



AGATA Week 2024

Improvements in the Post – PSA analysis

- Neutron Damage correction
- Time dependent energy calibration



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Neutron Damage correction - Introduction



- Neutron Damage (ND) from fast neutrons (E > 1 MeV) in HPGe detectors causes charge trapping and therefore energy resolution degradation
- The loss in signal amplitude is dependent on the **path length**, the **electric field** and the **trap density**
- In highly segmented HPGe detectors, we can use Pulse Shape Analysis (PSA) to determine the **position of** interaction of the gamma – rays inside the crystal
- The current model is optimized for fast neutrons only and employs a 1st order Taylor expansion in its calculations

$$\frac{E_{meas}(x)}{E_{corr}(x)} = 1 + \frac{t_e(x)}{\lambda_e} + \frac{t_h(x)}{\lambda_h}$$

 $\lambda_{e,h}$ = inverse electron / hole trap density $t_{e,h}(x)$ = sensitivity to electron / hole trapping

B. Bruyneel et al, EPJ A 49 (2013)

Charge trapping

CATHODE

ANODE



Critical for lineshape analysis in DSAM lifetime measurements





Algorithm code: https://gitlab.in2p3.fr/ip2igamma/agapro/-/blob/preprod/zPrograms/SortPsaHits.cpp



LOGARITMIC GRID – SEARCH

Fixed 50 x 100 grid



• Source data (usually ⁶⁰Co)

- Only segment multiplicity 1 events
- Optimization of SG, CC or SG+CC
- Estimation of a FOM
- Computation in each point of the grid, for each segment of each crystal

Ease of use, speed and robustness are key

The objective is to obtain the $\lambda_{e,h}$ parameters that yield the

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Various methods of estimating a FOM implemented

- 1. By fitting (Gaussian + tails): **amplitude** of Gaussian part
 - Slower, computation heavy
 - Less robust, works well with slightly damaged detectors

2. FOM = **h** / σ

- Introduced recently (2023) as the "-std" option
- Very fast
- · More robust, works with badly damaged detectors



OBS: for SG+CC, the algorithm just sums the FOMs of the two spectra



Issues:

- **h** is the height of a single bin
 - Large statistical variance
 - Binning dependent
- σ is largely dependent on background
 - Not very significant
 - Range dependent



We want to keep ease of use, robustness, speed and backwards compatibility.

Two main improvements:

- For **SG+CC**, we employ a **normalization** of the whole grid of FOMs (such that the highest FOM has value 1) before calculating the average
- Implementation of an Adaptive Grid Search
 - "-size" option: variable grid size, default is kept at 50 x 100
 - "-algo" option: user must choose which spectra to optimize (SG, CC or SG+CC) and number of iterations. Default is kept at 1 iteration only
 - "zoom" option: user can specify the magnification factor M between iterations. Default is 0.25



Neutron Damage correction – Improvements



- ⁶⁰Co source data taken in November 2023
- Optimization of CC only with 3 iterations and 30 x 50 grid
- Better performance especially for badly damaged detectors and for front and back segments





*crystals are defined "problematic" if at least one segment shows a visible tail in the CC spectrum after ND correction





- **Core energy obscillations** were observed in a long run taken with a high energy source
- Previously observed during the LNL Agata demonstrator campaign
 - Pseudo periodic with 6h period
 - More pronounced in detectors placed in the lower part of Agata, pointing upwards
 - Gain obscillations up to 2.91‰





Implementation of time – dependent core gain correction before RecalCC



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Time dependent energy calibration - algorithm used



Cross – correlation Correction Method (CCM)

- Divide the matrix in slices, for each a projection
- Calculate dot product between a reference specturm and the shifted projection
- Perform a scan for each slice

Main characteristics

- Extremely fast and robust, no fitting involved
- No need for a peak, any feature can be used





Matus Balogh NIM paper: <u>https://doi.org/10.1016/j.nima.2021.165368</u> GitHub code: <u>https://github.com/matLogh/CCM</u>





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Thank you for your attention



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