

# AGATA Analysis Workshop 2025

## PostPSA Calibration. Hands on

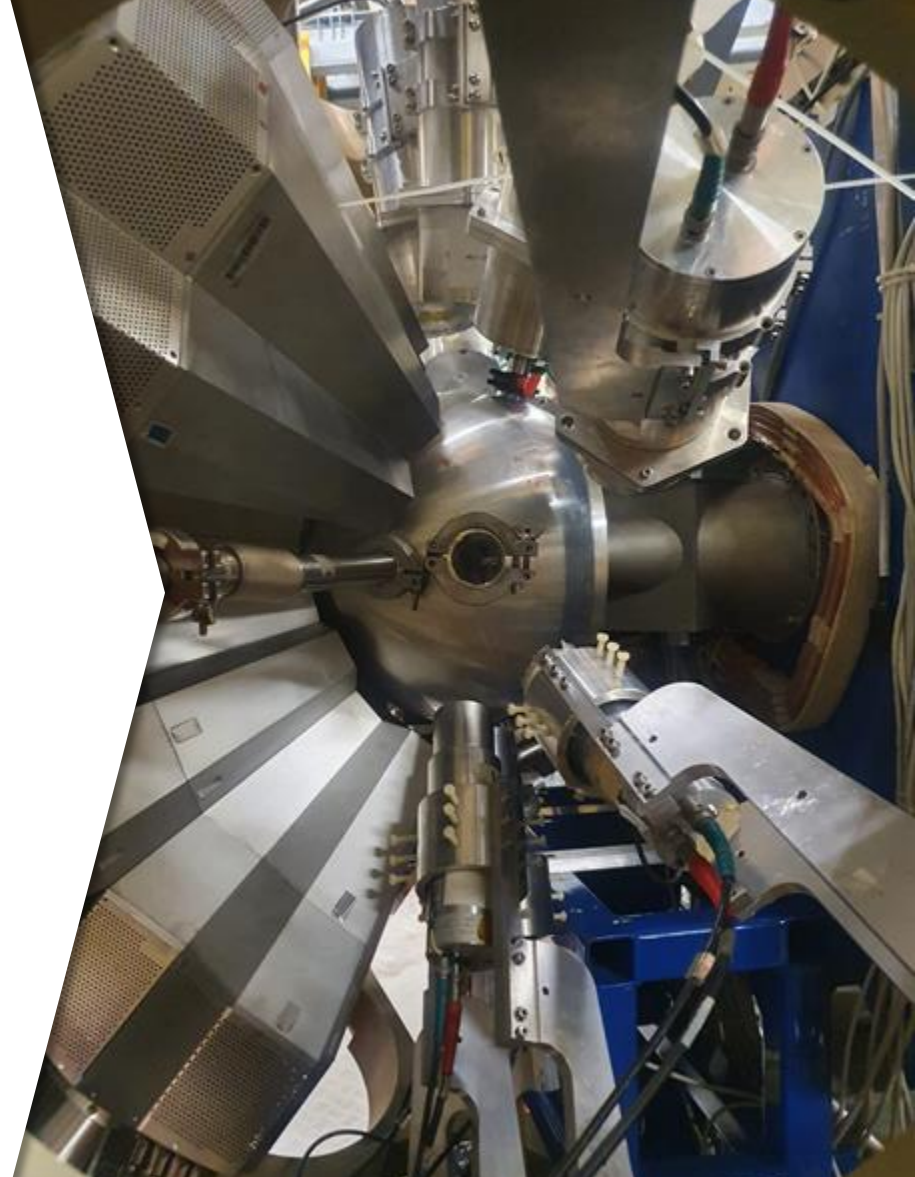
**R.M. Pérez-Vidal**

14/01/2025, Lyon

# Hands On

## AGATA configuration

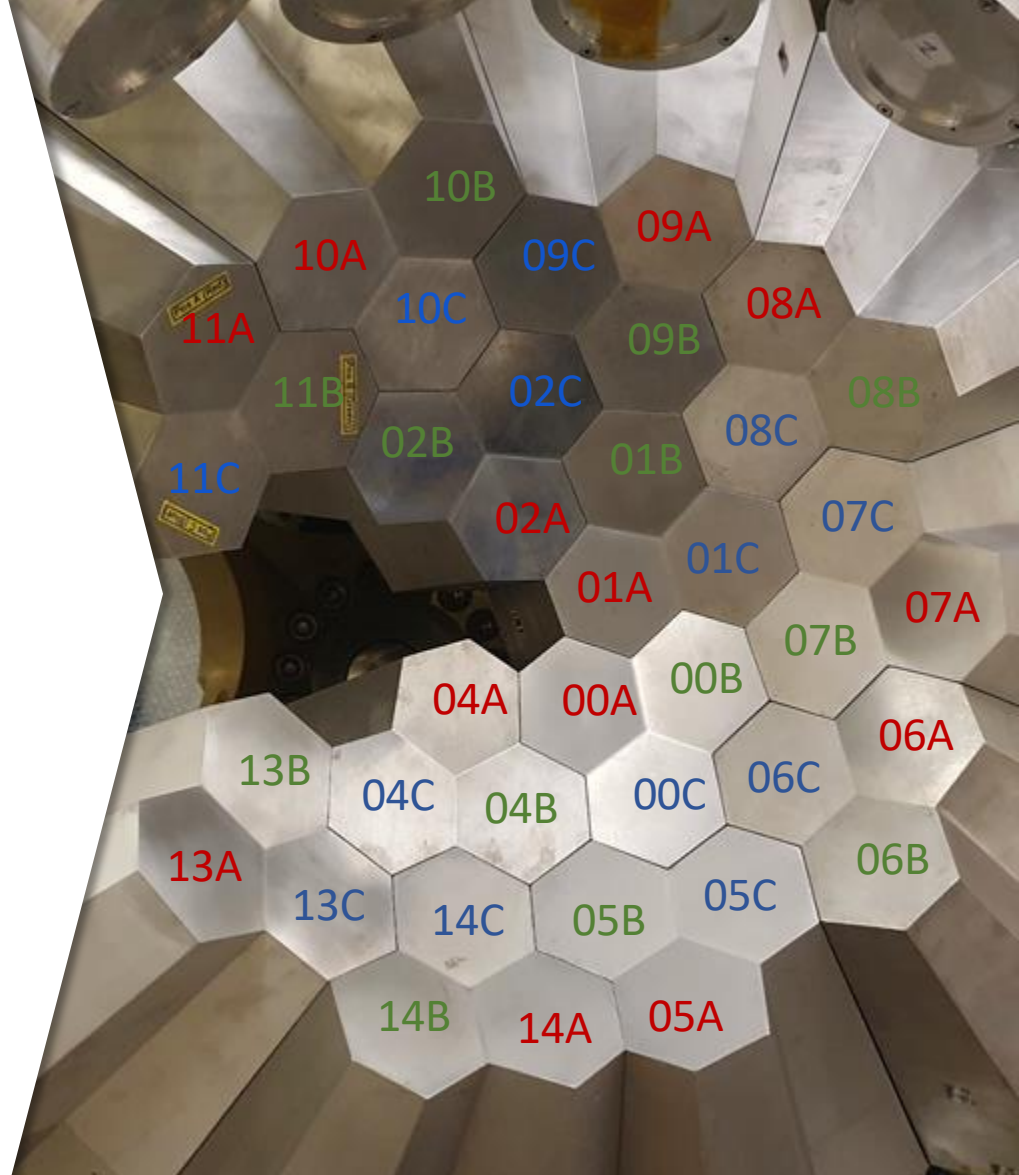
- AGATA data set taken in July 2024
- Counting rate per crystal : 1-2 kHz
- MWD-risetime: 2.5  $\mu$ s
- Chamber closed, Without absorbers
- Position: Nominal
- Data sets with  $^{60}\text{Co}$ ,  $^{152}\text{Eu}$  and target ( $^{66}\text{Ga}$ )



# Hands On

## Data sets in example directory

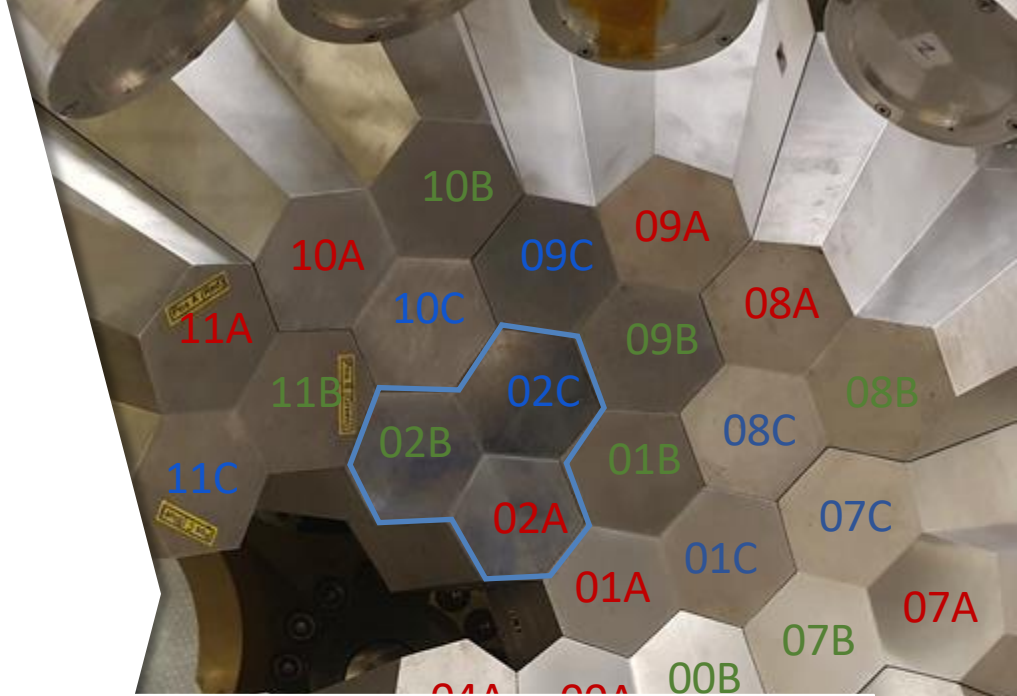
- **run\_1003\_60Co\_ATC02 (33Gb)**
  - Neutron damage correction
  - 1 ATC: 02A 02B 02C with traces
- **run\_1002\_152Eu\_ATC02 (3.7Gb)**
  - Energy recalibration
  - 1 ATC: 02A 02B 02C
- **run\_1010\_60Co (Gb)**
  - Global time alignment
  - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 14
- **run\_1010\_66Ga\_3ATC (8.4Gb)**
  - Tracking
  - 3 ATC: 02 10 11



# Hands On

## Data sets in example directory

- **run\_1003\_60Co\_ATC02 (33Gb)**
  - Neutron damage correction
  - 1 ATC: 02A 02B 02C **with traces**
- **run\_1002\_152Eu\_ATC02 (3.7Gb)**
  - Energy recalibration
  - 1 ATC: 02A 02B 02C
- **run\_1010\_60Co (3.2Gb)**
  - Global time alignment
  - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 14
- **run\_1010\_66Ga\_3ATC (8.4Gb)**
  - Tracking
  - 3 ATC: 02 10 11



To be taken into account:

- 02C: seg. missing **recovered** as a lost seg. 9

POS	ATC	Crys	Capsule
2	17	A	16
		B	17
		C	13

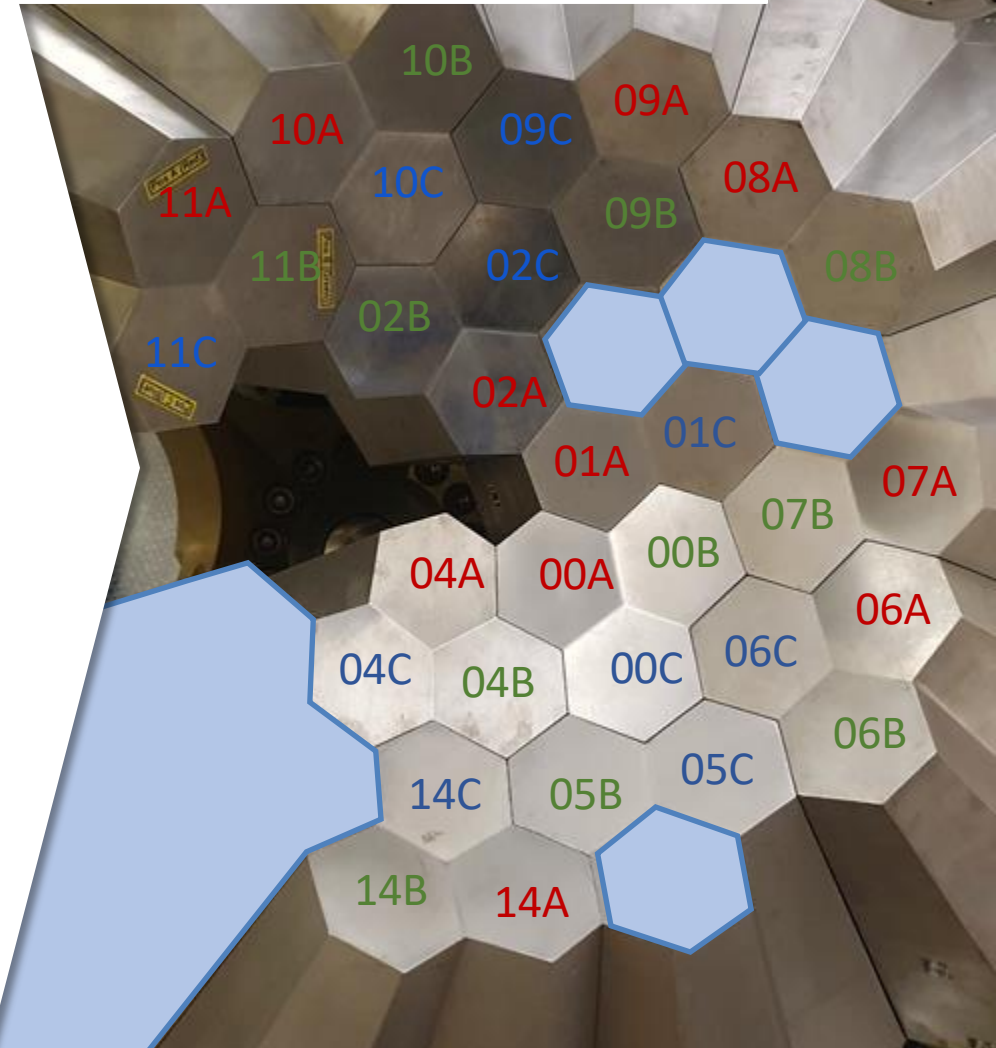
# Hands On

## Data sets in example directory

- **run\_1003\_60Co\_ATC02 (33Gb)**
  - Neutron damage correction
  - 1 ATC: 02A 02B 02C with traces
- **run\_1002\_152Eu\_ATC02 (3.7Gb)**
  - Energy recalibration
  - 1 ATC: 02A 02B 02C
- **run\_1010\_60Co (3.2Gb)**
  - Global time alignment
  - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 14
- **run\_1010\_66Ga\_3ATC (8.4Gb)**
  - Tracking
  - 3 ATC: 02 10 11

### To be taken into account:

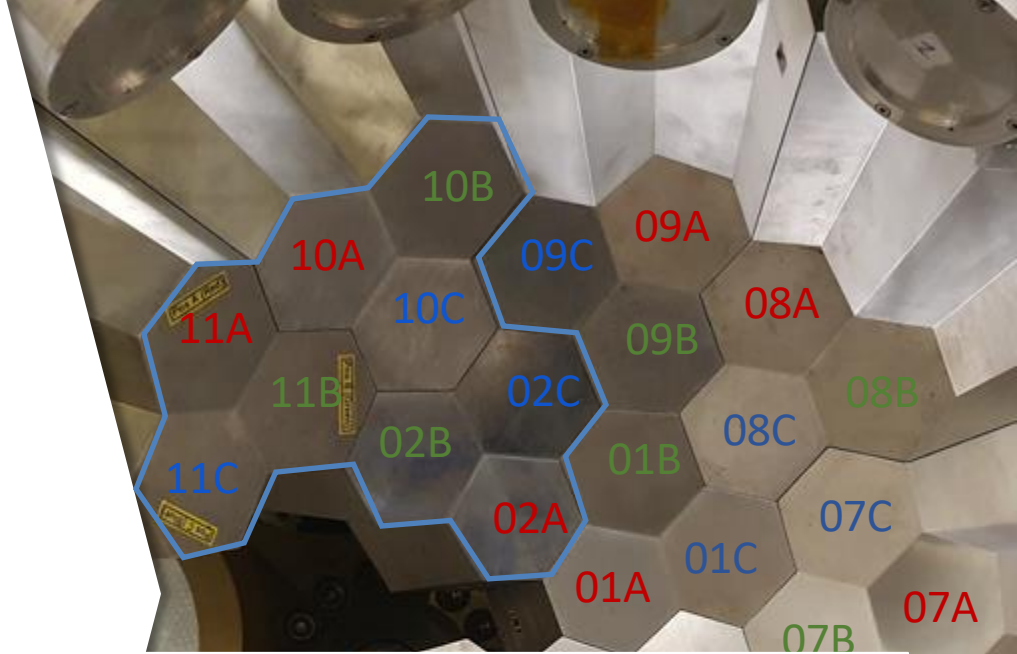
- 13ABC out of the structure
- 01B, 05A, 07C, 08C excluded from the acquisition



# Hands On

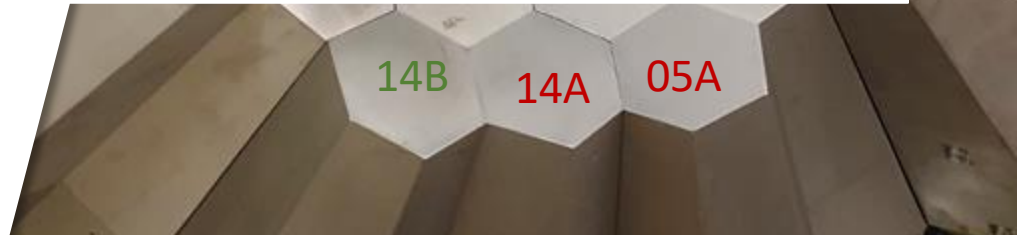
## Data sets in example directory

- **run\_1003\_60Co\_ATC02 (33Gb)**
  - Neutron damage correction
  - 1 ATC: 02A 02B 02C with traces
- **run\_1002\_152Eu\_ATC02 (3.7Gb)**
  - Energy recalibration
  - 1 ATC: 02A 02B 02C
- **run\_1010\_60Co (3.2Gb)**
  - Global time alignment
  - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 14
- **run\_1010\_66Ga\_3ATC (8.4Gb)**
  - Tracking
  - 3 ATC: 02 10 11



### To be taken into account:

- 02C: **seg. missing** **recovered** as a lost seg. 9
- 10B: **very neutron damaged**
- 11A: **seg. F1 missing** **recovered** as a lost seg. 30  
all segments with an extra peak  
(electronic noise filtered in the Prep level)
- 11C: column E and F swapped.  
**Corrected** in CrystalProducerATCA.conf (mapping).



# Hands On

## Files in example directory

- **In run\_\*:**
  - Conf/ configurations, read during initialization
  - Data/ part of original data and full PSA hits
  - Out/ created by gen\_conf.py
- **bases\_ADL/** calculated signal basis for 02A (A016), 02B (B017) 02C (C013)
- **In TemplatePrep/**
  - Topology\_Local.conf using Data/\*/event\_mezzdata.cdat.0000
  - Topology\_Global.conf using Data/\*/psa\_0000.adf
  - Topology\_Global\_Tree.conf using Data/\*/psa\_0000.adf
  - gen\_conf.py generator of configuration
  - ADF.conf definition of adf frames used for this analysis
  - 1-Neutron.sh some analysis scripts
  - 2-RecalSegs.sh
  - 3-Recal\_CC-SumSeg.sh
  - 4-GlobalTime.sh
  - solveTT.py
  - colupdate.py

# Hands On

## Files in example directory: run\_1003\_60Co

Neutron Damage correction

### ○ Conf/ 02A

- BasicAFC.conf
- BasicAFP.conf
- CrystalProducerATCA.conf
- CrystalProducer.conf
- PostPSAFilter.conf
- PreprocessingFilter.conf
- PreprocessingFilterPSA.conf
- PreproHist.conf
- ProdHist.conf
- PSAFilter.conf
- PSAHist.conf
- RecalEnergy2.cal
- Trapping.cal
- xdir\_1325-1340.cal
- xinv\_1325-1340.cal

### ○ Data/02A

- **event\_mezzdata.cdat.0000**
- **event\_energy.bdat.0000**
- **psa\_0000.adf**
- Prod\_\_100-42-100-S\_\_Traces.samp
- Prod\_\_38-16384-UI\_\_Baseline.spec
- Prod\_\_4-38-32768-UI\_\_Ampli.spec
- Prep\_\_100-44-100-S\_\_Traces.samp
- Prep\_\_2000-2000-UI\_\_EsEs.matr
- Prep\_\_2-1000-1000-US\_\_EeEtrCC.matr
- Prep\_\_2-10-16384-UI\_\_Esum.spec
- Prep\_\_2-2000-1000-US\_\_EcTc.matr
- Prep\_\_2-40-16384-UI\_\_Ener.spec
- Prep\_\_36-36-UI\_\_IsIs.matr
- Prep\_\_6-40-1000-UI\_\_TT.spec
- Psa\_\_100-2-42-60-F\_\_Traces.samp
- Psa\_\_2-38-37-60-F\_\_AverSingles.samp
- Psa\_\_3-100-100-100-US\_\_XYZ.matr
- Psa\_\_37-37-60-F\_\_Base.aver
- Psa\_\_37-37-60-F\_\_Base.aver\_raw
- Psa\_\_40-1000-UI\_\_RedChi.spec
- Psa\_\_40-1000-UI\_\_Tzero.spec
- Psa\_\_40-100-UI\_\_Stat.spec
- Psa\_\_40-16384-UI\_\_Ener.spec
- Psa\_\_524288-F\_\_DistanceMetric.spec
- **Psa\_\_0-16-F\_\_Hits.fdat**



# Hands On

## Files in example directory: run\_1011\_152Eu

### Recalibration

#### ○ Conf/ 02A

- BasicAFC.conf
- BasicAFP.conf
- CrystalProducerATCA.conf
- CrystalProducer.conf
- PostPSAFilter.conf
- PreprocessingFilter.conf
- PreprocessingFilterPSA.conf
- PreproHist.conf
- ProdHist.conf
- PSAFilter.conf
- PSAHist.conf
- RecalEnergy2.cal
- Trapping.cal
- xdir\_1325-1340.cal
- xinv\_1325-1340.cal

#### ○ Data/02A

- **psa\_0000.adf**
- Prod\_\_38-16384-UI\_\_Baseline.spec
- Prod\_\_4-38-32768-UI\_\_Ampli.spec
- Prep\_\_2000-2000-UI\_\_EsEs.matr
- Prep\_\_2-1000-1000-US\_\_EeEtrCC.matr
- Prep\_\_2-10-16384-UI\_\_Esum.spec
- Prep\_\_2-2000-1000-US\_\_EcTc.matr
- Prep\_\_2-40-16384-UI\_\_Ener.spec
- Prep\_\_36-36-UI\_\_IsIs.matr
- Prep\_\_6-40-1000-UI\_\_TT.spec
- Psa\_\_2-38-37-60-F\_\_AverSingles.samp
- Psa\_\_3-100-100-100-US\_\_XYZ.matr
- Psa\_\_37-37-60-F\_\_Base.aver
- Psa\_\_37-37-60-F\_\_Base.aver\_raw
- Psa\_\_40-1000-UI\_\_RedChi.spec
- Psa\_\_40-1000-UI\_\_Tzero.spec
- Psa\_\_40-100-UI\_\_Stat.spec
- Psa\_\_40-16384-UI\_\_Ener.spec
- Psa\_\_524288-F\_\_DistanceMetric.spec

# Hands On

## Files in example directory: run\_1010\_60Co

Global Time Alignment

- Conf/ 00A

- BasicAFC.conf
- BasicAFP.conf
- CrystalProducerATCA.conf
- CrystalProducer.conf
- PostPSAFilter.conf
- PreprocessingFilter.conf
- PreprocessingFilterPSA.conf
- PreproHist.conf
- ProdHist.conf
- PSAFilter.conf
- PSAHist.conf
- RecalEnergy2.cal
- Trapping.cal
- xdir\_1325-1340.cal
- xinv\_1325-1340.cal

- Data/00A

- **psa\_0000.adf**

# Hands On

## Files in example directory: Topologies

### Topology\_Local.conf

```
LOOP CRY5 02A 02B 02C
```

```
Chain 4      CRY5  
Producer    CrystalProducerATCA  
Filter      PreprocessingFilterPSA  
Filter      PSFilter  
Consumer    BasicAFC  
ENDLOOP
```

# Hands On

## Files in example directory: Topologies

### Topology\_Global.conf

```
LOOP CRY5 02A 02B 02C
```

```
Chain 3      CRY5  
Producer    BasicAFP  
Filter      PostPSAFilter  
Dispatcher  EventBuilder  
ENDLOOP
```

```
Chain 2      Builder/  
Builder     EventBuilder  
#Consumer   BasicAFC  
Dispatcher  EventMerger
```

```
Chain 3      Merger/  
Builder     EventMerger  
Filter      TrackingFilterOFT  
Consumer    None
```

# Hands On

## Files in example directory: Topologies

### Topology\_Global\_Tree.conf

```
LOOP CRY5 02A 02B 02C
```

```
Chain 3      CRY5  
Producer    BasicAFP  
Filter       PostPSAFilter  
Dispatcher  EventBuilder  
ENDLOOP
```

```
Chain 2      Builder/  
Builder      EventBuilder  
#Consumer    BasicAFC  
Dispatcher  EventMerger
```

```
Chain 3      Merger/  
Builder      EventMerger  
Filter       TrackingFilterOFT  
Consumer     TreeBuilder
```

# Data Processing

## Useful programs. PostPSA Filter

The number of channels (38 x number of detectors) to be calibrated and checked at each analysis level is too large to be done one by one: **automatic tools and procedures are distributed**

- **TkT & Mat spectrum viewer:** to plot any spectrum produced all along the actors chain
- **RecalEnergy:** Analysis of spectra looking for peaks
- **SortPsaHits:** Sort of PSA hits (special format) to determine neutron damage correction parameters
- **solveTT.py:** Optimize time alignment of “equal” detectors
- **colupdate.py:** change columns of coefficients between files

# TkT & Mat

## Download the software

After installing the agata software by following this procedure (<https://agata.pages.in2p3.fr/handbook/install-P1/install> ) you will find the software to visualize the spectra dumped by the actors in the AGAPRO package (**GammaSoftware/agapro/zUseful**):

**TkT** spectrum viewer (**TkT64.exe TkT.exe**)

**Mat** matrix viewer (**Mat64.exe, Mat.exe**)

To execute them:

```
wine TkT64.exe
```

```
wine Mat64.exe
```

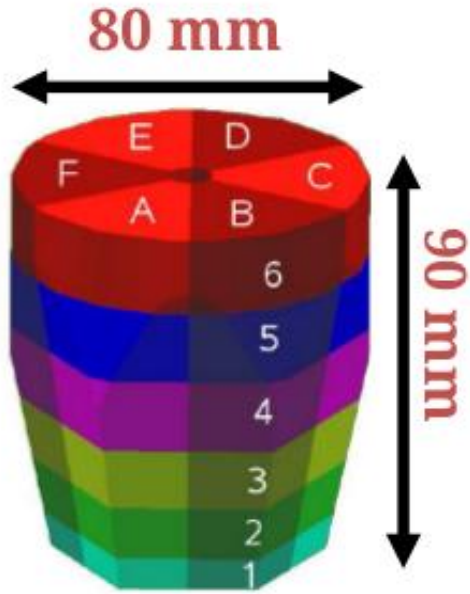
To the access to the program you can make an alias in your .bashrc

```
alias TkT='wine ~/YourPath/gammaSoftware/agapro/zUseful/TkT64.exe'
```

```
alias Mat='wine ~/YourPath/gammaSoftware/agapro/zUseful/Mat64.exe'
```

# TkT

## Channels correspondence after Replay



Back segs. 6

5

4

3

2

Front segs. 1

	A	B	C	D	E	F	CC
6	5	11	17	23	29	35	
5	4					34	
4	3					33	
3	2					32	
2	1					31	37
1	0	6	12	18	24	30	36

1 Low gain

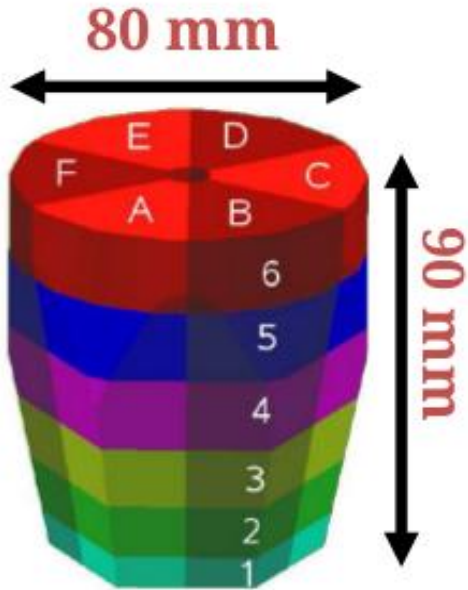
0 High gain & trigger

6x6 segmented cathode



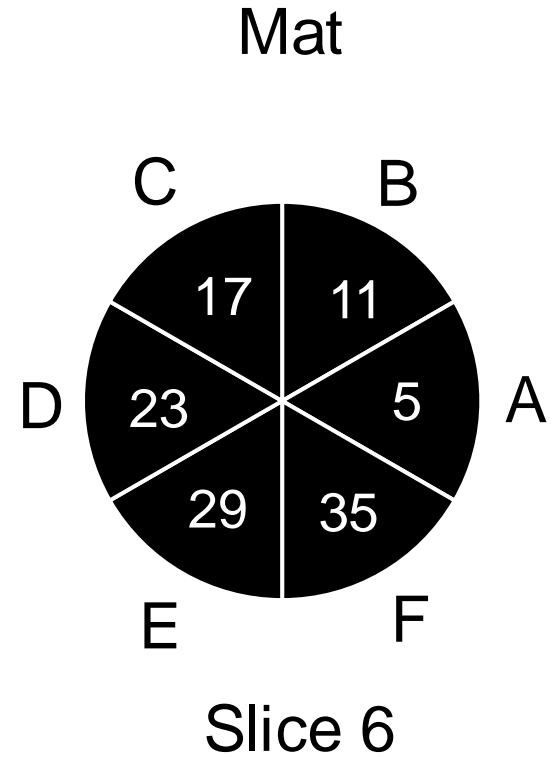
# TkT & Mat

## Channels correspondence after Replay



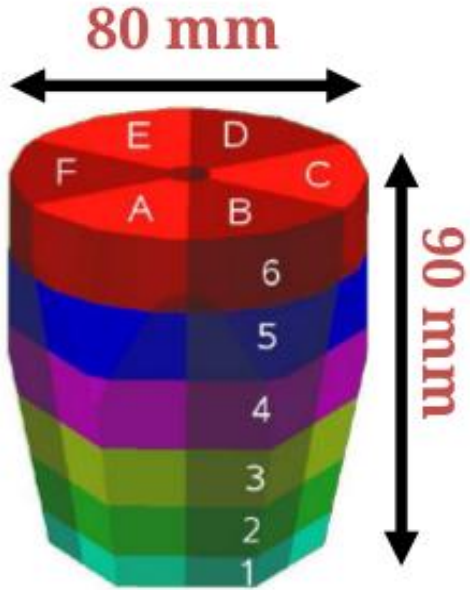
TkT

	A	B	C	D	E	F
6	5					
5	4					
4	3					
3	2					
2	1					
1	0	6	12	18	24	30



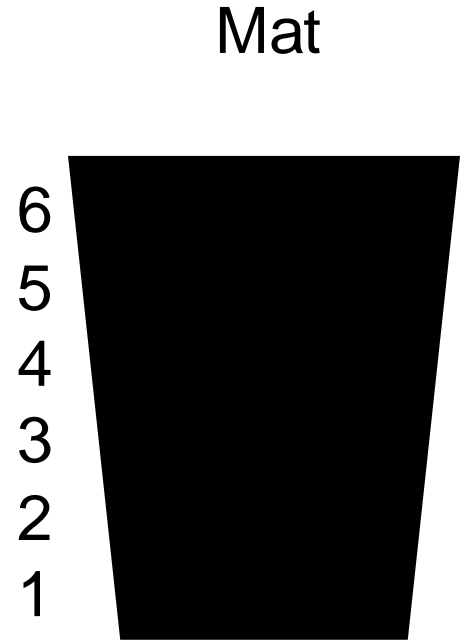
# TkT & Mat

## Channels correspondence after Replay



TkT

	A	B	C	D	E	F
6	5					
5	4					
4	3					
3	2					
2	1					
1	0	6	12	18	24	30



6x6 segmented cathode

# TkT

## Reading the format

**[Library][NbSpectra][Length][Format]** **PathToFile**

Y	0	3.26e+07	D	0	0	4	38	32768	UI	Z:\home\romapevi\AGATAWorkshop2025\AGATA\data\run_1003_60Co_ATC02\Data\02A
X	0	32767	#	0	0	0	0	38	1	Prod_4-38-32768-UI_Ampli.spec

**To choose what to display**  
Library 0 to 3  
Spectrum 0 to 37  
Display 38 spectra/ Display 1 spectrum

**FileName**

The viewers TkT and Mat can decode and interpret the format and length from the file name:

**Actor\_\_Library-NbSpectra-Length-Format\_\_Type.spec** dump file of an array defined as:  
**Format ActorSpecType[Library][NbSpectra][Length]**

E.g. **Prod\_\_4-38-32768-UI\_\_Ampli.spec** is a file dumped by the **Crystal Producer actor** containing the **amplitude spectra** of segments and cores organized in **4 libraries of 38 spectra** written in **32768 unsigned integer bins**

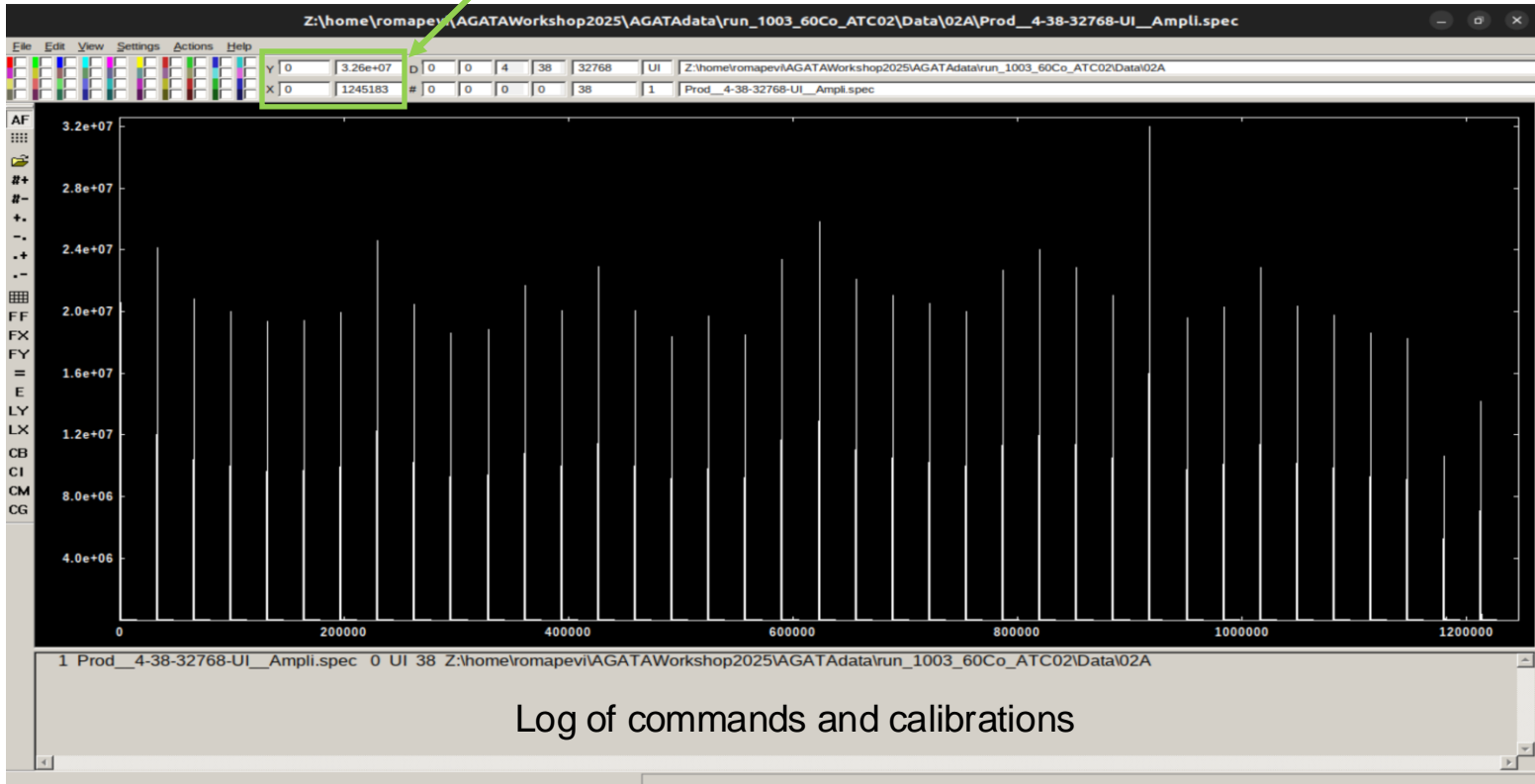
# TkT

## Some basics to get started

Boxes to store spectra

Axis ranges

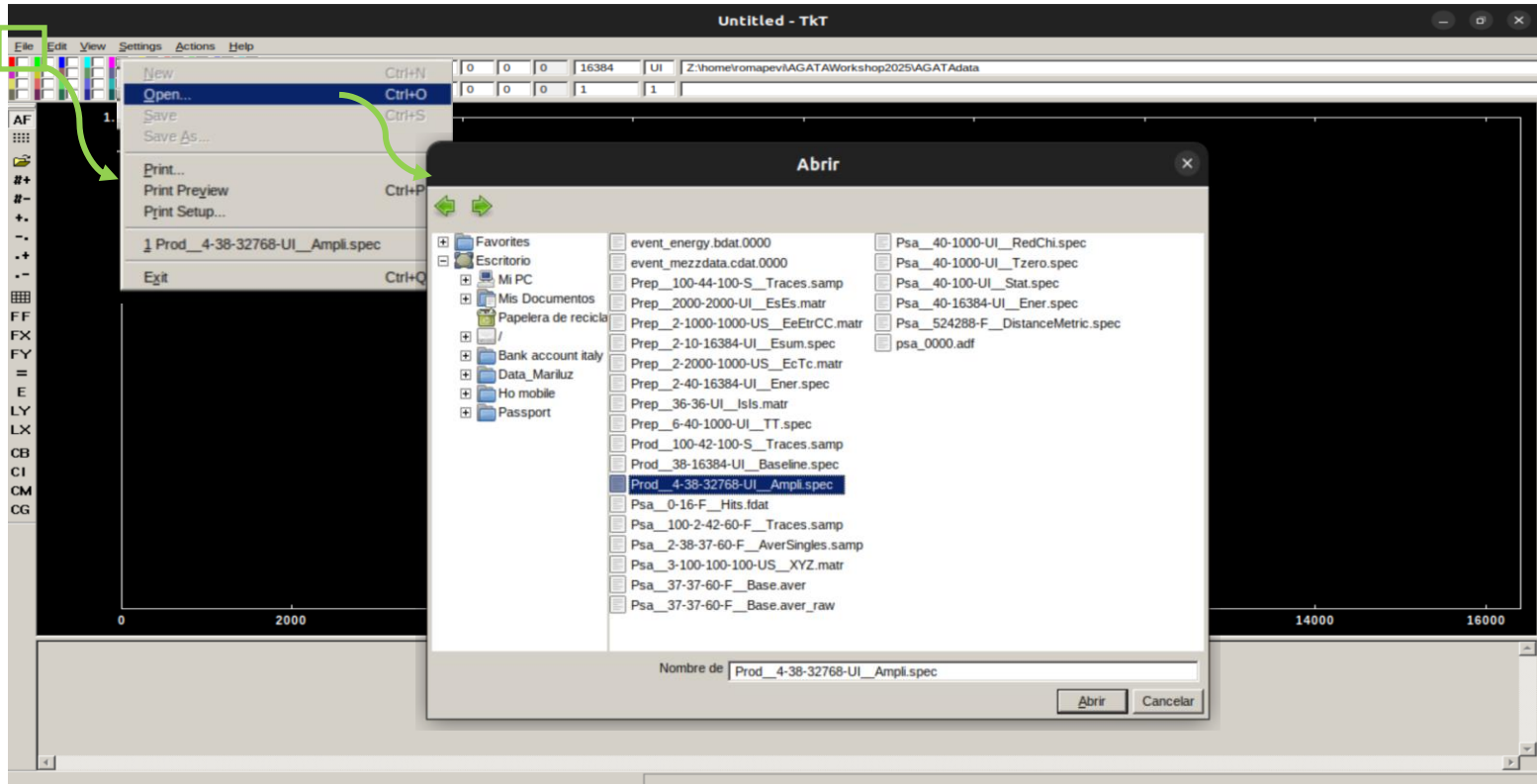
Shortcuts





## How to open a file

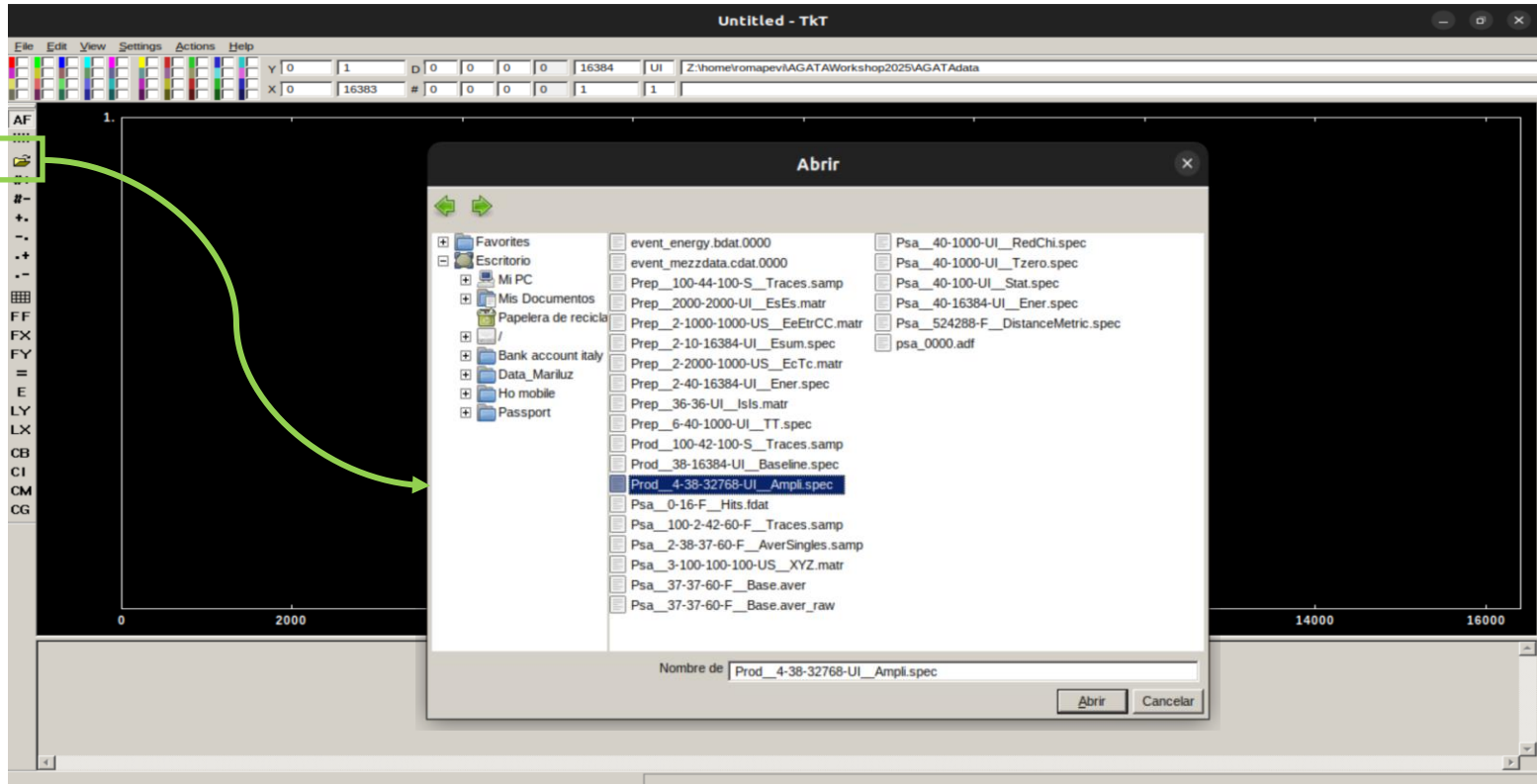
Execute Tkt → go to File → Open → select a file (e.g. Prod\_\_4-38-32768-UI\_\_Ampli.spec)





## How to open a file

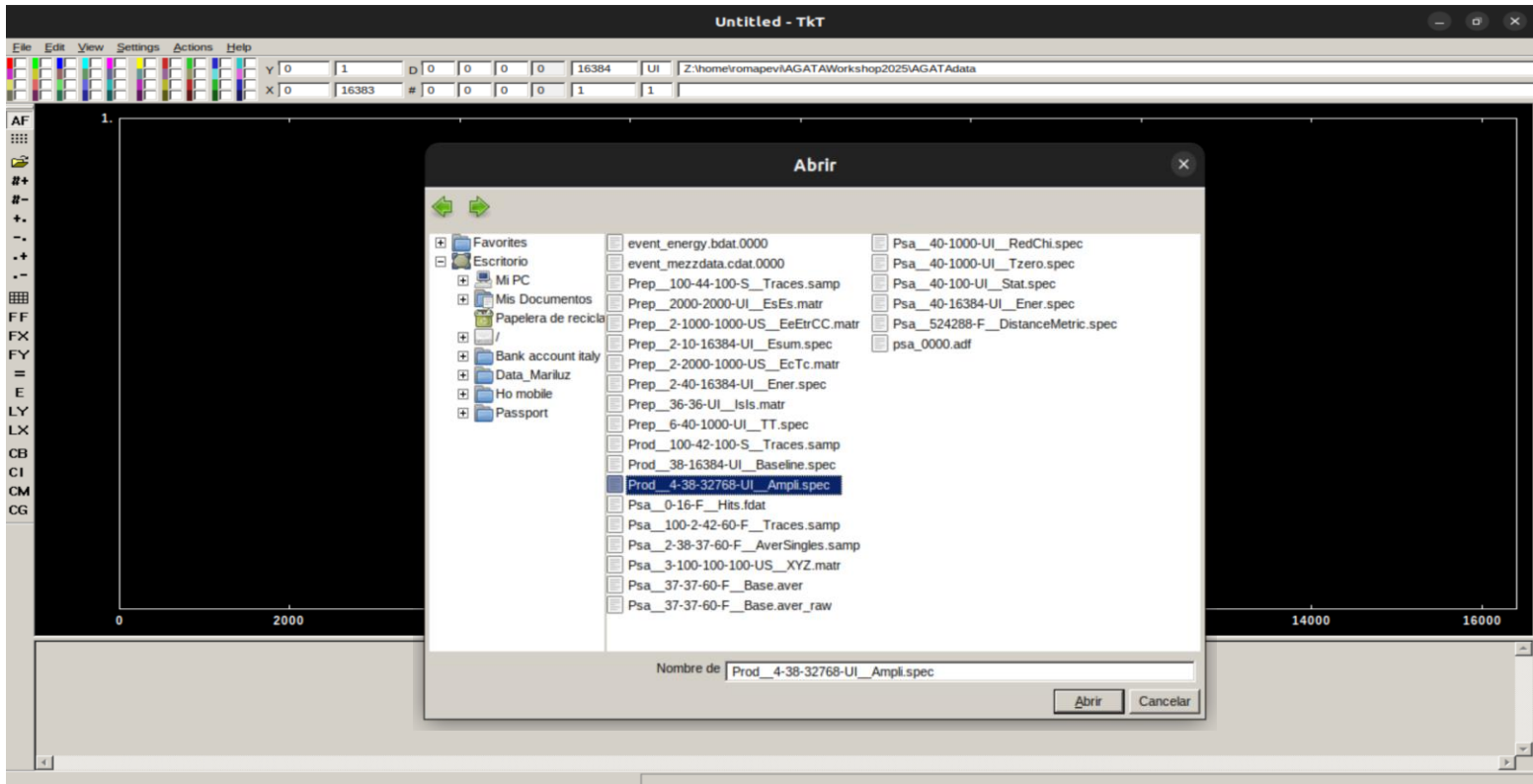
Execute TkT → click on the icon  → select a file (e.g. Prod\_\_4-38-32768-UI\_\_Ampli.spec)



# TkT

## How to open a file

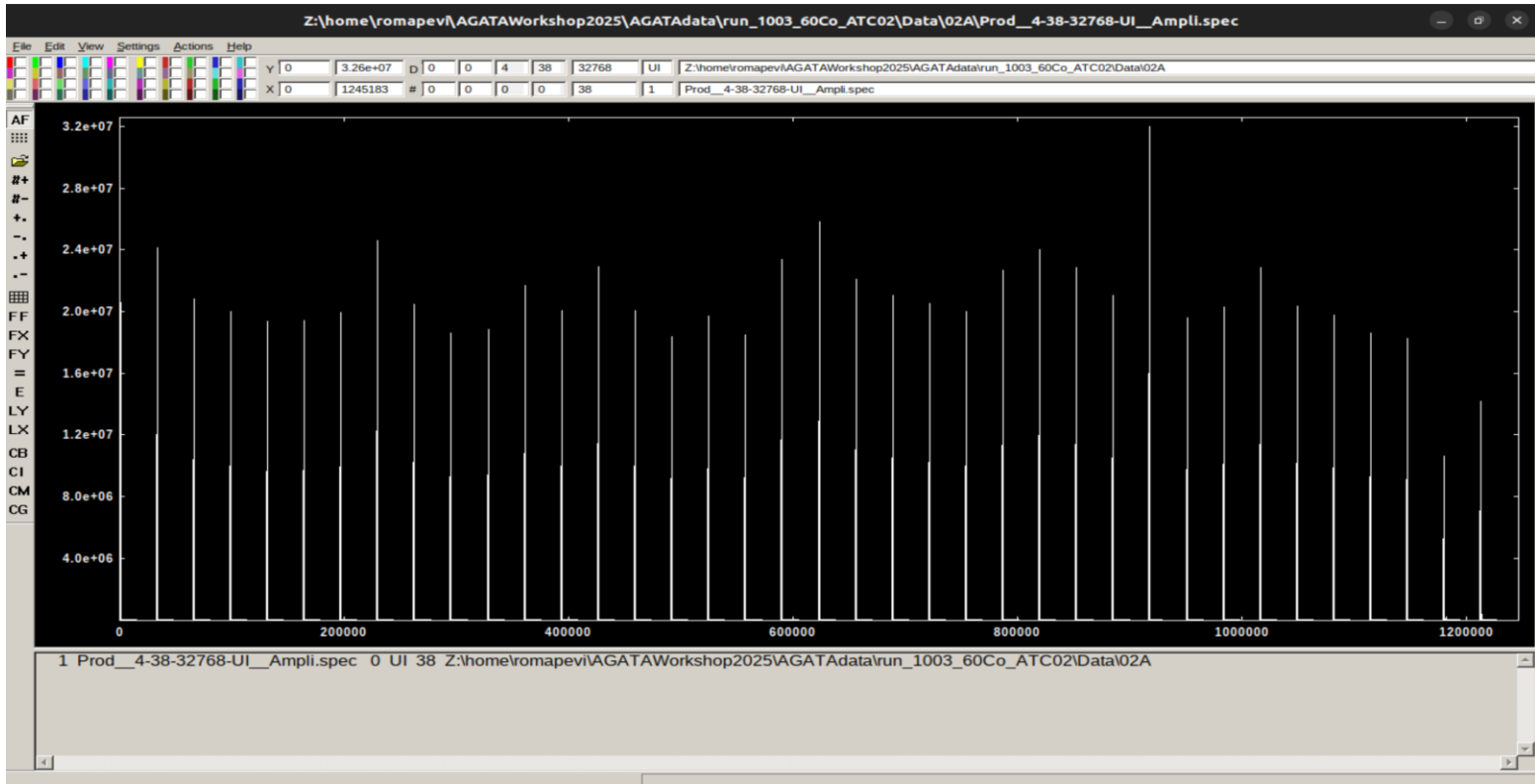
Execute TkT → CTRL+O → select a file (e.g. Prod\_\_4-38-32768-UI\_\_Ampli.spec)



# TkT

## How to open a file


Execute: `TkT PathToFile/Prod__4-38-32768-UI__Ampli.spec`





# TkT

## How to split the screen

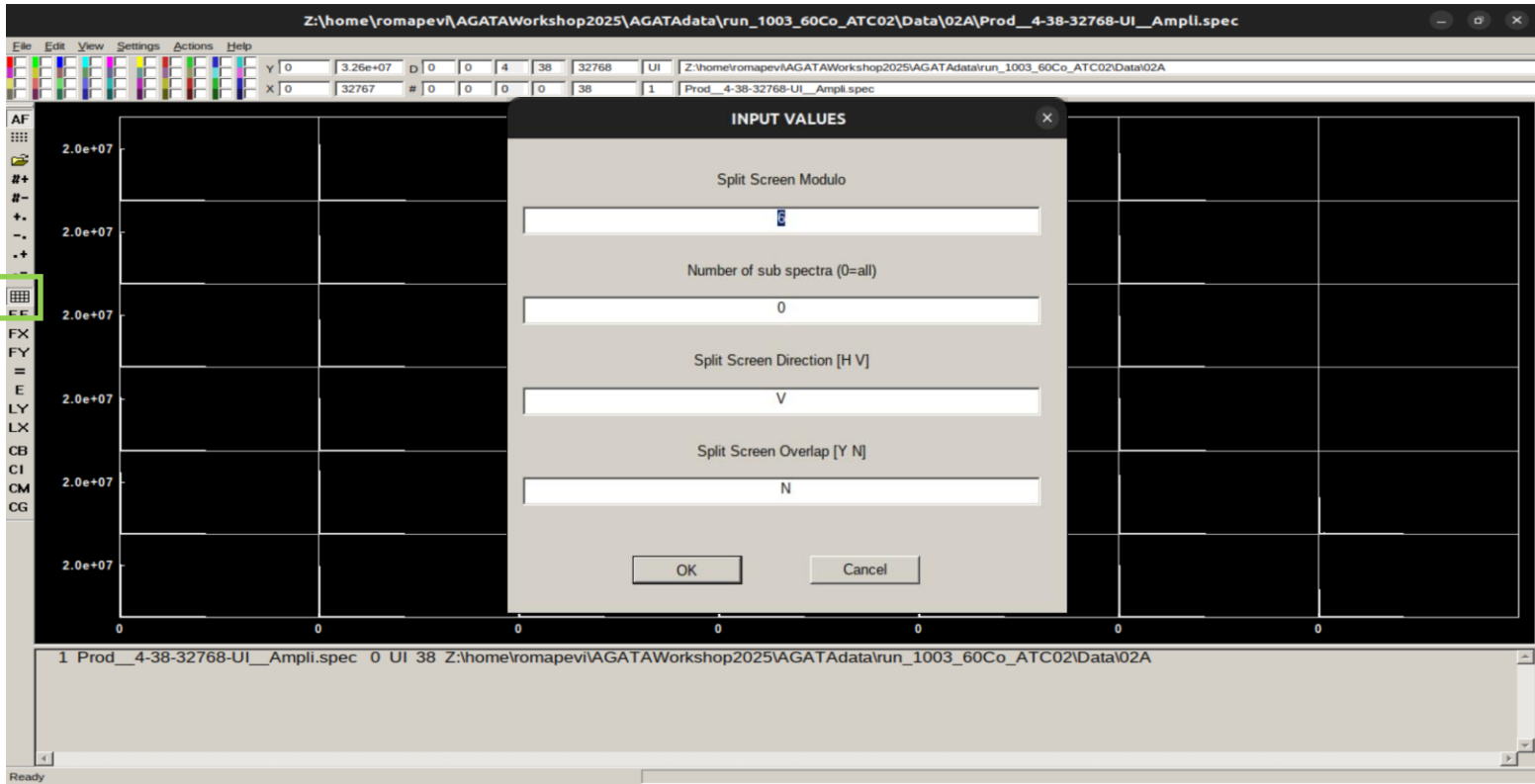
Click on the icon 



## How to split the screen

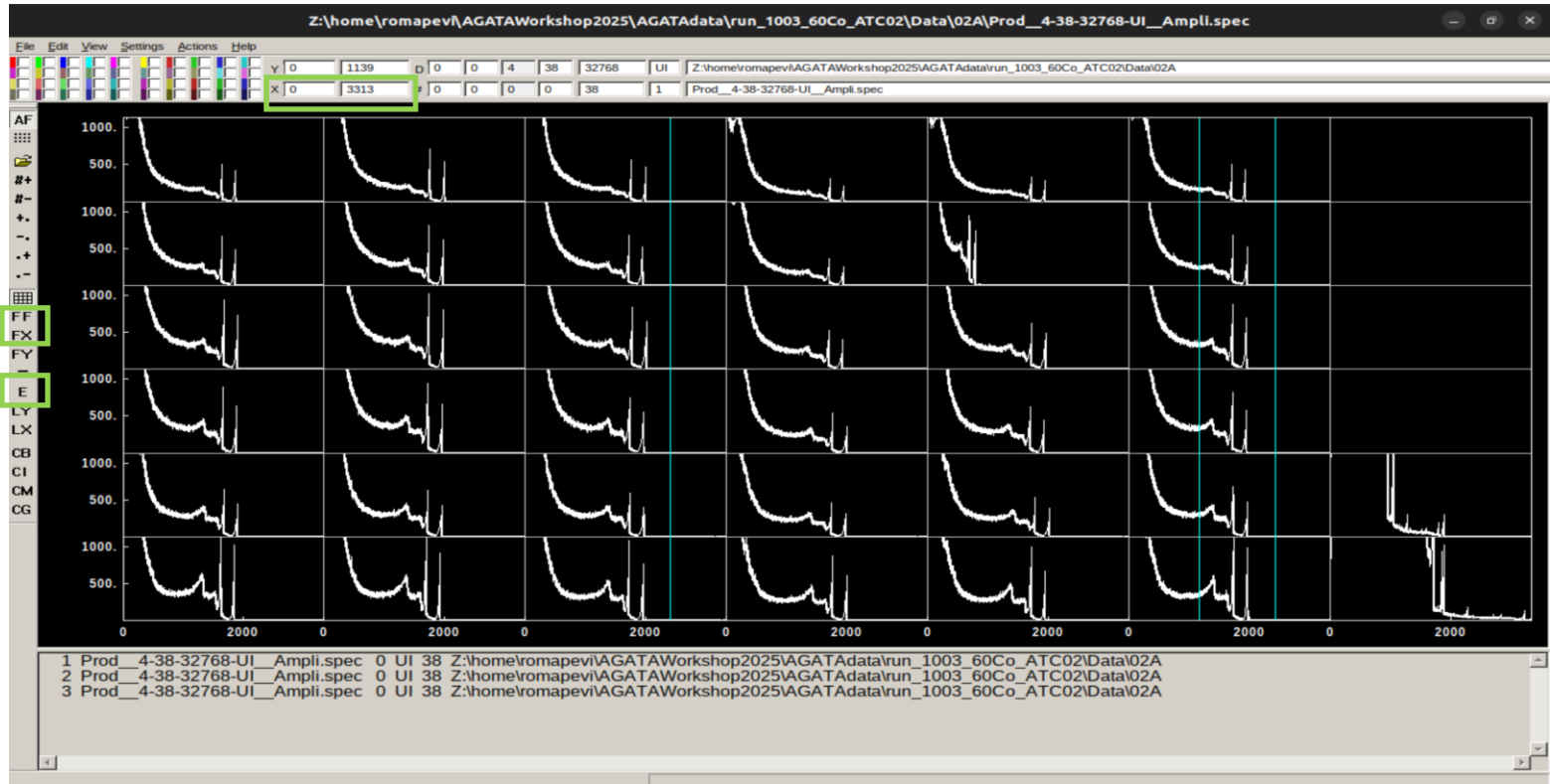
By default, the screen splits in modules of 6 and displays all the spectra. You can change the settings with the command CTRL+ 

CTRL+









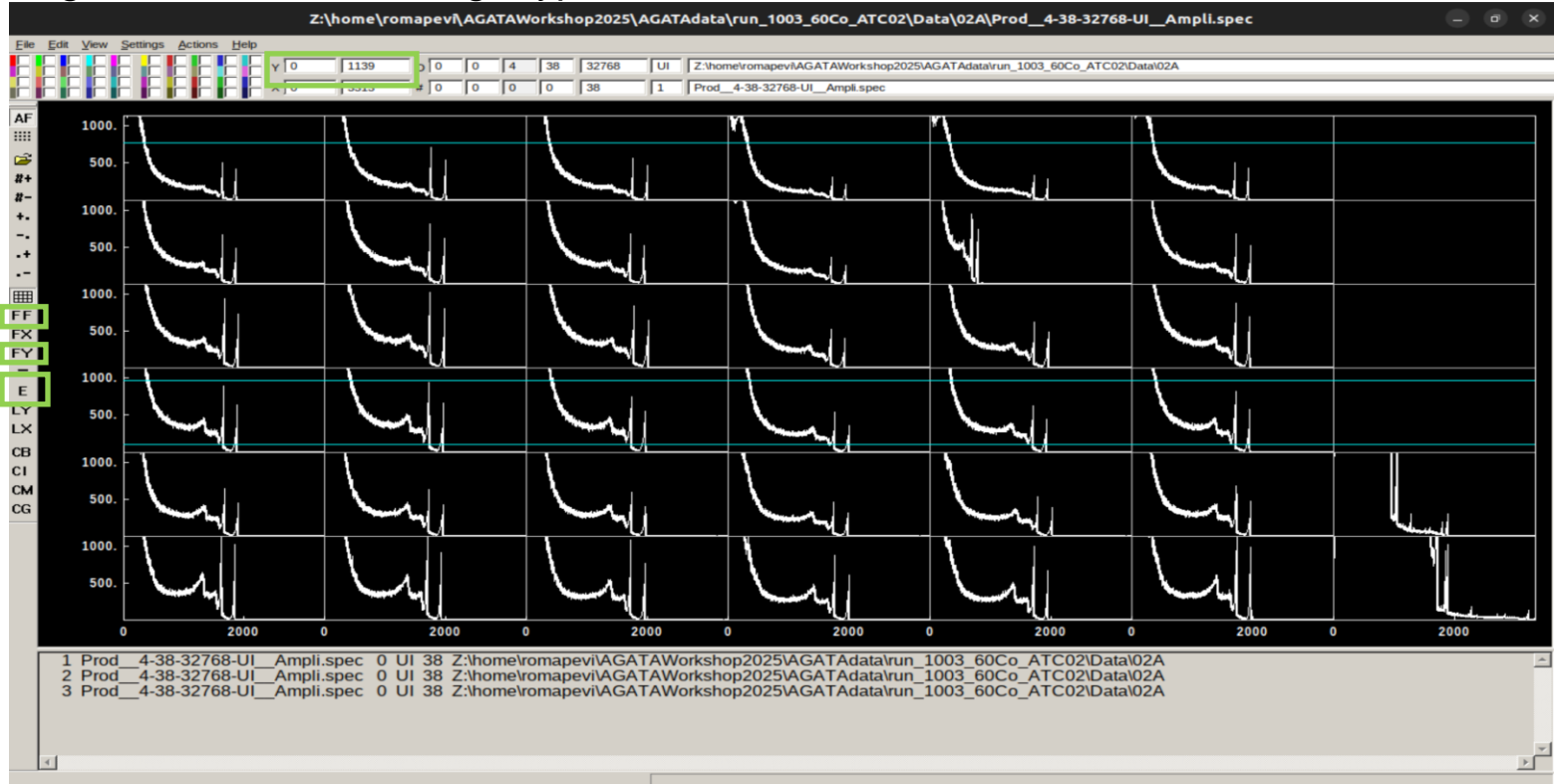
## How to zoom in the X axis

Change the X axis ranges / click on one part of the Energy spectra | and type E or click **E** / click on both sides of the region to zoom || and type E or click **E** . To go back to the full range type FX or FF **FF**  
**FX**



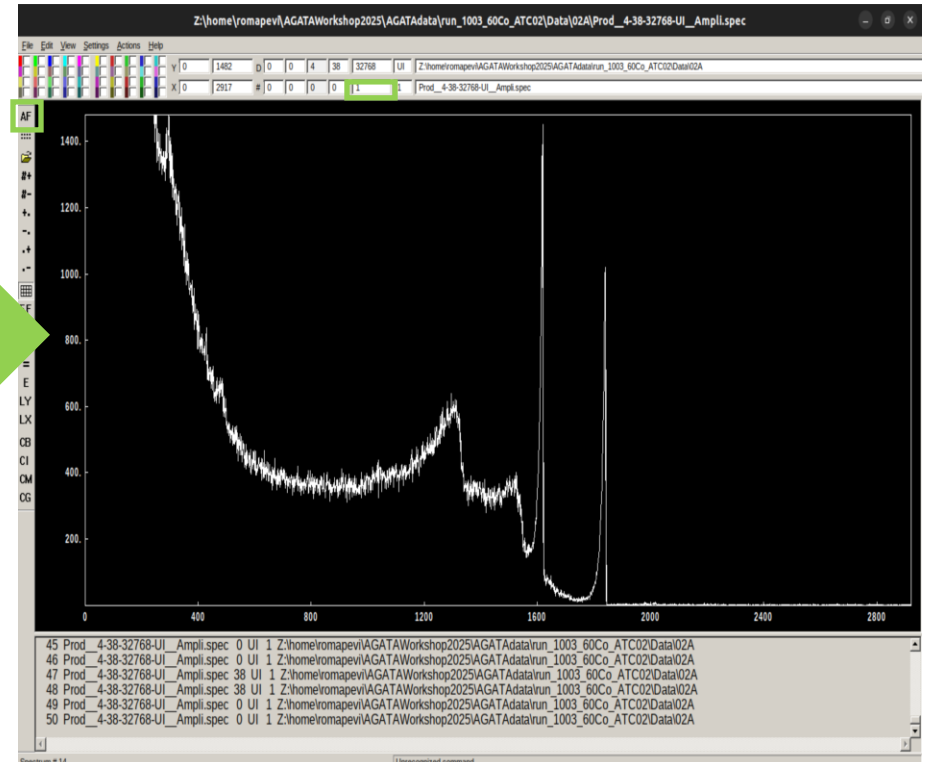
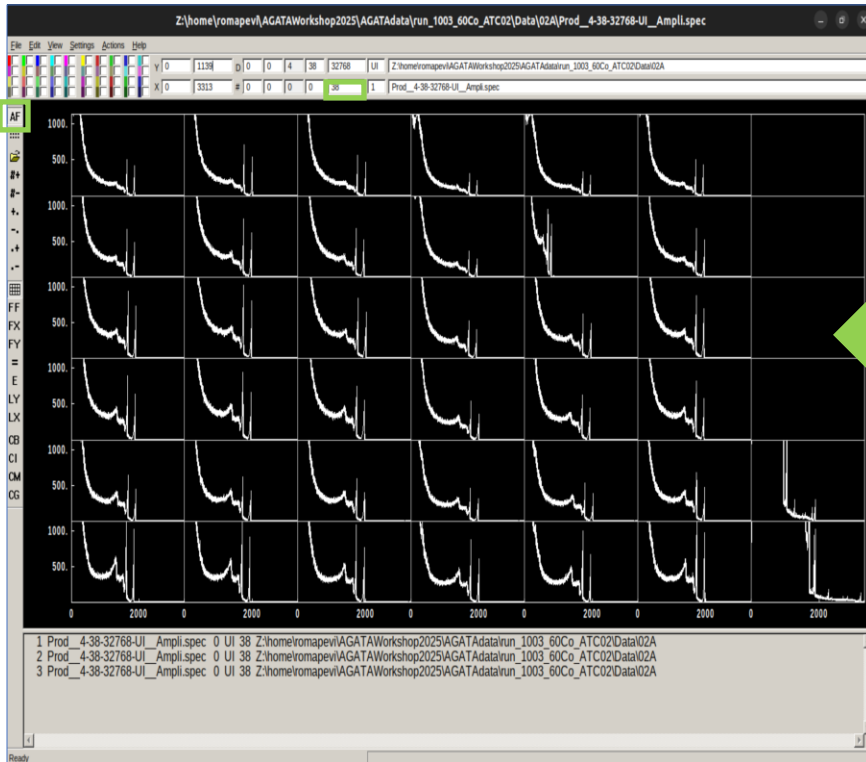
## How to zoom in the Y axis

Scroll up and down with the mouse / change the Y axis ranges / SHIFT + click on one part of the Energy spectra  and type E or click  / SHIFT+ click on both sides of the region to zoom  and type E or click . To go back to the full range type FF or FF  



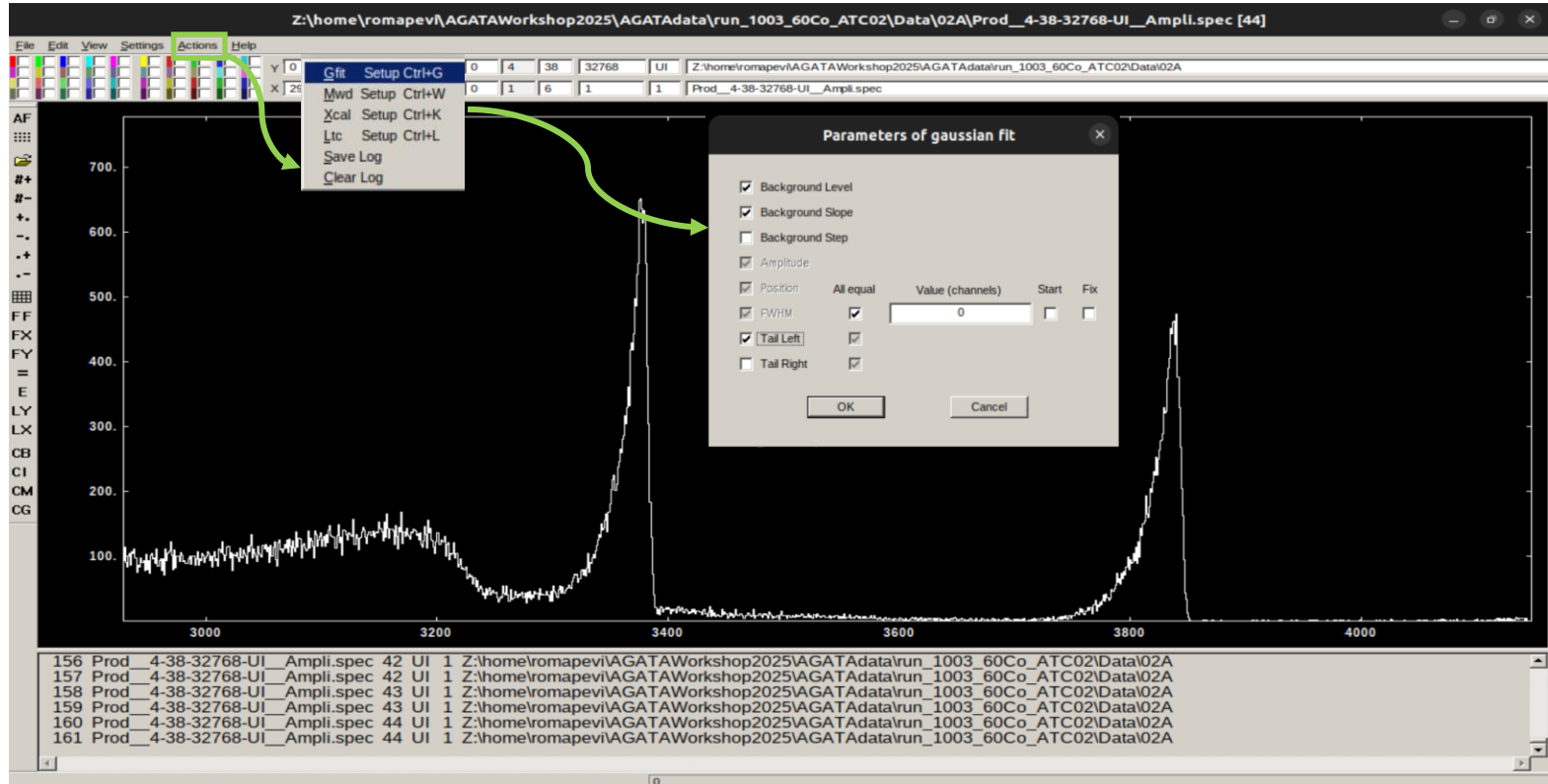
## Display modes, all spectra <-> one spectrum

Deactivate AutoFormat (CTRL+ **AF** ). Switch between the display modes with CTRL + click on the box or by changing the number in the box



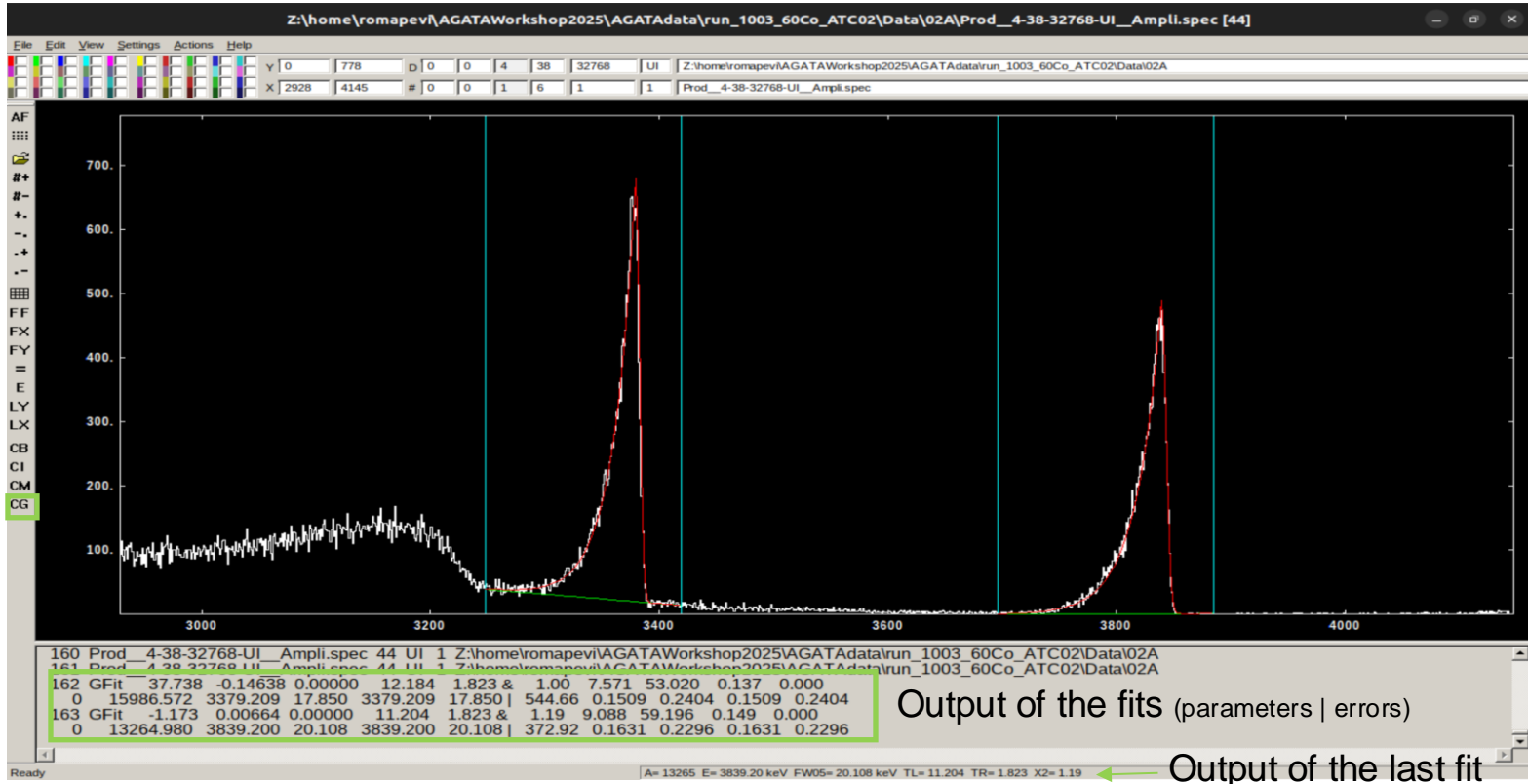
## Fit

Display 1 spectrum. Set up the parameters of the gaussian fit (Actions → Gfit or CTRL+G).



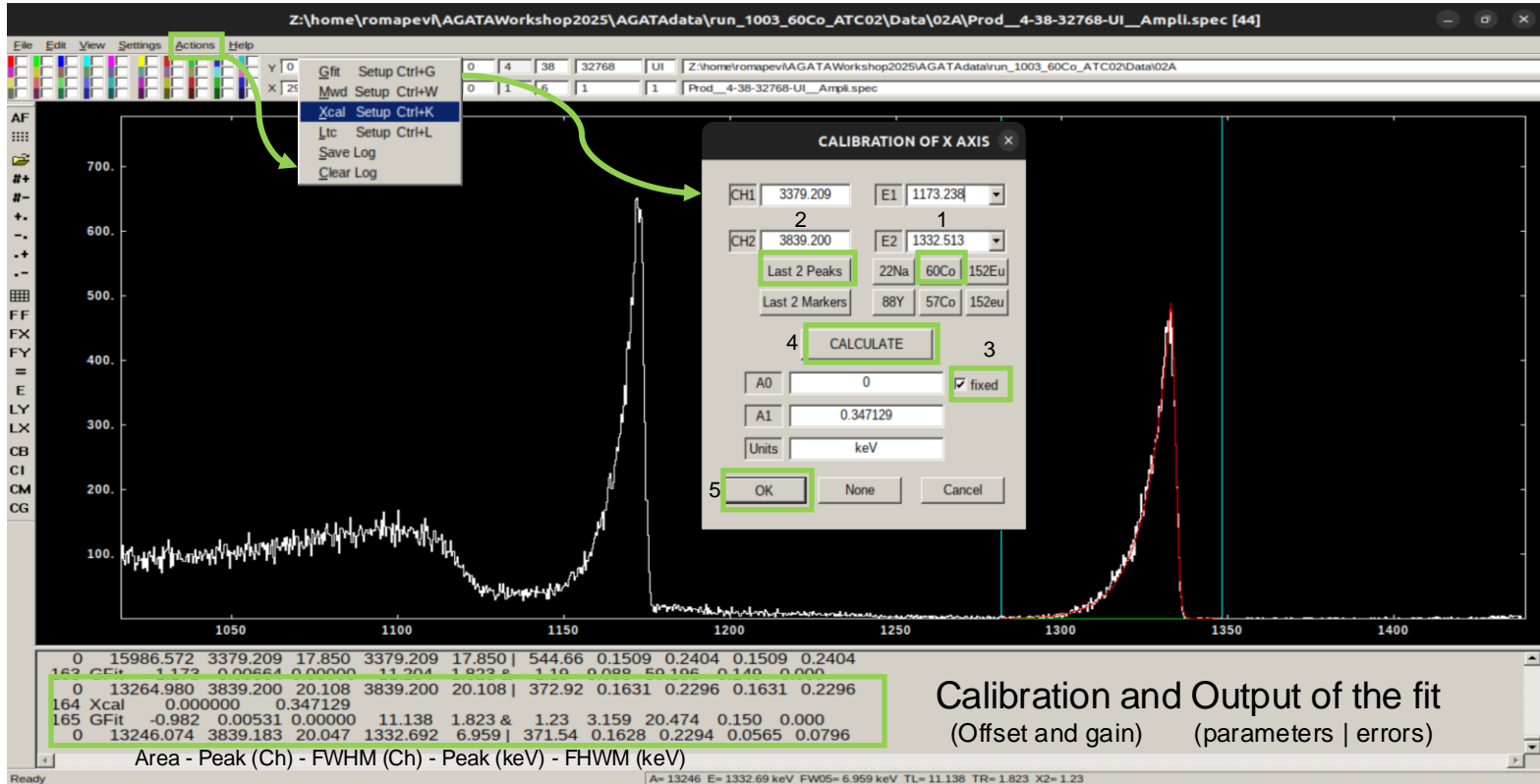
## Fit

Display 1 spectrum. Set up the parameters of the gaussian fit (Actions → Gfit or CTRL+G). Click on both sides of the peak to select the range | |, type CG or click **CG**. To erase the fit type CZ



## Fit + Calibration

After fitting the 2 peaks, open the calibration panel (Actions → Xcalt or CTRL+K).  
 Select the source (1), select Last peaks (2), fix the offset (3) and calculate (4). OK to apply the calibration factor (5).

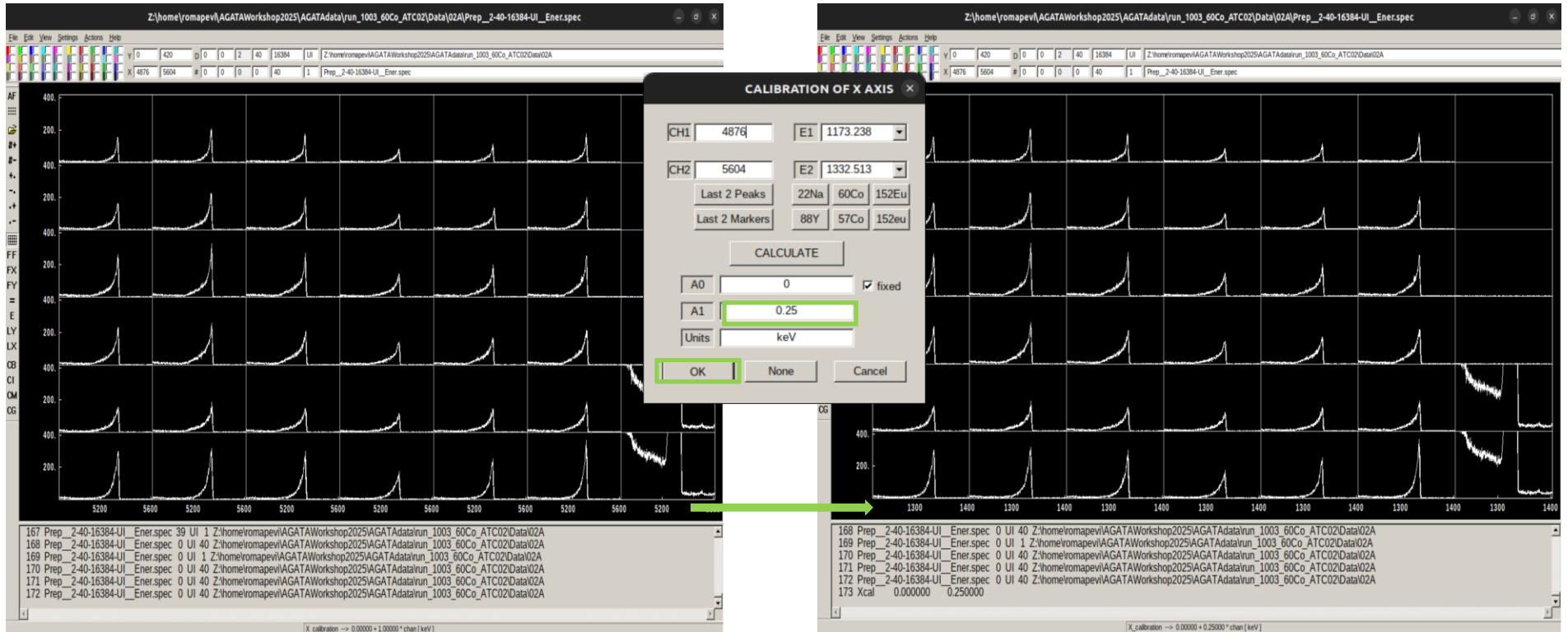


Calibration and Output of the fit  
 (Offset and gain) (parameters | errors)



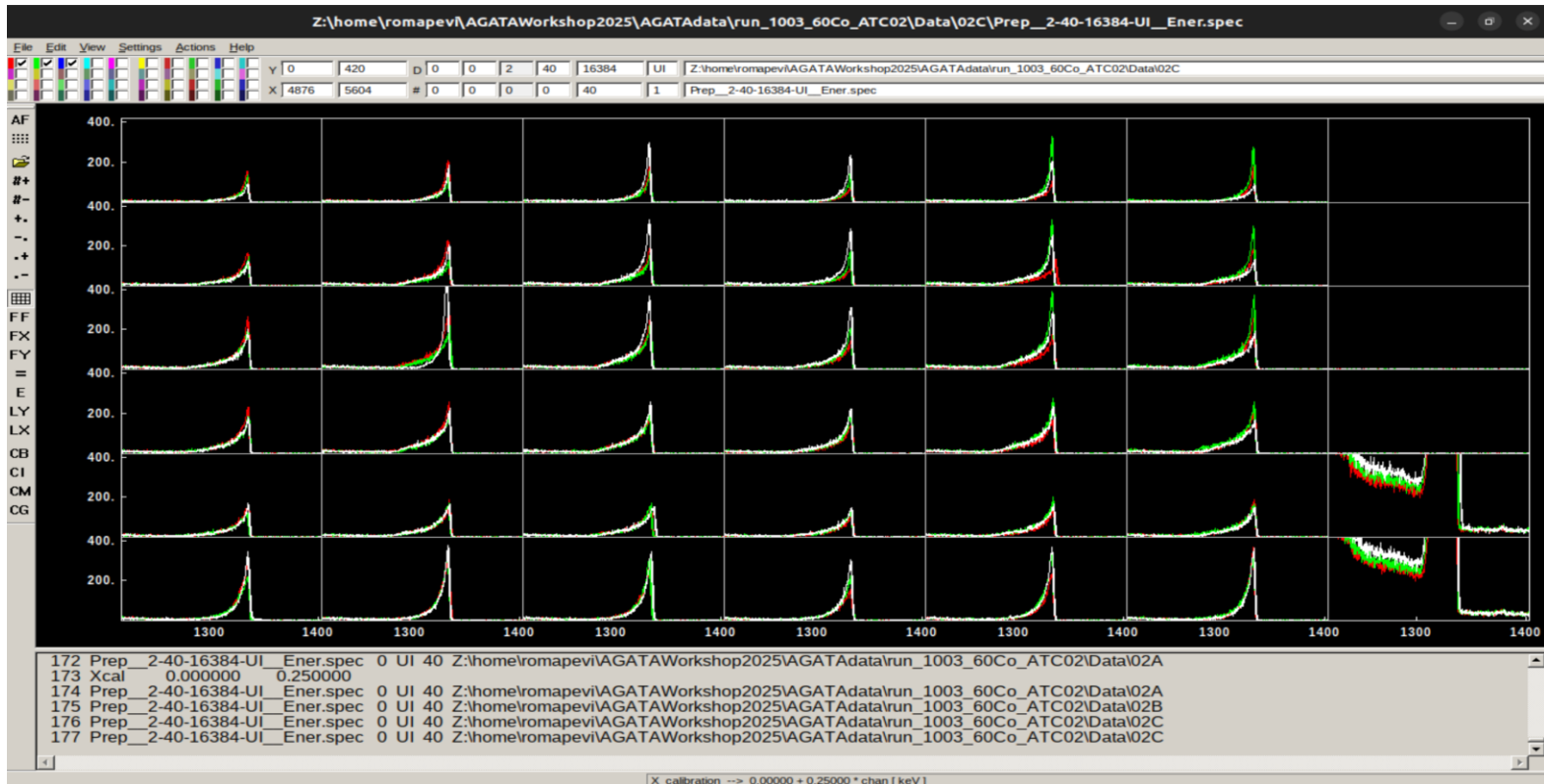
## Calibration factor

For the spectra files calibrated in energy apply the gain factor selected in the gen\_con.py. Typically gain 4 (0.25) or gain 2 (0.5).

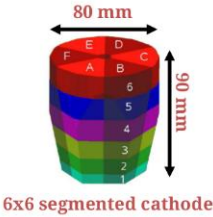


## Overlap spectra

Store spectra for different detectors in the boxes by clicking in the box after changing to another detector.  
**Shortcut:** type in the path box 02 [A, B, C] and press enter

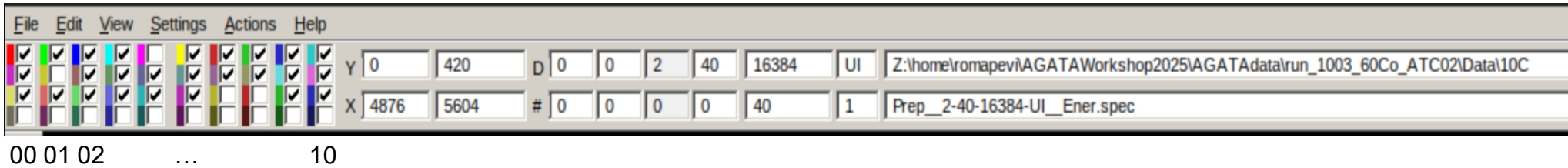


## Overlap spectra



### Shortcuts

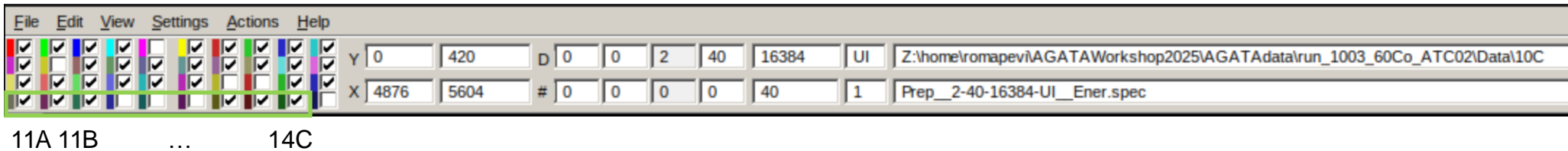
- `[,00,01,02,....,10](ABC)` → `[]` store #dets in the same raw, `()` store ABC in the same column



- **Export the command in your .bashrc and use it typing \$+the number assigned**

`export TKTSYMBOL_8='[,00,01,02,04,05,06,07,08,09,10](ABC)'` → `$8`

`export TKTSYMBOL_9='[,11A,11B,11C,13A,13B,13C,14A,14B,14C]( , , ,.)'` → `$9`



# TkT

## Short keys:?,s?,w?

## w? → Help for operations on waves

TkT

The multi-key commands Wx[n] perform Moving Window operations on sampled waveforms (Wx spectra)  
Parameters given in n'th column of Moving Window Dialog  
CTRL<W> opens the Moving Window Dialog

WAn	Average [n=1,2]	
	Sum [n=3,4]	
WB	Base line follower (continuous average)	
WC	CFD timing filter	
WDn	Gast's Deconvolution [n=1,2]	
WE	Trapezoidal filter = WD1+WA1	
WH	Info and calcs. for top and background windows	
Win	IIR CR [n=1,2]	
	RC [n=3,4]	
WJn	Jordanov's DSU [n=1,2]	HPD [n=3,4]
	DDC [n=5,6]	DPZ [n=7,8]
WLn	Leading edge discr. on Left [n=1,2]	
	on Max [n=3,4]	
WM	Median filter passed over the waves	
WO	Subtract offset based on BL width parameter	
WS	Slope condition counting	
WW	Test of SortMars analysis	
WZ	Test of SortMars analysis with integer arithmetic	

Acceptar

TkT

## s? → Help for operations on spectra

The multi-key commands Sxn perform operations on the spectrum  
SR means selected region (expanded or between L2Ms)

SAC	Autocorrelation of SR (very slow for large spectra)
SADC	Test ADC non-linearity, starting from channel-width spectrum
SCx	Constant operation 'SR x= C'; x one of '+ - * /   & ^ < > %' (%0-->int); constant C taken from dialog
SCM	Masking of data digitized with 14 bits (SC& 16383 SC^ 8192)
SD	running Differentiation in SR
SFn	Recursive Digital Filter n applied to SR
SF[AS]	Moving Average or Moving sum in SR
SFC	Complex conjugation of SR interpreted as (Ampl,Phase)
SFD	Dir. Discrete Fourier Transform of SR
SFI	Inv. Discrete Fourier Transform of SR
SF[MN]	Median Filter in SR. Filter width 3(M) or 5(N) points
SF[RZ]	Set to random or zero the phases of SR interpreted as (Ampl,Phase)
SFWn	0-Square 1-Blackmann 2-Hamming 3-Hanning 4-Bartlett windowing in SR
SF[*]	Operation on 'Fourier' spectra in the (Ampl,Phase) representation
SGx	Generate function in SR: Band Delta Exp Gauss bRestorer Sin Trapezoid
SI	running Integration of SR
SKnn	Convolution with kernel taken from stored spectrum nn
SLB	transform to Bode plot SR
SL[XY]	x or y axis mapped to log10 in SR
SLI	10^y in SR
SMn	Smooth SR n times [n=0-9, (0->10)]
SN[UGS]	add Noise to SR [Uniform Gaussian SpecSigma]
SN[A]	Normalize spectrum to totArea or maxAmplitude given in dialog
SP[CDERS]	Pack, Decimate, Filter-out, Reshuffle channels or spectra
SPn	pow(spectrum, n) of SR [n=0 --> ask, n=1 --> abs[]]
SQn	pow(spectrum, 1/n) of SR [n=0 --> (int), n=1 --> 1/0]
SRnn	Recall stored spectrum nn
SR[BED]	Running linear[B] fit, exponential[E] fit, standard deviation[D]
SR[LR]	Rotate Left/Right channels of SR
SSnn	Store as spectrum nn
SSA	Spectrum shift and accumulate
SSBn	Swap bytes [n=2,4]
SS[LR]	Shift Left/Right channels of SR
SS[OS]	Swap Segments Order, Swap incrStep Sets
SSK	Shift spectrum according to calibration coefficients
SS[HV]	Swap spectrum Horizontally/Vertically
SS?	List of stored spectra
SX	Distribute y values on x axis SR
SY	Distribution of y-values of SR
SV[AN?]	Set/reset/query visibility of spectra
SVnn	Toggle visibility of stored spectra
SWA[IF]	Write spectrum as ASCII text in Integers or Floats
SW[CSILFD]	Write spectrum as binary 8-,16-,32-,64-bit integers, Float or Double
SZnn	Remove Stored Spectrum nn
S0	Value of first channel in SR set equal to value of second one
S[+*/@S&]nn	Spectrum operation SR = ss[nn] op SR @=LUT \$=accumulate &=nonzero
S-[ST]	Sum of Stored Spectra or Segment Traces
S=[01FB]	Replace channels between L2Ms with 0, 1, the last fit or background

TkT

## ? → Help

Set Vertical Markers with <spacebar> or <left\_mouse>  
Set Horizontal Markers with SHIFT<spacebar> or SHIFT<left\_mouse>  
LnMs stays for 'last n markers' n stays for an integer number between 0 and 9

A	Auto fit a peak between L2Ms
B	Set a B marker
CA	Fit a flat line between L2Ms or in B regions
CB	Fit a stright line between L2Ms or in B regions; CRTL<CB> --> flat line
CE	Exponential fit between 2 markers (with 4 markers, last 2 define background region)
CG	Gauss fit between L2Ms; CRTL<CG> --> toggle left tail; SHFT<CG> --> show fit results
C[J]	Integration between L2Ms; Background subtracted if 2 R-markers and pairs of B-markers
CM	First moments between L2Ms; background from CB or first-last; CRTL<CM> no background
C[PO]	Peak search using first(P) or second(Q) derivative of spectrum
CR	Residues of last fit between L2Ms
CS	Sinusoidal fit; initial period given by 2 markers (if 4 markers, last 2 define fit region)
CZ	Remove drawing of last fit
DD	Toggle direction of increments/decrements
DFF	Data format float F 100 1000
DFK	Toggle value of K between 1024 and 1000
DFL	Resize spectrum K-length
DFT	Data format WT2 1 100000
DMn	Define macro n (0...9) DM? list macros
DM[GHV]	Define marker G, H or V
DSn	Define symbol n (0...9) DS? list symbols
DSS	Define Split Screen
DUG	Define User Grid
D[XV]	Print L2Ms and their difference
E	x-Expand between L2Ms
F[XV]	Expand to full-X, full-Y, full-XY
F[KLN]	FX for the 1st, last or dialog-given "k"
G	Set a G marker
KA	X-calibration from 2 largest peaks
K[MP]	X-calibration from L2Ms or last 2 peaks
K[ZL]	X-calibration remove or reset to last
LT	Long traces capture from ggp (same as SCC)
L[X]	Toggle Lin/Log X or Y scale (<LX> <LY>)
Mn	Execute macro n (0...9) SHFT<M> execute macro 1
N	Re-read spectrum
T	Trigger on next peak; SHFT<T> on previous peak (level from cursor)
V[LV]	List channels or View value of current channel
Y0	Ymin set to zero
Y[OU]n	Y-expand to selected decimal fraction
Z[BGHV]	Delete B G H V markers
ZS	Erase spectrum between L2Ms
^K	Open X-calibration dialog
^O ^R	Read again present spectrum and redraw or overdraw
<ALT>+n	Execute macro n
1 2	Next/Previous #spectrum (<#++> <#-->) (<SHFT> redraws)
3 4	Increment/Decrement Spectrum Name (<SHFT> redraws)
5 6	Increment/Decrement Spectrum Type (<SHFT> redraws)
= . < >	Redraw, Center spectrum at mouse position, Left / Right shift
Sx	Operations on spectra (S? --> help)
Wx	Operations on waves (W? --> help)

**Thank you!**

**AGATA Analysis Workshop 2025  
PostPSA Calibration. Hands on**

**R.M. Pérez-Vidal**

14/01/2025, Lyon

**Questions?**

# AGATA Crystal lookup table

Position	ATC	Crystal			Installation date
		A	B	C	
00	12	006	005	001	01/04/2022-15/01/2024
	<b>18*</b>	017	018	018	15/01/2024
01	10	011	006	012	01/02/2022- 03/04/2024
	<b>13</b>	003	016	015	03/04/2024
02	<b>17</b>	016	017	013	01/02/2022
04	11	004	004	010	01/09/2022-04/04/2024
	<b>7*</b>	007	014	003	04/04/2024
05	<b>09</b>	001	001	004	?
				006	01/05/2022
06	06	008	009	002	?
	<b>19</b>	018	012	014	01/09/2022-14/12/2023
07	14	014	010	016	01/03/2022-18/09/2023
	<b>2</b>	019	019	020	18/09/2023
08	<b>3</b>	002	007	007	01/03/2022
09	18	017	018	018	01/03/2022-19/10/2023
	<b>14*</b>	014	010	016	19/10/2023
10	15	013	015	011	01/03/2022
11	<b>1</b>	010	011	009	01/04/2022
13	<b>19</b>	018	012	019	01/09/2022
14	7	015	014	008	?
	<b>20</b>	007	014	003	01/09/2022-01/01/2023
		009	020	005	01/01/2023

\*2nd time in the array

