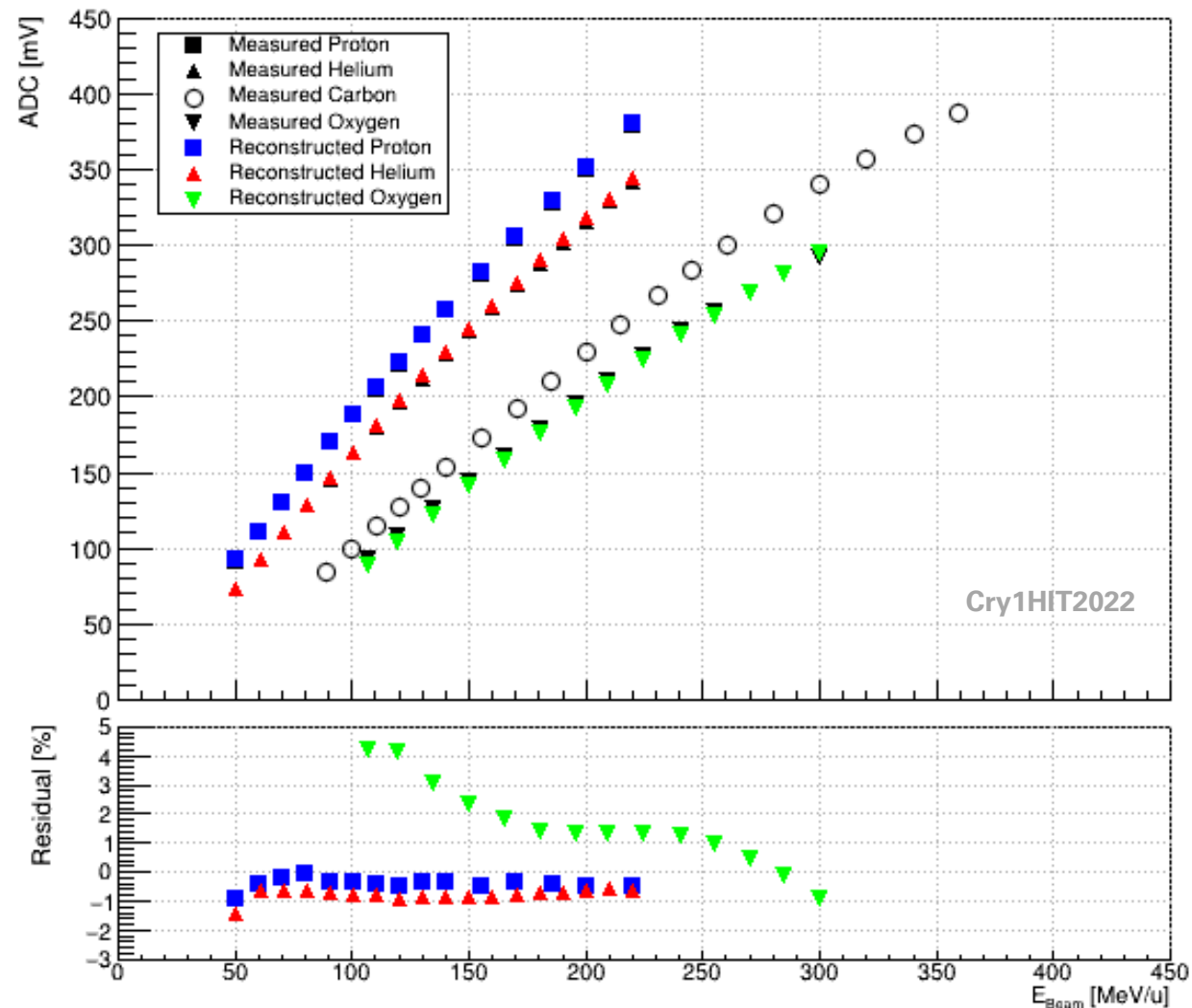


Energy Calibration

With the power law function from previous slide, it is possible to calculate **proton**, **helium** and **oxygen** point starting from carbon MBF parameter

Superimposition and residual to verify the performance of this method

It is now possible to measure the kinetic energy by knowing the charge of the ion (TW)



Applicability

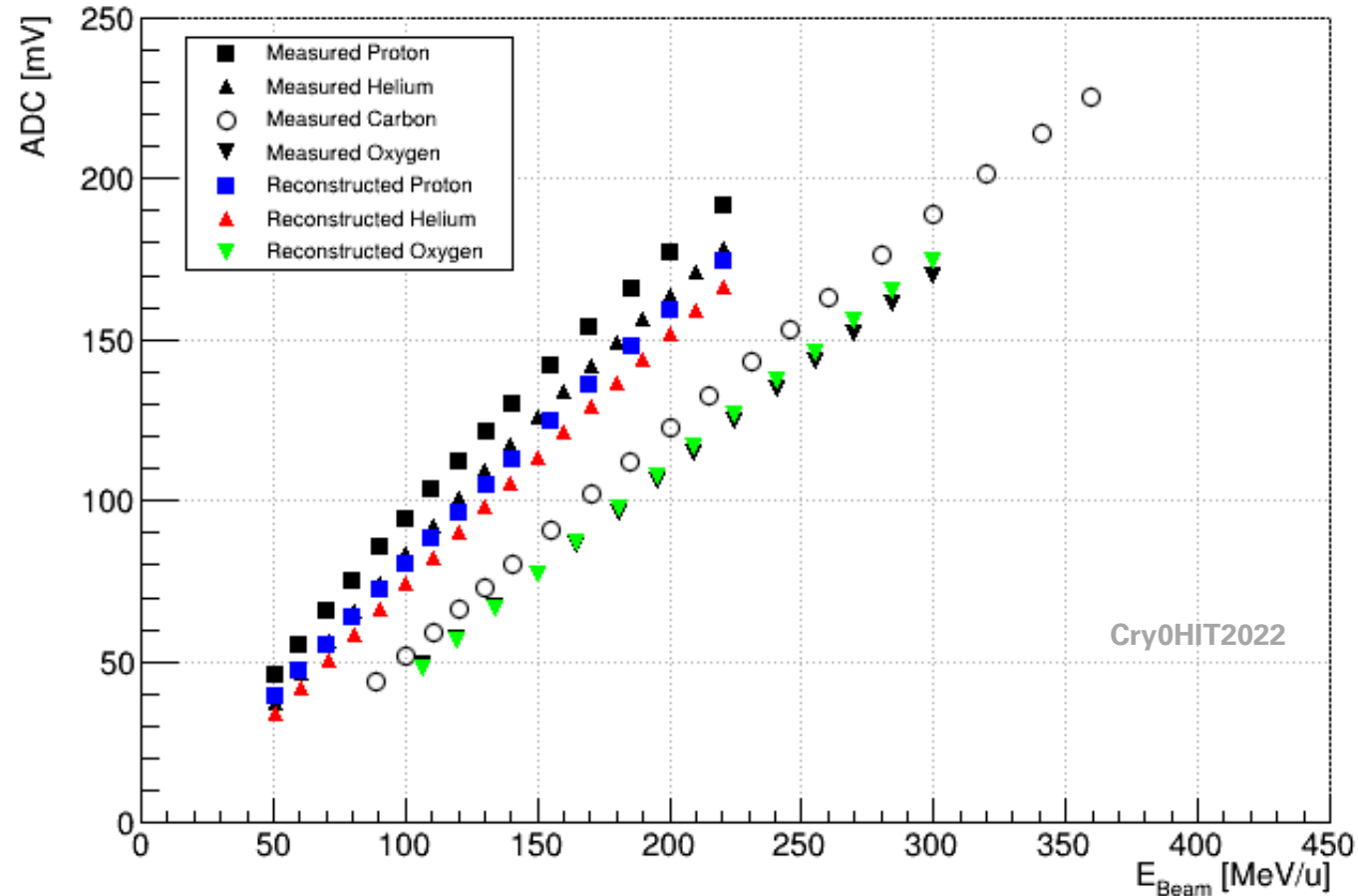
Are the power law parameters the same for every crystal?



Reconstruct proton, helium and oxygen point of Cry0HIT2022 with power law parameter calculated from Cry1HIT2022



Each crystal has its own power law parameter



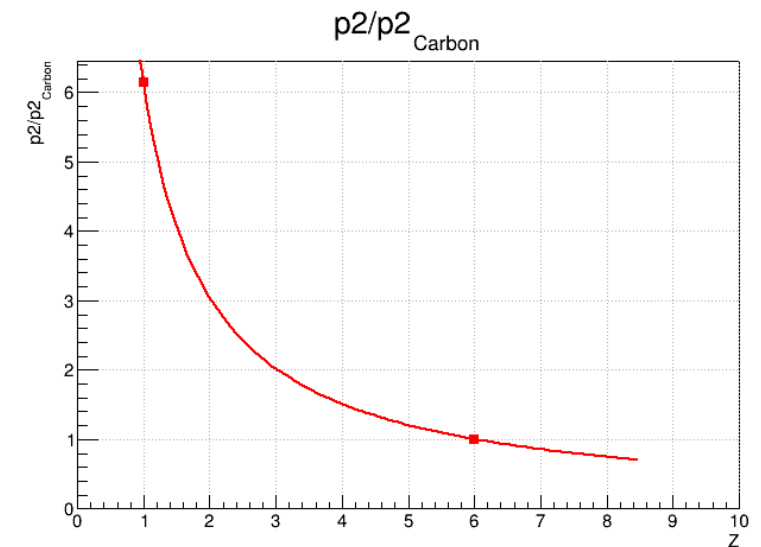
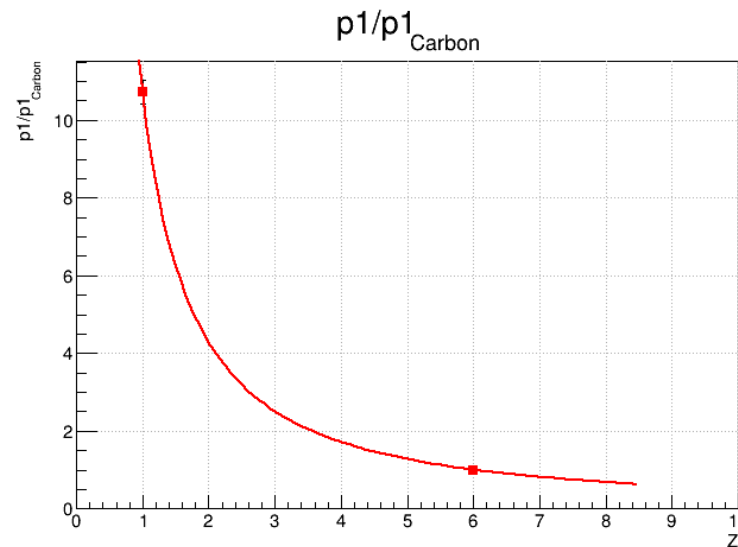
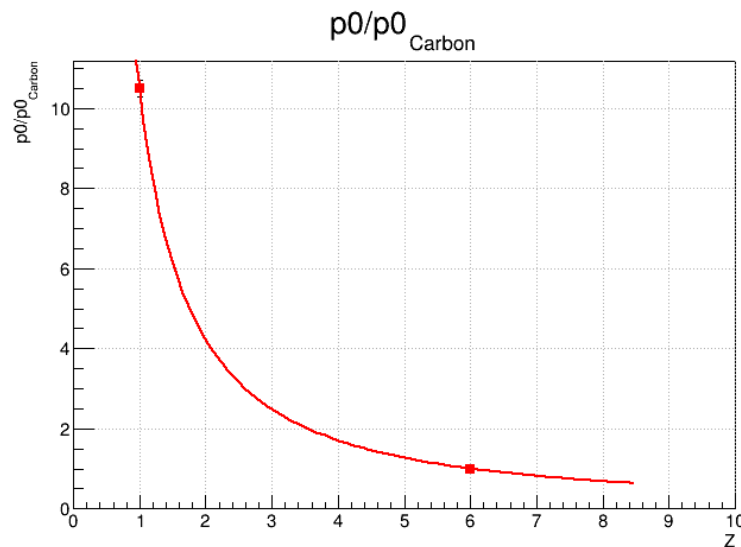
Energy Calibration – CNAO feasibility

- ✓ Taking data at HIT is a challenging task: not enough beam time available to us to measure the response of all the crystals
- ✓ Performing calibration procedure at CNAO is possible but CNAO provides only **Proton** and **Carbon** ions



What equalization strategy at CNAO?

$$\frac{P_x}{P_{x_{carbon}}} = a_0 Z^{a_1}$$

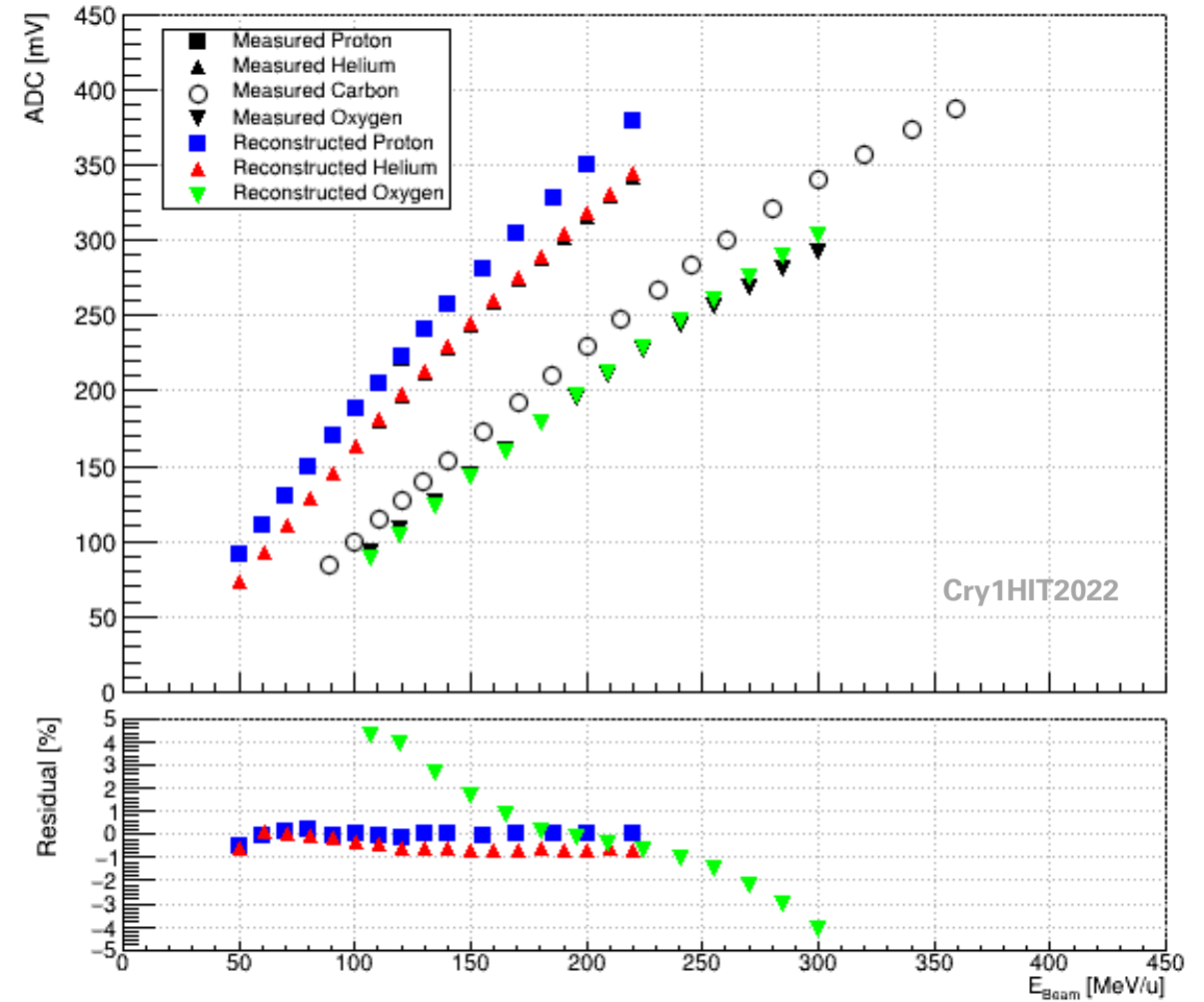


Energy Calibration – CNAO feasibility

Same validation procedure used with 4-ions-
power law



Seems possible to calibrate all crystals at
CNAO with the provided ions



Summary and Results

- ✓ Calibration run at HIT with proton, helium, carbon and oxygen ions
- ✓ BGO calorimeter is affected by Birks effect (Modified Birks function achieve $<1\%$ residual distribution)
- ✓ Modified-Birks-Parameter-function based Energy Calibration method has been identified, tested and validated on a single crystal
- ✓ Each crystal has different power-law parameters
- ✓ CNAO allows calibration for all crystals with only two ions (proton and carbon) using screensaver run
- ✓ Screensaver run meets experimental requirement: **resolution is well below $< 2\%$**
- ✓ Next step: full calorimeter calibration

