



AGATA Analysis Workshop 2023

PostPSA calibration. Hands on

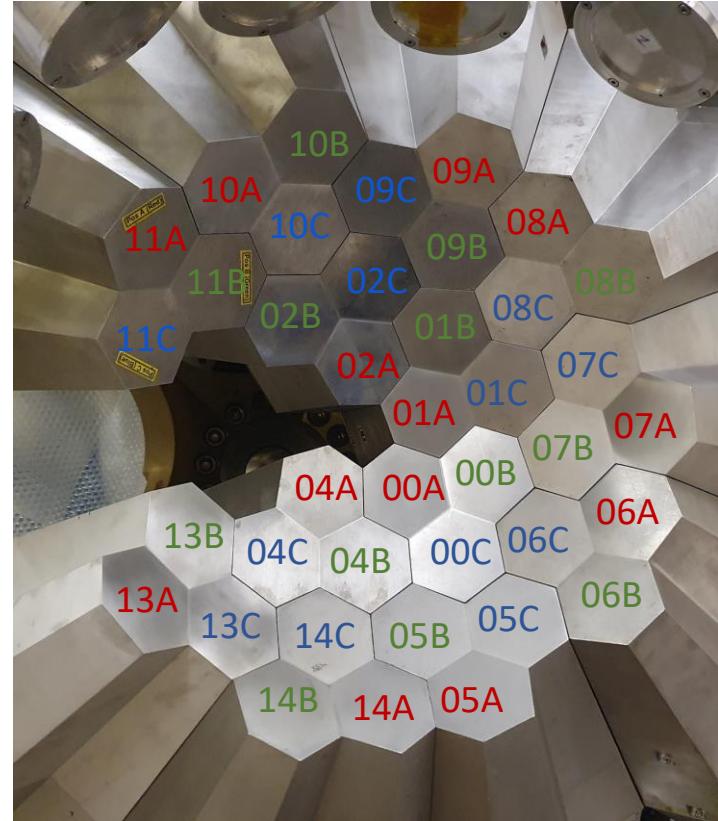
R.M. Pérez-Vidal

12/09/2023, Legnaro

Hands On

AGATA configuration

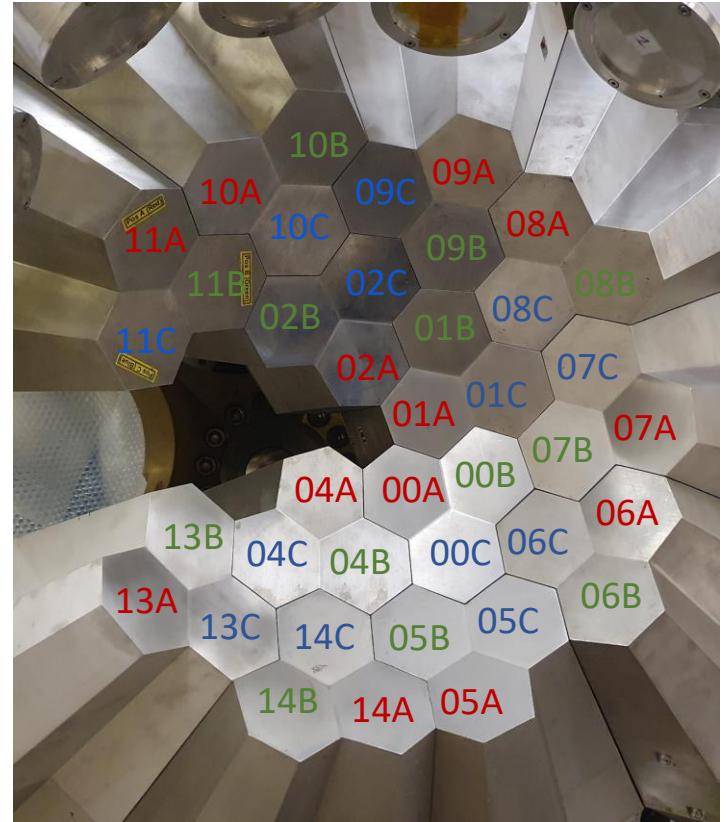
- AGATA data taken in May
- Counting rate per crystal : 500-900Hz
- MWD-risetime: 2.5 μ s
- Position: Nominal
- Backward angles
- Data sets with ^{60}Co and ^{152}Eu



Hands On

Data sets in example directory

- **run_1005_60Co (40Gb)**
 - Neutron damage correction
 - 4 ATC: 00 01 02 11
- **run_1002_152Eu (36Gb)**
 - Energy recalibration
 - 4 ATC: 00 01 02 11
- **run_1010_60Co (16Gb)**
 - Global time alignment
 - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 13



Hands On

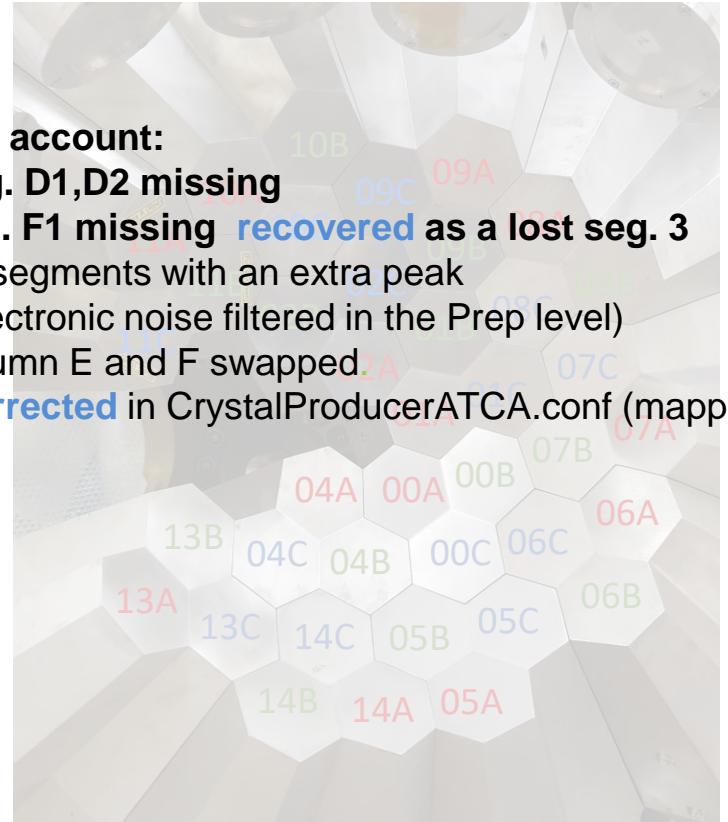
Data sets in example directory

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To be taken into account:

- 01B: **seg. D1,D2 missing**
- 11A: **seg. F1 missing** **recovered as a lost seg. 3**
all segments with an extra peak
(electronic noise filtered in the Prep level)
- 11C: column E and F swapped.
Corrected in CrystalProducerATCA.conf (mapping).

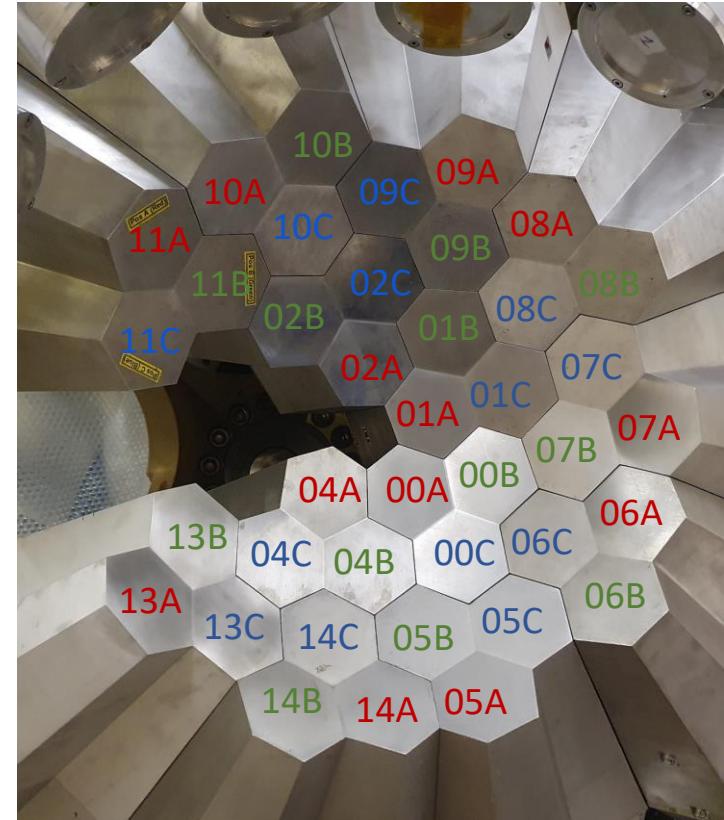


Hands On

Data sets in example directory

- **run_1005_60Co_ATC00 (14Gb)**
 - Neutron damage correction
 - 1 ATC: 00A (traces) 00B 00C
- **run_1002_152Eu_ATC00 (7.9Gb)**
 - Energy recalibration
 - 1 ATC: 00A (traces) 00B 00C
- **run_1010_60Co (16Gb)**
 - Global time alignment
 - 36 ATC: 00 01 02 04 05 06 07 08 09 10 11 13

POS	ATC	Crys	Capsule
		A	6
0	12	B	5
		C	1



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 00A (A006)
- 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
 - Neutron.sh
 - Recalibration.sh some analysis scripts
 - GlobalTime.sh

Hands On

Files in example directory: run_1005_60Co

Neutron Damage correction

- 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

- **event_mezzdata.cdat.0000**
- **event_mezzdata.cdat.0001**
- **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_1002_152Eu

Recalibration

- 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**
- **psa_0001.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 CRYS 00A 00B 00C  
  
    Chain 4      CRYS  
    Producer  
    CrystalProducerATCAFitter  
    PreprocessingFilterPSAFilter  
    PSAFilter  
    Consumer BasicAFC  
ENDLOOP
```

Hands On

Files in example directory: Topologies

Topology_Global.conf

```
LOOP CRYS 00A 00B 00C
```

```
    Chain 3      CRYS
    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 Crys 00A 00B 00C

 Chain 3 Crys
 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

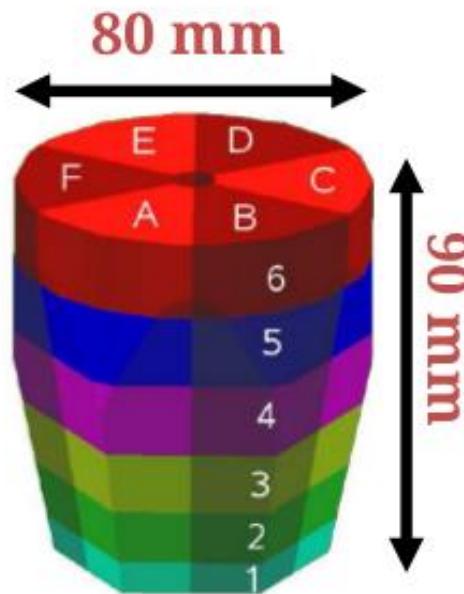
Useful programs. PostPSA Filter

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

TkT

Channels correspondence after Replay



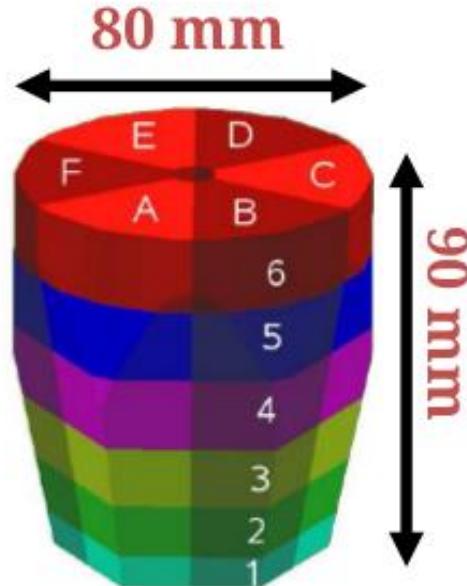
A	B	C	D	E	F	CC
6	5	11	17	23	29	35
5	4				34	
4	3				33	
3	2				32	
2	1				31	37
1	0	6	12	18	24	30
0						36

1 Low gain

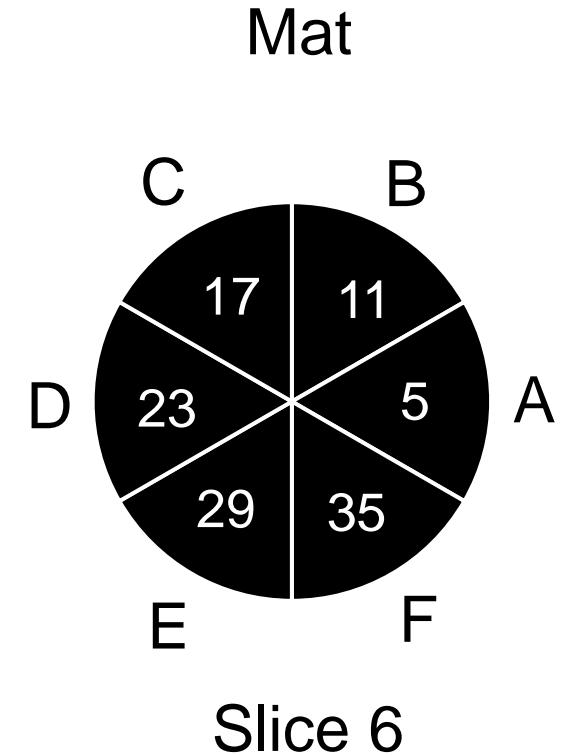
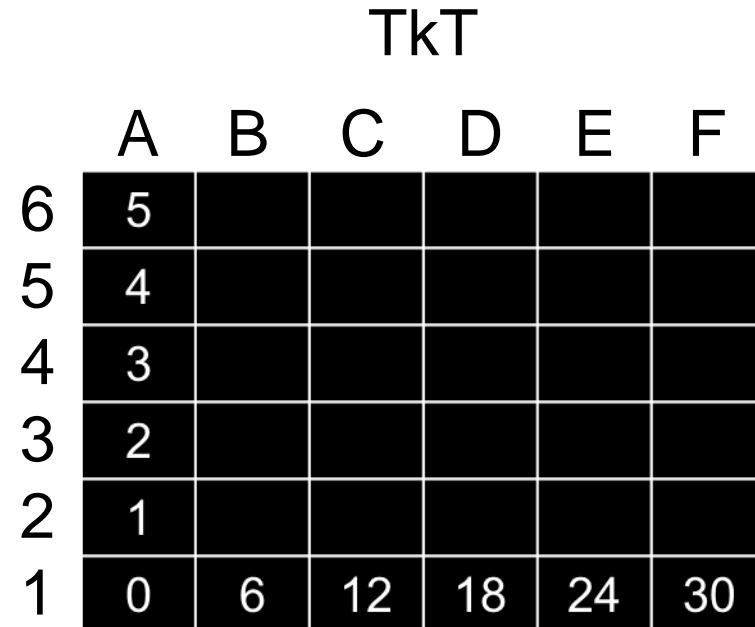
0 High gain
& trigger

TkT & Mat

Channels correspondence after Replay

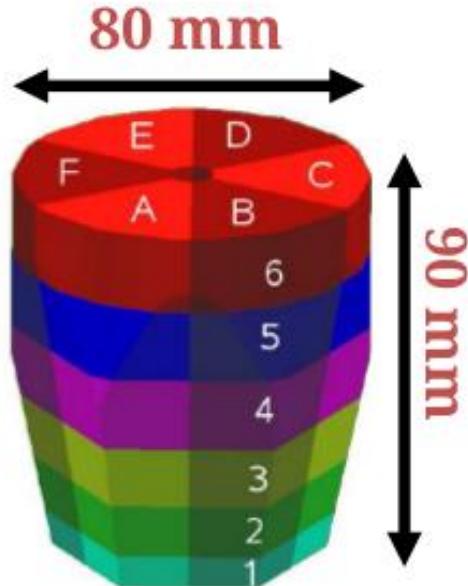


6x6 segmented cathode



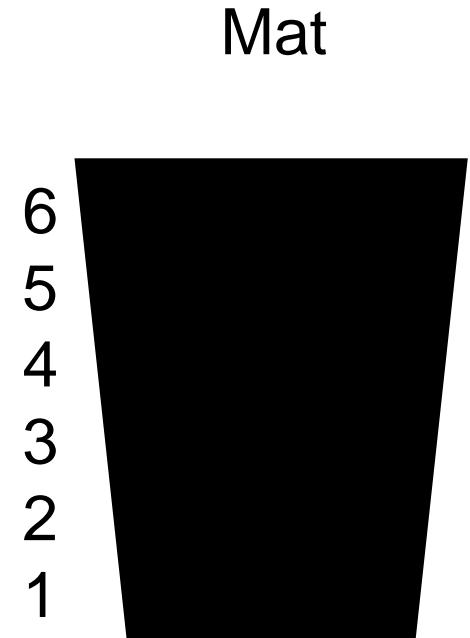
TkT & Mat

Channels correspondence after Replay



6x6 segmented cathode

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



TKT

Short keys:?:s?,w?

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)

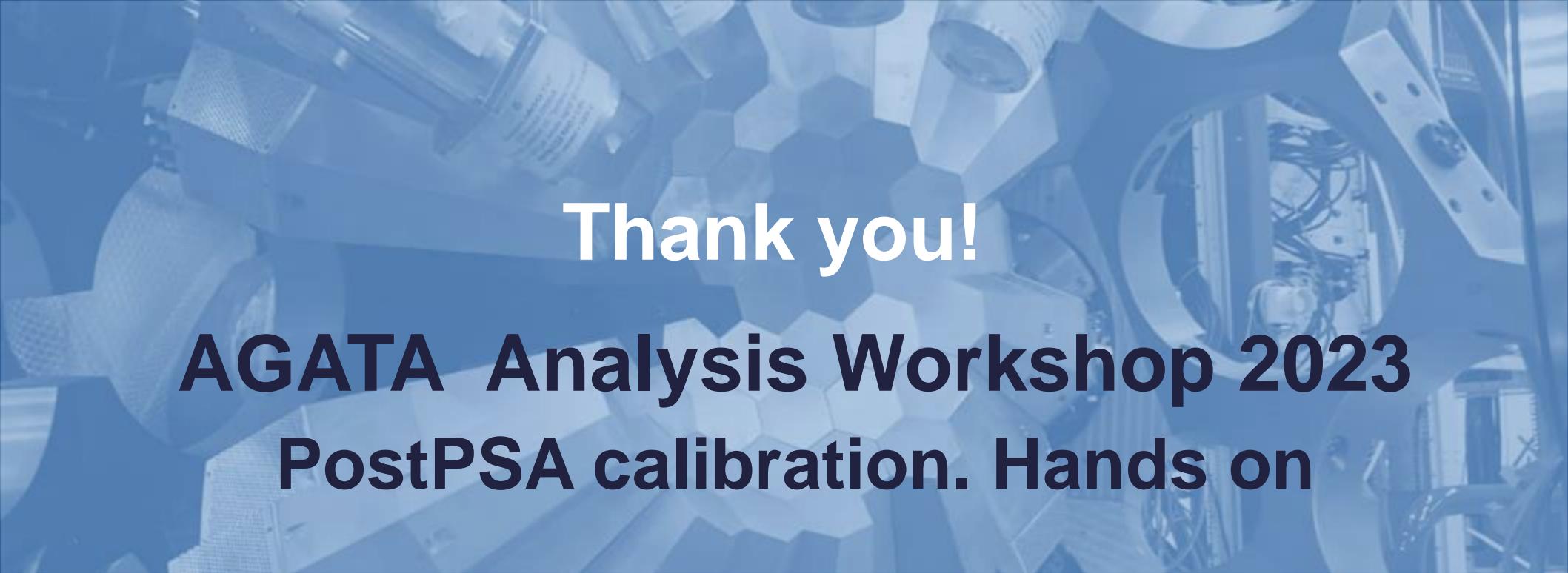
SAC	Autocorrelation of SR (very slow for large spectra)
SADC	Test ADC non-linearity, starting from channel-width spectrum
SCx	Constant operation 'SRx=C; x one of '+ - * / & ^ < > %' (%0-->int); constant C taken from dialog
SCM	Masking of data digitized with 14 bits (SC& 16383 SC^ 8192)
SD	running Differentiation 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[V]	Operation on 'Fourier' spectra in the (Ampl,Phase) representation
SGx	Generate function in SR: Band Delta Exp Gauss blRestorer 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[JUGS]	add Noise to SR [Uniform Gaussian SpecSigma]
SN[AY]	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[+.*/@\$]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

Aceptar

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 straight 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; SHIFT<CG> --> show fit results
CIJJ	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[PQ]	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(XYF)	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[XY]	Toggle Lin/Log X or Y scale (<LX> <LY>)
Mn	Execute macro n (0...9) SHIFT<M> execute macro 1
N	Re-read spectrum
T	Trigger on next peak; SHIFT<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 (##+> <#->) (<SHIFT> redraws)
3 4	Increment/Decrement Spectrum Name (<SHIFT> redraws)
5 6	Increment/Decrement Spectrum Type (<SHIFT> redraws)
= . < >	Redraw, Center spectrum at mouse position, Left / Right shift
Sx	Operations on spectra (S? --> help)
Wx	Operations on waves (W? --> help)



Thank you!

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Questions?