### **GRETINA / GRETA Signal Decomposition**

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## Outline

- The Algorithms
  - Signal basis generation
    - siggen
    - xtalk
    - Grid
  - Signal decomposition algorithm
    - One hit segment
    - Two hit segments
    - More...
    - Adaptive Grid Search
- Strengths and weaknesses, possible improvements
  - Signal basis
  - Decomposition with more resources
- Summary



# **Signal Decomposition**

- Digital signal processing to determine the *number, positions,* and *energies* of gamma interactions in the crystal
- Uses a "signal basis"; a set of simulated signals
- Position resolution is crucial for energy resolution, efficiency, and peak-to-total ratio in tracking
- But getting the *number* of interactions correct may be harder, and is at least as important
- Speed is critical as it determines overall gamma throughput of array





## **Expected Distribution of Hits**

GEANT simulations; 1 MeV gamma into GRETA

Most hit crystals have one or two hit segments

Most hit segments have one or two interactions





## What Can Affect the Signals?

- Field and Weighting Potential:
  - Overall impurity concentration

(Two values (~20%) from maker, one at each end; depletion voltage)

- Longitudinal impurity gradient (Linear? Nonlinear?)
- Radial impurity gradient?
- Hole diameter; hole depth; etching cycles; lithium thickness
- Neutron damage (p-type)
- Charge carrier mobilities as a function of electric field
- Crystal axis orientation (~ 5 degrees from maker)
- Crystal temperature (Some info from RTD)
- Cross-talk (differential and integral)
- Neutron damage (trapping)
- Impulse response of 37 preamps
- Charge cloud size
- Digitizer nonlinearity



## Signal Generation: AGATA (ADL)



B. Bruyneel, P. Reiter, G. Pascovici, Characterization of large volume HPGe detectors. Part I: Electron and hole mobility parameterization B. Bruyneel, P. Reiter, G. Pascovici, Characterization of large volume HPGe detectors. Part II: Experimental results



## Signal Generation: fieldgen and siggen







#### Signal Generation: xtalk

#### fieldgen siggen plus... Geant MC Positions Crystals axis Geometry, orientation Material **Calculate Traces** Poisson Electrical Path of charge Space **Mobilities** Solver Field Charge carriers Induced charges Transfer Weighting in electrodes Function Potential Convolute Signal Crosstalk superpulses xtalk xtalk parameters

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## **Fitting to Extract Cross-Talk Parameters**

- 36 "superpulses": averaged signals from many single-segment events (red)
- Monte-Carlo simulations used to generate corresponding calculated signals (green)
- ~ 996 parameters fitted (integral and differential cross-talk, delays, rise times) (blue)
- Calculated response can then be applied to decomposition "basis signals"



## Signal Generation: xtalk superpulses

- Collect "flood-field" data from <sup>60</sup>Co source at known location (target location
- Select events with single hit segment and 1332 keV energy
- Time-align using digital CFD on both CC and hit segment signals
- Make an averaged "super-pulse" for each of the 36 segments
- Run GEANT simulation of <sup>60</sup>Co data
- Again select single-segment, full-energy events
- Generate signals, including multi-site (but single-segment) events
- Add noise, time-align as for measured data



## xtalk Fitted Parameters

Fit function and included parameters were determined empirically (by trial and error)

- (2 x 36) delays (channel-to-channel)
- (37) preamp rise-times plus (2) extra times
  - segment-segment and segment-CC xtalk
- (5 x 36 + 36) integral cross-talk coefficients (selected pairs)
  - asymmetric
- (17 x 36 / 2 + 36) differential cross-talk coefficients
  - or (35 \* 36 / 2 + 36); all pairs
  - symmetric except for CC
- Total of 669 or 993 parameters
- Parameters can be fixed at specified values if desired





## **Regular Basis Grid**

 The GRETINA Signal Decomposition originally made use of a Cartesian grid

Different colors show active regions for different segments



- An irregular quasi-cylindrical grid has several important advantages:
  - The possibility to optimize the spacing of points in the grid based on separation in "Chi-squared space"
  - Reducing the number of grid points for improved speed
  - Constructing the grid around the real segment volumes allows much better and faster constraints to be programmed into the least-squares search algorithms

## **Optimized Quasi-Cylindrical Grid**





## **Decomposition Basis (Signal Library)**

- Signal decomposition **algorithm** appears to work very well
  - Validated using simulated signals
- Most issues with the decomposition results appear to come from the fidelity of the signal basis
- Poor fidelity results in
  - Too many fitted interactions
  - Incorrect positions and energies
- We already include effects of
  - Integral cross-talk
  - Differential cross-talk
  - Preamplifier rise-time
- *Differential cross-talk* signals look like image charges, so they strongly affect position determination



## **Data Flow: Preparation for Decomposition**

• Raw waveform timing derived from online CC LED





## **Decomposition Fits**

- Shows two typical multi-segment events measured in prototype triplet cluster (red) (concatenated signals from 36 segments, 500ns time range)
- Linear combination of basis signals, as fitted by decomposition algorithm (blue)
- Includes differential cross talk from capacitive coupling between channels



## **Decomposition Fits**



## **Decomposition Fits**





## **GRETINA Decomposition Algorithm**

Current algorithm is a *hybrid* 

- Adaptive Grid Search with Linear Least-Squares (for energies)
  - Finds best-fit pair of interactions in one segment
  - Limited to positions on grid
  - $t_0$  of signals fixed
- Non-linear Least-Squares
  - Interpolate off-grid
  - $t_0$  of signals fitted
- Have also tried Singular Value Decomposition as replacement for AGS
  - Had slightly poorer performance for same CPU time
  - CPU time required goes as
    - Adaptive Grid Search:  $\sim O(n + \delta n^*t)$
    - Nonlinear Least-Squares : ~ O(n\*t + δn<sup>2</sup>) for n interactions and t time steps



## Why is it Hard?

- Very large parameter space to search
  - Average segment ~ 6000 mm<sup>3</sup>, so for ~ 1 mm grid search,
    - two interactions in one segment:  $\sim 2 \times 10^6$  possible positions
    - two interactions in each of two segments:  $\sim 4 \times 10^{12}$  positions
    - two interactions in each of three segments:  $\sim 8 \times 10^{18}$  positions
    - PLUS additional dimensions; energy sharing, time-zero, ...
- Under-constrained fits, especially with > 1 interaction/segment
  - For one segment, the signals provide only ~ 6 x 40 = 240 nontrivial numbers
- Strongly-varying, nonlinear sensitivity
  - $\delta \chi^2 / \delta(\theta z)$  is much larger near segment boundaries



#### **Problems with the Basis have a Big Effect**

Same data set, decomposed with old and new bases Crystal Q4P4



## **Overall Strategy: One hit segment**

- 1. Start by finding approximate  $t_0$ 
  - Fit a single interaction and time offset using nonlinear least-squares
  - Five parameters:  $\mathbf{x}_1$ ,  $\mathbf{e}_1$ ,  $\mathbf{t}_0$
  - Potential problem for multi-site events!
- 2. Then find best *two-interaction* solution
  - Adaptive grid search using ~  $10^5$  pairs of grid points and best-fit energies
    - Much more detail about this later
  - Then interpolate off basis grid using nonlinear least-squares
    - Two interactions, nine parameters:  $\mathbf{x}_1$ ,  $\mathbf{e}_1$ ,  $\mathbf{x}_2$ ,  $\mathbf{e}_2$ ,  $\mathbf{t}_0$
- 3. Try adding a third interaction (if total energy is > 400 keV and chisq is bad)
  - Insert extra interaction in middle of segment, with 1/3 of the energy
  - Re-do nonlinear least-squares  $\mathbf{x}_1, \mathbf{e}_1, \mathbf{x}_2, \mathbf{e}_2, \mathbf{x}_3, \mathbf{e}_3, \mathbf{t}_0$
- 4. Try coalescing two interactions into one
  - Re-do nonlinear least-squares  $\mathbf{x}_{1}, \mathbf{e}_{1}, \mathbf{t}_{0}$
- Choose best overall solution, with penalty factor for extra parameters 5 (i.e. extra interactions). End up with 1, 2, or 3 interactions.



## **Overall Strategy: Two hit segments**

- 1. List hit segments in order of decreasing energy;  $e_a > e_b$
- 2. Start by finding approximate fit, with one interaction per segment
  - Nonlinear least-squares  $\mathbf{x}_{a1}, \mathbf{e}_{a1}, \mathbf{x}_{b1}, \mathbf{e}_{b1}, \mathbf{t}_{0}$
- 3. Subtract calculated signals for  $(\mathbf{x}_{b1}, \mathbf{e}_{b1}, \mathbf{t}_0)$  from the measured signals.
  - Use adaptive grid search to find best *two-interaction* solution for the remainder (segment a).
- 4. Now have three interactions:  $\mathbf{x}_{a1}$ ,  $\mathbf{e}_{a1}$ ,  $\mathbf{x}_{a2}$ ,  $\mathbf{e}_{a2}$ ,  $\mathbf{x}_{b1}$ ,  $\mathbf{e}_{b1}$ ,  $\mathbf{t}_0$ 
  - Re-fit full signal using nonlinear least-squares, 13 parameters
- 5. Use the same trick: Subtract calculated signals for  $(\mathbf{x}_{a1}, \mathbf{e}_{a1}, \mathbf{x}_{a2}, \mathbf{e}_{a2}, \mathbf{t}_0)$  from the measured signals.
  - Use adaptive grid search to find best *two-interaction* solution for the remainder (segment b).
  - Re-fit full signal using nonlinear least-squares, with 4 interactions

 $\mathbf{x}_{a1}, \, \mathbf{e}_{a1}, \, \mathbf{x}_{a2}, \, \mathbf{e}_{a2}, \, \mathbf{x}_{b1}, \, \mathbf{e}_{b1}, \, \mathbf{x}_{b2}, \, \mathbf{e}_{b2}, \, \, \mathbf{t}_0$ 



## **Overall Strategy: Two hit segments**

- 6. For both segments, try coalescing the pairs of interactions into one
  - Re-do nonlinear least-squares each time
- 7. Choose best overall solution, with penalty factor for extra parameters. End up with 2, 3, or 4 interactions.



## **Overall Strategy: Three or more hit segments**

- 1. List hit segments in order of decreasing energy;  $e_a > e_b > e_c$
- 2. Start by finding approximate fit, with one interaction per segment
  - Three interactions, plus t<sub>0</sub>
- 3. Subtract calculated signals for segments b and c from the measured signals.
  - Use adaptive grid search to find best *two-interaction* solution for the remainder (segment a).
  - Re-fit full signal using nonlinear least-squares with 4 interactions
- 4. Repeat step 3 twice more, to get pairs of interactions in segments b and c.
- 5. For all segments, try coalescing the pairs of interactions into one, re-doing nonlinear least-squares each time
- 6. Choose best overall solution, with penalty factor for extra parameters. End up with 3 – 6 interactions.



## **Overall Strategy: Post-processing**

Independent of number of hit segments

- Optional recombination of interactions if their decomposed positions are too close together
- Energy-dependent separation limit
- Try to include cross-segment-boundary pairs
- Never been studied or optimized?



## **Adaptive Grid Search Least-Squares**

Adaptive grid search fitting:

- Critical that the signals start at  $t_0 = 0$  for reliable results!
- Use for only one segment at a time
- Start on a coarse grid, every second point in each direction (2x2x2)
  - All the in-segment basis dot products are pre-calculated on this coarse grid
- Loop over all pairs of positions inside the segment,
  - Energies  $e_i$  and  $e_j$  are constrained, such that 0.1 ( $e_i + e_j$ ) <  $e_i$  < 0.9 ( $e_i + e_j$ )
- Once the best pair of positions (lowest χ2) is found, then all neighbor pairs are examined on the finer (1x1x1) grid. This is 26x26 = 676 pairs. If any of them are better, the procedure is repeated.
  - Here the signal dot-products cannot be pre-calculated
- Finally, nonlinear least-squares (SQP) can be used to interpolate off the grid. This improves the fit ~ 50% of the time.



#### **Adaptive Grid Search Least-Squares**

#### Linear Least-Squares

For two interactions of energies  $e_i, e_j$  at locations i and j, the calculated signal is  $C_{kt} = (e_i s_{ikt} + e_2 s_{jkt})$  where k is the segment and t the time step.  $s_{ikt}$  is the basis signal calculated at point i.

If the observed signal is  $O_{kt}$ 

$$\chi^{2} = \sum_{kt} \frac{(O_{kt} - C_{kt})^{2}}{\sigma_{kt}^{2}} = \frac{\sum_{kt} (O_{kt} - e_{i}s_{ikt} - e_{j}s_{jkt})^{2}}{\sigma^{2}}$$
(1)

where  $\sigma_{kt} = \sigma$  is the uncertainty (noise) in  $O_{kt}$ , assumed independent of k, t.

We want a minimum in  $\chi^2$ , *i.e.* 

$$\frac{\partial \chi^2}{\partial e_i} = \frac{\partial \chi^2}{\partial e_j} = 0 \tag{2}$$

$$\frac{\partial \chi^2}{\partial e_i} = \frac{2\sum_{kt}(O_{kt}s_{ikt} - e_i s_{ikt}^2 - e_j s_{ikt} s_{jkt})}{\sigma^2} = 0 \tag{3}$$

#### **Adaptive Grid Search Least-Squares**

Thus we get two equations in two unknowns:

$$\sum_{kt} O_{kt} s_{ikt} - e_i \sum_{kt} s_{ikt}^2 - e_j \sum_{kt} s_{ikt} s_{jkt} = 0$$
(4)  
$$\sum_{kt} O_{kt} s_{jkt} - e_j \sum_{kt} s_{jkt}^2 - e_i \sum_{kt} s_{ikt} s_{jkt} = 0$$
(5)

kt

kt

$$\sum_{kt} s_{ikt}^2$$

kt

and

$$\sum_{kt} s_{ikt} s_{jkt}$$

once for all events, and

$$\sum_{kt} O_{kt} s_{jkt}$$

once per event.



## **Adaptive Grid Search: Some Numbers**

(Cartesian grid for illustration purposes):

- ~35000 grid points in 1/6 crystal (one column, 1x1x1 mm)
- ~ 2x2x2 mm (slices 1-3) or ~ 3x3x3 mm (slices 4-6) coarse grid gives N ≤ 600 course grid points per segment.
- For two interactions in one segment, have N(N-1)/2 ≤ 1.8 x 10<sup>5</sup> pairs of points for grid search. This takes < 3 ms/cpu to run through.</li>
- Two segments:
  - (N(N-1)/2)<sup>2</sup> ~ 3.2 x 10<sup>10</sup> combinations for two interactions in each of 2 segments; unfeasible!
  - Limit N to only 4<sup>3</sup> = 64 points; then (N(N-1)/2)<sup>2</sup> ~ 4 x 10<sup>6</sup>
     This may be possible? But is it worthwhile?
- Three segments:
  - But (N(N-1)/2)<sup>3</sup> ~ 8 x 10<sup>9</sup> combinations for two interactions; impossible even for N = 64.



## **Adaptive Grid Search: Some Numbers**

- What about 1-interaction x 1-interaction in two segments, on the coarse grid?
  - Requires a very large number of pre-calculated dot-products
  - We now calculate ~ 2e5 sums for each of 36 segments
  - For all pairs of segments, would need ~ 4e5 for 630 pairs
    - 35 times the storage is required
    - But still only  $\sim$  1.5 GB, roughly the same as the basis signals
    - Entirely feasible today
- But would this be useful?
- Remember that the grid search relies on knowing t<sub>0</sub> accurately...



## **Strengths and Weaknesses**

- Able to identify up to 2 interactions per segment (three for a single segment)
- Finds correct solution in simulation tests
- Fast
- Modest memory requirements
- Optimized, irregular grid makes a very significant difference
  - Took some serious coding and a lot of time on the part of K. Lagergren
- Poor determination of number of interactions
- Strong covariance between reported interaction positions and t<sub>0</sub>
  - Fitted t<sub>0</sub> distribution is wider than normal CFT distribution



## **Covariance:** t<sub>0</sub> and position

• Some results from varying *fixed* t<sub>0</sub> in very early tests:



## **Covariance:** t<sub>0</sub> and position

## **Clipping (time windows) - 1 interaction**



Mario Cromaz AGATA-GRETINA Tracking Arrays Collaboration Meeting, Dec. 5-7, 2016



## Time windowing overview

- DAQ uses LED trigger [Timestamp] to make an imperfect selection of trace. Attempt to capture "fit region" + excess trace before and after. 186 trace samples collected.
- Preprocessor (CPU farm) runs a CFD filter and selects a subset for fitting [T<sub>offset</sub>]. 50 samples fit.
- Decomposition fits basis traces to experiment providing the final timing [T<sub>fit</sub>].
- $T_{crystal} = Timestamp + T_{offset} + T_{fit}$



What I call  $t_0$  - fitted, not fixed



#### Testing setup

- LaBr scintillator + <sup>166m</sup>Ho @ Gretina center
- Collect coincidences between LaBr and a single Gretina crystal
- Drive local trace output by delayed LaBr trigger for that same Gretina crystal
- So, the collected traces for that crystal will be read out in a window set by the LaBr timing
- If time windowing is perfect, expect a ~fixed T<sub>offset</sub>
- After Decomp, expect  $T_{offset} + T_{fit}$  to give a constant timing w.r.t LaBr
- Deviations can show problems which will arise in all gamma ray data. Here we simply have a handle on good/bad time(s).
- Examine how timing can be mistaken and the effect on Decomp, i.e. interaction points



## Timing observed





#### Investigating timing issue Gated on 184keV



CAK RIDGE

### Interaction Points (XY) [crystal coordinates]

Gated 184keV





Interaction Points (ZR) [crystal coordinates] Gated 184keV





Interaction point multiplicity Gated 184keV









- fieldgen code improvements from MJD
  - Adaptive grid
  - Undepleted voxels
  - More parameters in impurity profile
    - Quadratic, radial, passivated surface
  - Capacitance calculation
- siggen improvements
  - Cloud size (not directly applicable  $\ensuremath{\mathfrak{S}}$  )
- xtalk improvements
  - Dead layers in simulation



- Extra timing information to constrain t<sub>0</sub>
  - External fast detectors or RF signal
  - Ge-Ge coincidences (i.e. global t<sub>0</sub> fitting)
    - Requires event building prior to decomposition; hard!
- Better initial estimate of t<sub>0</sub>
  - Especially important for multi-site events in a segment
- Better interaction-number penalty algorithm
  - e.g. position-dependent penalty
  - Tuning of penalty factors crystal-by-crystal or segment-by-segment
- Improved irregular grid
  - Somewhat reduced chisq-dependence
  - Does grid depend on impurity profile?
- Try to include charge-sharing at segment boundaries in basis



- Other improvements in basis fidelity
  - Preamplifier impulse response function
  - Include charge cloud size and charge-sharing in signal generation
    - Especially important at small radius, near segment boundaries
    - Also dead layers at outside of segments
    - But charge size is energy-dependent?
  - Better field determination
    - From segment capacitance measurements as a function of bias
  - <sup>241</sup>Am surface-scan "superpulse" fitting for field, WP, electron drift, and preamp parameters
- Compressed basis signals to save RAM and memory bandwidth?
- 50 MS/s basis to save CPU time? Or...



#### What more could be done with more RAM & CPU?

- AGS with multiple guesses for t<sub>0</sub>
- Longer basis and signal time windows
- Better time-interpolation algorithms, and/or higher-frequency basis signals
- Two-segment AGS?
  - Requires additional 1.5 GB RAM for pre-calculated products
  - Is it worthwhile? (t<sub>0</sub>)
- More...



- Preamplifier impulse response function
- Ben Shanks for point-contact detector



## Some conclusions from Mario, Dec 2016

- First/primary/isolated/... hit position resolution as measured through source tests and Doppler reconstruction is very good [1-2 mm], secondary hits appears degraded, clustering on boundaries
- Simulations with basis signals indicate current algorithm performs well for 1-, 2-interactions in the presence of Gaussian noise

Areas for focus:

- Impact of timing errors, especially for low-E signals
- Better time determination, longer basis windows)
- Better understanding of noise characteristics (correlated noise)
- Examine impact, find sources



## Summary

- The algorithm is very complex
- Desired result is computationally under-determined
- But the method works reliably when the basis is known perfectly
- Fast, relatively modest memory requirements
- Basis fidelity is crucial
- Strong covariance between position and t<sub>0</sub>
- Tendency to overestimate the number of interactions
  - Requires penalty factors



#### Acknowledgements

Karin Lagergren (ORNL / UTK)

- Signal calculation code in C
- Optimized pseudo-cylindrical grid
- I-Yang Lee
  - Original signal calculation code
- C. Campbell, H. Crawford, M. Cromaz, M. Descovich, P. Fallon, A. Machiavelli, ...
  - Basis calculations, cross-talk fits, in-beam data analysis, simulations, electric field calculations, and much more

Tech-X Corp, especially Isidoros Doxas

• SVD development



#### Backup



## **Calculated Signals: Sensitivity to Position**



National Laboratory

## **Singular Value Decomposition**

Very roughly:

- The full signal -vs.- grid position matrix can be decomposed into the product of three matrices, one of which contains the correlations (eigenvalues)
- By neglecting the small eigenvalues, the length of the signal vectors (and hence computation with them) can be greatly reduced
- The more eigenvalues kept, the higher the quality of the fit





## **Singular Value Decomposition**

Very roughly:

- The full signal -vs.- grid position matrix can be decomposed into the product of three matrices, one of which contains the correlations (eigenvalues)
- By neglecting the small eigenvalues, the length of the signal vectors (and hence computation with them) can be greatly reduced
- The more eigenvalues kept, the higher the quality of the fit
- Measured signals can be compressed the same way as, and then compared to, the calculated library signals
- Different similarity measures can be used to emphasize different aspects

Dot Product 
$$x.y = \sum_{i=1}^{N} x_i y_i$$
  
Cosine  $\cos(\theta_{xy}) = \frac{x.y}{|x||y|}$   
uclidean Distance  $euclid(x, y) = \sqrt{\sum_{i=1}^{N} (x_i - y_i)^2}$ 



F