



Calorimeter calibration @ CNAO 2024: preliminary results

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Calibration runs

First calibration day: right crystals (back view)
>> CALO was shut down, then turned on for the second calibration day on left crystals
>> no further shutdown for the rest of data taking (and physics runs)



Х	1	2	9	10	11	18	19	20	27	28	29	36	37	38	45	46	X
3	4	5	12	13	14	21	22	23	30	31	32	39	<mark>40</mark>	41	48	49	50
6	7	8	15	<mark>16</mark>	17	24	25	26	33	<mark>34</mark>	<mark>35</mark>	42	43	44	51	52	<mark>53</mark>
<mark>54</mark>	55	<mark>56</mark>	63	64	65	72	73	74	81	82	83	90	91	92	99	<mark>100</mark>	<mark>101</mark>
57	58	59	66	67	68	75	76	77	<mark>84</mark>	85	86	93	94	95	<mark>102</mark>	<mark>103</mark>	<mark>104</mark>
60	61	<mark>62</mark>	<mark>69</mark>	70	71	78	79	80	87	88	89	96	97	98	<mark>105</mark>	<mark>106</mark>	<mark>107</mark>
<mark>108</mark>	<mark>109</mark>	<mark>110</mark>	<mark>117</mark>	<mark>118</mark>	<mark>119</mark>	<mark>126</mark>	<mark>127</mark>	<mark>128</mark>	<mark>135</mark>	<mark>136</mark>	<mark>137</mark>	<mark>144</mark>	<mark>145</mark>	<mark>146</mark>	<mark>153</mark>	<mark>154</mark>	<mark>155</mark>
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<mark>219</mark>	<mark>220</mark>	<mark>221</mark>	<mark>228</mark>	<mark>229</mark>	<mark>230</mark>	<mark>237</mark>	<mark>238</mark>	<mark>239</mark>	<mark>246</mark>	<mark>247</mark>	<mark>248</mark>	<mark>255</mark>	<mark>256</mark>	<mark>257</mark>	<mark>264</mark>	<mark>265</mark>	<mark>266</mark>
<mark>222</mark>	<mark>223</mark>	<mark>224</mark>	<mark>231</mark>	<mark>232</mark>	<mark>233</mark>	<mark>240</mark>	<mark>241</mark>	<mark>242</mark>	<mark>249</mark>	<mark>250</mark>	<mark>251</mark>	<mark>258</mark>	<mark>259</mark>	<mark>260</mark>	<mark>267</mark>	<mark>268</mark>	<mark>269</mark>
<mark>270</mark>	271	<mark>272</mark>	<mark>279</mark>	<mark>280</mark>	<mark>281</mark>	<mark>288</mark>	<mark>289</mark>	<mark>290</mark>	<mark>297</mark>	<mark>298</mark>	<mark>299</mark>	<mark>306</mark>	<mark>307</mark>	<mark>308</mark>	<mark>315</mark>	<mark>316</mark>	<mark>317</mark>
<mark>273</mark>	274	<mark>275</mark>	<mark>282</mark>	<mark>283</mark>	<mark>284</mark>	<mark>291</mark>	<mark>292</mark>	<mark>293</mark>	<mark>300</mark>	<mark>301</mark>	<mark>302</mark>	<mark>309</mark>	<mark>310</mark>	<mark>311</mark>	<mark>318</mark>	319	320
×	277	<mark>278</mark>	<mark>285</mark>	286	287	<mark>294</mark>	<mark>295</mark>	<mark>296</mark>	<mark>303</mark>	<mark>304</mark>	<mark>305</mark>	<mark>312</mark>	<mark>313</mark>	<mark>314</mark>	<mark>321</mark>	322	×







Calibration runs

First calibration day: right crystals (back view) >> CALO was shut down, then turned on for the second calibration day on left crystals

>> no further shutdown for the rest of data taking (and physics runs)

Operational notes:

42 43 44

85 86 93 94 95 102 103 104

88 89 96 97 98 105 106 107

35 136 137 144 145 146 153 154 155

138 139 140 147 148 149 156 157 158

141 142 143 <mark>150 151 152</mark> 159 160 161 189 190 191 198 199 200 207 208 209

92 193 194 201 202 203 210 211 212

195 196 197 204 205 206 213 214 215

43 244 245 252 253 254 261 262 263

246 247 248 255 256 257 264 265 266

249 250 251 258 259 260 267 268 26**9**

97 298 299 306 307 308 315 316 31

00 301 302 309 310 311 318 319 320

303 304 305 <mark>312 313 314</mark> 321 32

- LV @ 6V, HV @ 33V;
- WD threshold set to -5 mV;
- CALO scanned in quarters as usual.

During the second calibration round: - M32-9 unplugged due to tripping; - M9-6 threshold was set to -10 mV (noisy channel). Scanning with:

- p @ 100, 125, 150, 170, 230 MeV - C @ 115, 200, 260, 300 MeV/u (also 400 MeV/u was attempted, but then discarded due to beamsharing issues)





1
)
9
5
4

Reminder(s):

- CNAO scanning magnets cover a 20x20 cm^2 surface \rightarrow one CALO quarter at the time;
- CALO front face is concave, not flat;
- higher energies \rightarrow higher range, but increase for p is faster than for C.



At higher energies, due to CALO-beam geometry:

>> particles cross multiple crystals with higher probability

>> charge-sharing phenomena between neighbor crystals become not negligible

>> peak events collection for p @ 230 MeV and C @ 330/400 MeV/u is harder.





Sometimes, higher statistics is required (or more scanning precision)...





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...sometimes it's not enough!

Even though the range of p in BGO is \approx 7 cm, very few particles cross this full length due to the scanning geometry. >> p @ 230 MeV/u were discarded from subsequent analysis, due to the lack of "full energy peak" events.





With p @ 230 MeV, statistics should be 4 times higher than with C @ 330 MeV/u in order to actually separate peak events from beam sharing ones.

Temperature analysis

Map_run6874





Temperature analysis





T corrections have small impact on calibration runs (in this case the difference between the average T on single crystal and full CALO was $\approx 10^{\circ}$ C).

Single crystal variations between calibration run 6874 and physics run 7029 (3 days later) are $\approx 0.8^{\circ}C \rightarrow expected variation in ADC$ response < 1%.

Peak analysis



Amplitude

First step: cutting out all events with amplitude > 0 in other crystals >> events with multiplicity > 1 are discarded

Peak analysis





First step: cutting out all events with amplitude > 0 in other crystals >> events with multiplicity > 1 are discarded



Sometimes, ~ all peak events feature small charge sharing with other crystals \rightarrow underestimation of ADC response vs energy >> what to do in such cases? How big is this underestimation?

Amplitude

First correction attempt





Cross energy

In cry 1

In cry 126, if we only preserve multiplicity = 1 events, there's no peak event for C @ 260 MeV/u.

However, if a neighbor crystal i is already calibrated, we can try reconstructing the energy in 126 as:

 $E_{126} = E_{beam}$ (nominal energy) – E_i \rightarrow (ADC, E)₁₂₆ scatter plot to be fitted with a MBF.

First correction attempt





Unfortunately, events with multiplicity = 2 below 200 MeV/u are denser, thus "privileged" by MBF fit >> the peak position is not well reconstructed (280 MeV/u instead of 260 MeV/u)

>> new multiplicity cut to be attempted.

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$E_{126} = E_{beam}$ (nominal energy) – E_i \rightarrow (ADC, E)₁₂₆ scatter plot to be fitted with a MBF.



Second correction attempt





In this plot, the amplitude in Ch126 is plotted against the sum of amplitudes in other crystals \rightarrow let's select the events where sum in other crystals / sum in single crystal is < 5%.

Second correction attempt





In this plot, the amplitude in Ch126 is plotted against the sum of amplitudes in other crystals \rightarrow let's select the events where sum in other crystals / sum in single crystal is < 5%.

Now, the peak appears, and its position is underestimated of < 5% \rightarrow this correction method is meaningful.

Peak analysis

Peaks are fitted with a crystal ball function \rightarrow central gaussian region and decreasing "tail" under a certain threshold, to model small energy release events.

In some cases, small energy releases are dominated by beam-sharing effects \rightarrow fit with Gauss function.

Amplitude

 $f(x) = a \exp\left[-\frac{1}{2} \exp\left[-\frac{$



Minuit2 fit status, covariance matrix, reduced chisquare, etc. are always checked and the limits on range and σ are changed subsequently. Some data points were taken in both rounds, in order to check stability \rightarrow in almost all analyzed crystals, shifting in peak position was < 2%.

40

30

10

20



histoclone 6824 83

990

30.08

11.55

Entries

Std Dev

60

Amp (mV)

Mean

50

Peak analysis





Proton_82

ADC response vs energy is then modeled with a modified Birks function (MBF):



At least 3 points per crystal are needed.

Calibration status (front view)





Let's focus on the central 4x4 ones >> more useful for fragmentation analysis.

Energy correction

For green/light green single crystals (among the 4x4 central ones), energy could be reconstructed via the inverse equation:

 $E(ADC) = \frac{-p_1 ADC - \sqrt{(p_1 ADC)^2 - 4ADC(p_2 ADC - p_0)}}{2(p_2 ADC - p_0)}$

>> equalized energy spectra corresponding to a single nominal beam energy are summed up

>> a histogram is obtained, then again fitted with a crystal ball function

>> integral resolution is calculated.



Energy



Integral resolution (protons)





Resolution is around 3%; in addition to this, the ratio between peak and low energy events progressively builds up @ higher beam energies, due to beam sharing phenomena.

Integral resolution (carbon)





In carbon, energy resolution is well below 2%. In this case, low energy events due to beam sharing are limited within 0 and 50 MeV/u >> geometrical reasons to be further investigated.

Residual distributions for MBF

60

50

40

30

20

10

-Ŭ.05



In both cases, residual distributions are limited within 1% >> MBF reproduces nominal energy with the expected precision.

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



MBF extrapolation using **HIT2022** parameters







We tried using the power-law coefficient $(a_0, a_1)_{pi}$ extrapolated from HIT2022, to reconstruct proton MBF parameters (p_0, p_1, p_2) for each crystal @ CNAO2024 (see Alessandro's presentation).



MBF extrapolation using HIT2022 parameters



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Blue: RecEnergy in ch124 via direct MBF fit Red: RecEnergy in ch124 via p_i extrapolation Green: RecEnergy in ch127 via p_i extrapolation.



MBF extrapolation using HIT2022 parameters





Blue: RecEnergy in ch124 via direct MBF fit Red: RecEnergy in ch124 via p_i extrapolation Green: RecEnergy in ch127 via p_i extrapolation.

Energy hEnergy ch124 220 Entries 4204 p @ 125 MeV 99.69 Mean 200 Std Dev 28.16 180 160 Energy is wrongly reconstructed, with no 140 systematic over/underestimation 120 100 >> power-law coefficient not only depend 80 on work conditions (i.e. HV power supply, 60 gain), but are crystal-specific! 40 20 0 50 100 250 150 200 E (MeV/u)

Conclusion



• By applying a new multiplicity cut, an integral resolution < 2% was achieved for the central 4x4 modules, in spite of beam sharing issues.



- For p, integral resolution was not below 2%: this was already known from the Heidelberg 2022 campaign, during which 2 crystals were calibrated singularly (no screen saver runs), thus excluding beam sharing issues.
- MBF parameters must be corrected with respect to Z for each crystal specifically. How to assess whether this is done effectively?

Upcoming tasks

In order to assess the CALO energy reconstruction performances we need to:

- obtain power-law coefficients (a_0, a_1) for all crystals suitable for p and C calibration;
- within CNAO2024 work conditions, try reconstructing the kinetic energy of a different ion species
 - \rightarrow ⁴He produced during C \rightarrow C fragmentation runs is the best choice;
- ⁴He mass is our "bench-test", to be obtained by the known formula:

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

In parallel:

- calibration must be completed for modules outside the central 4x4 subset;
- integral resolution of the detector must be measured with all functioning crystals;
- does energy reconstruction still show good resolution performances when multiplicity > 1? To be checked.





Reminders (T calibration)





Arduino reads temperatures in ADC channels, to be converted into T values via the Steinhart – Hart expression:

 $\ln(1/T) = a_0 + a_1 \ln(T_{ADC}) + a_2 \ln(T_{ADC})^3$





By using the so-obtained three coefficients a_0 , a_1 , a_2 , a custom script plots a T histogram for each crystal. T is known with a 0.1 °C sensitivity.

Reminders (T calibration)

Why is temperature correction necessary?



Analysis carried out on a single crystal shown that ADC signals due to beams with same energy decrease linearly with temperature.



only slightly changes the distribution of the parameters p_0 , p_1 and p_2 used for the energy calibration.





Reminders (simulated mass reconstruction)





Reminder(s):

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At higher energies, due to CALO-beam geometry:



>> charge-sharing phenomena between neighbor crystals become not negligible

>> peak events collection for p @ 230 MeV and C @ 330/400 MeV/u is harder.



In addition to this, readout amplitude depends on how particle-crystal interaction is close to the SiPM.

