

Excel XPS&XAS macro suite: EX3ms

- efficient and effective analysis ever experienced -

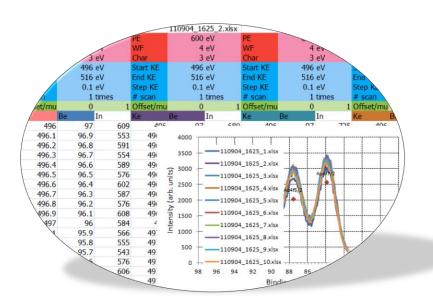
Hideki NAKAJIMA

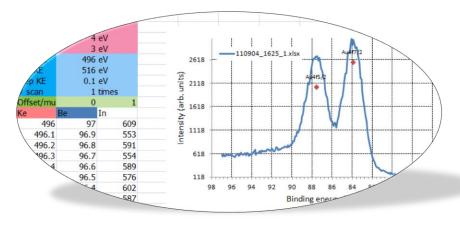
github: heitler/xps-excel-macro

26 Jun. 2019

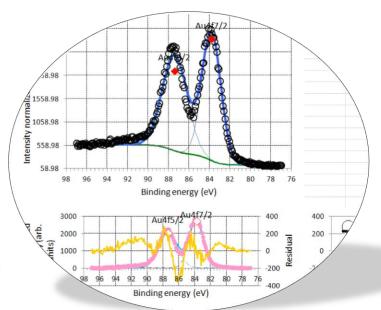
Brief

Plot with binding energy Identify elements and states





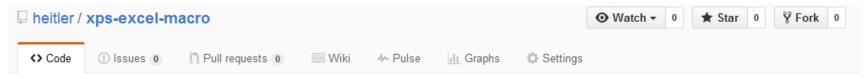
Energy and intensity calibrations



Curve fitting and BG subtraction

Preface

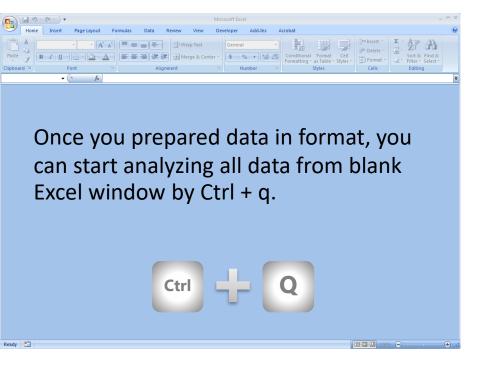
- Ctrl+Q is the first comprehensive data-analysis code for SR-XPS and XAS working on Windows Office Excel 2007 or later. Mac Excel 2016 works great now.
- Ctrl+Q will analyze the data by just pressing the shortcut key: Ctrl + q.
- It works great in publication quality.
 - Carbon, JPAP, ASS, ACS AMI, SAB, RCS Adv. etc. from the data obtained at BL3.2Ua and 5.3 in SPL
- GitHub: xps-excel-macro for updated info.



Data preparation

	Α	В
1	KE/eV	Scan#1
2	496	19.69872
3	496.1	17.95129
4	496.2	19.18263
5	496.3	18.03784
6	496.4	19.03867
7	496.5	18.80417
Q	406 G	10 22572

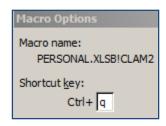
- Two-column data in the spreadsheet format are analyzed in the code.
- Worksheet is named after workbook filename.
- Syntax in A1 cell corresponds to the following.
- ISO (VAMAS) format is also acceptable.



Syntax in A1 cell	Data in column A	Data in column B
KE/eV	KE	XPS
BE/eV	BE	XPS
PE/eV	PE	XAS
GE/eV	PE	G scan
AE/eV	EE	Auger
QE/eV	mass	Q-mass
ME/eV	Position	Any

Installation of the code in Excel

- Generate Personal Workbook Macro
- Install Solver Add-in and setup its DLL
- Setup the Shortcut key in Macro Option



Compile your own database for BE and sensitivities on UD.xlsx

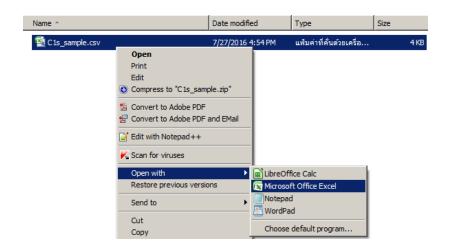
XPS worksheet

	Α	В	С	D
1	Element	Orbit	BE(eV)	ASF
2	C	1s	284.6	1
3	0	1s	532	2.93

AES worksheet

4	Α	В	С	Q
1	Element	Auger	KE(eV)	RSF
2	C	KLL	266	0.6
3	0	KLL	506	0.96

C1s sample csv or txt

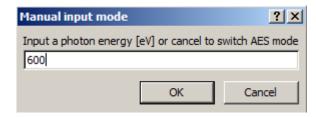


	Α	В
1	KE/eV	PE: 600 eV
2	290	1709
3	290.1	1798
4	290.2	1845
5	290.3	1722
6	290.4	1729
7	290.5	1699
8	290.6	1727
9	290.7	1693
10	290.8	1647
11	290.9	1684
12	291	1659



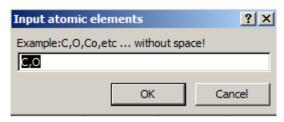
600 eV of photon energy



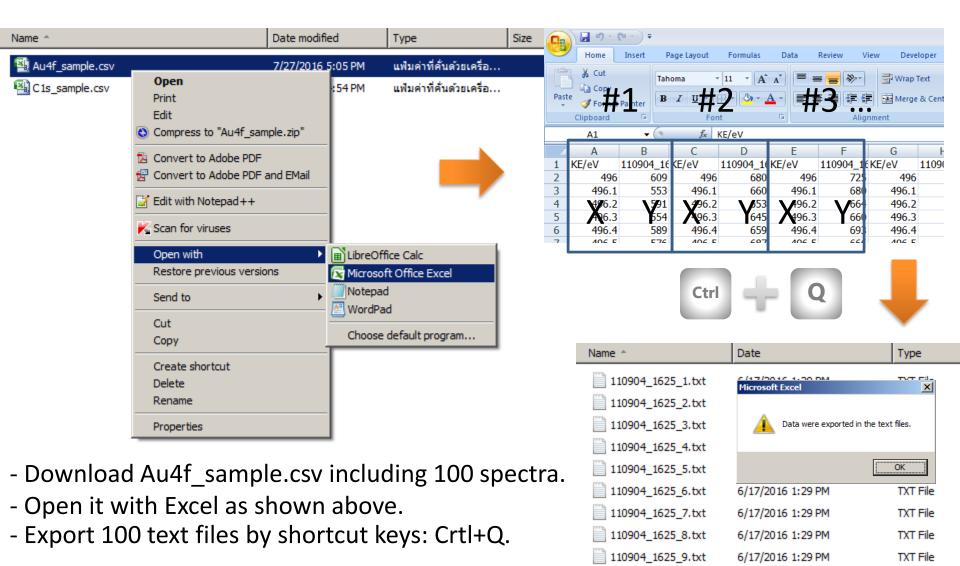


XPS: Carbon 1s spectrum





Sample file with multiple scans



110904_1625_10.txt

110904 1625 11.tvt

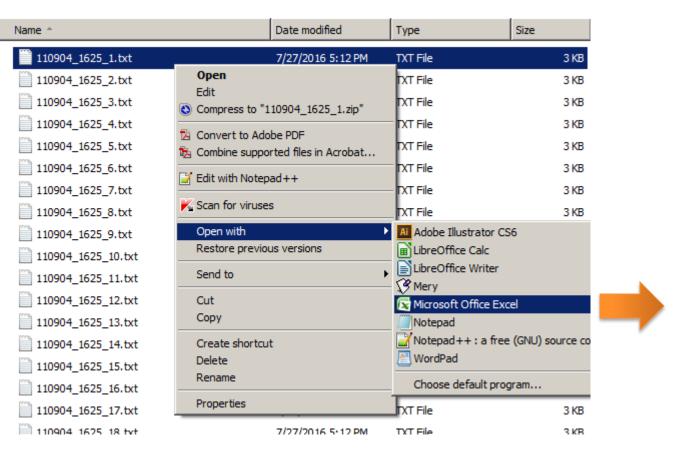
6/17/2016 1:29 PM

6/17/2016 1:29 PM

TXT File

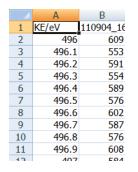
TXT File

Sample file includes single data



	Δ	В
1	KE/eV	110904_16
2	496	609
3	496.1	553
4	496.2	591
5	496.3	554
6	496.4	589
7	496.5	576
8	496.6	602
9	496.7	587
10	496.8	576
11	496.9	608
10	407	EOA

Single data to plot in Graph sheet

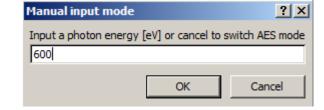


KE/eV represents a kinetic energy scale.



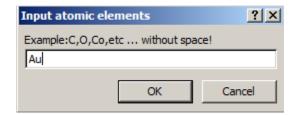
600 eV of photon energy





Sample: Gold metal foil





Note that Au4f sensitivity is not in database, so you have to input factors in database in a way below.

Au 4f BE and ASF in XPS worksheet on UD.xlsx



	Α	В	С	D
1	Element	Orbit	BE(eV)	ASF
2	C	1s	284.6	1
3	0	1s	532	2.93
4	Au	4f5/2	87.6	7.54
5	Au	4f7/2	84	9.58

Graph sheet

Adjustable factors

X axis

- **PE**: photon energy (eV)

- **WF**: work function (4)

- Char: Charging shift (0)

Y axis

- Offset: constant

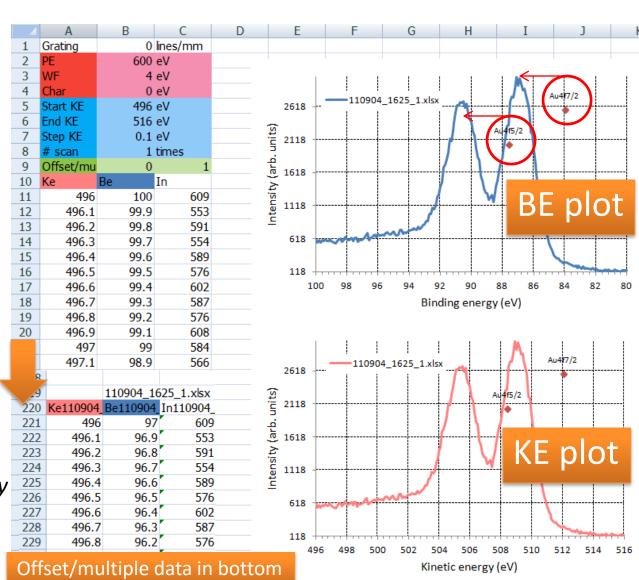
base line (0: default)

- Multiple: multiple

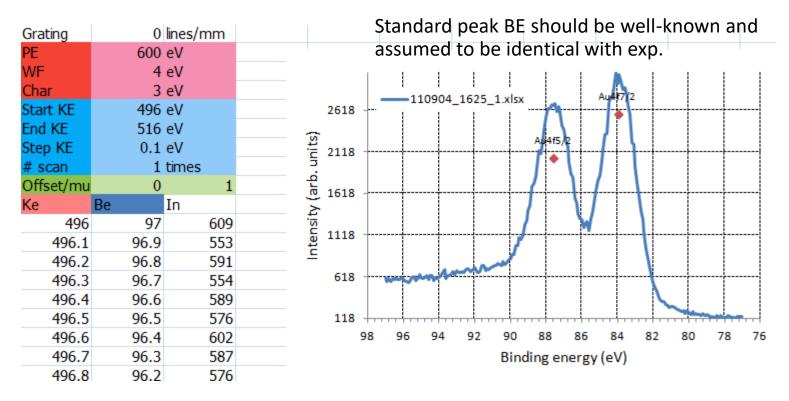
normalized factor (1)

Change **Char** to 3, and update by





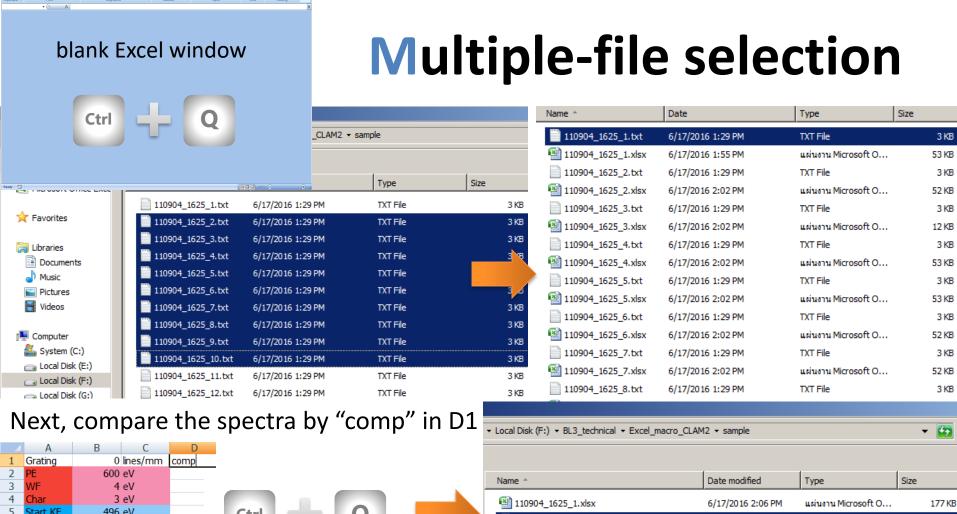
Reference spectrum



Next, apply these factors to another by "debug" in A1 cell then



	Α	В	С
1	debug	0	lines/mm
2	PE	600	eV
3	WF	4	eV
4	Char	3	eV
5	Start KE	496	eV
6	End KE	516	eV
7	Step KE	0.1	eV
8	# scan	1	times
9	Offset/mu	0	1
10	Ke	Be	In



496 eV Start KE End KE 516 eV Step KE 0.1 eV 1 times Offset/mu 0 10 In 496 97 609 11 496.1 96.9 553 12 13 496.2 96.8 591 96.7 554 496.3

14

15

16

17

18

496.4

496.5

496.6

496.7

96.6

96.5

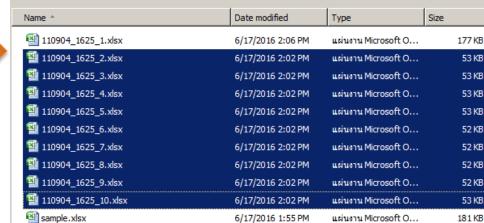
96.4

96.3

589

576

602 587



"comp" in D4 cell Ctrl Q

Fit_110904_1625_1

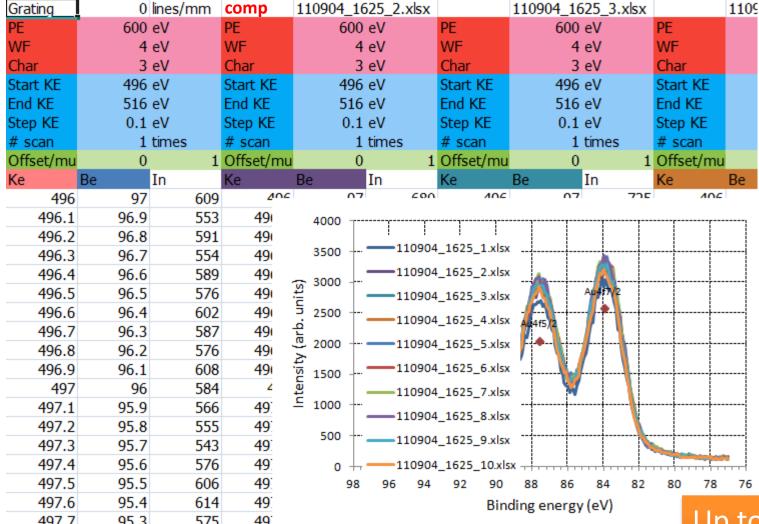
•

Ready



Graph_110904_1625_1

Compared spectra

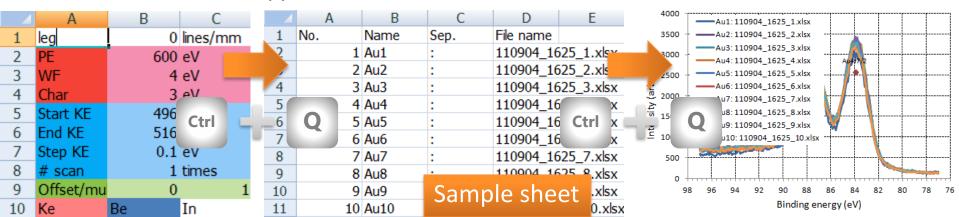


110904_1625_1

Up to 100 spectra for BE&KE plots

Analysis on spectra compared

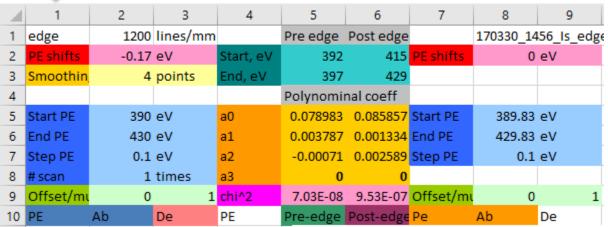
- Automatically scaled spectra for comparison
 - "auto" in A1 cell to scale at the both ends
 - "auto[x1:x2,x3:x4]" in A1 cell to scale in the ranges of (x1, x2) for offset (0) and (x3, x4) for multiple (1).
 - "auto{284.6}" in A1 cell to shift the maximum intensity to 284.6 eV in BE scale.
 - "auto'-7.8" in A1 cell to set all char factor to be -7.8 eV.
- Annotate legends in each plot
 - "leg" in A1 cell to generate Sample sheet
 - Describe sample or spectrum name in the sheet
 - "leg" in A1 cell to run the code again.
- Normalize spectra with ref. spectrum by "norm" in A1 cell.
 - Second set of data will be selected for ref. and third set to be normalized data.
 - Norm sheet appears from the data normalized.







Edge correction in NEXAFS



Post-edge region to be 1

New spline edge correction

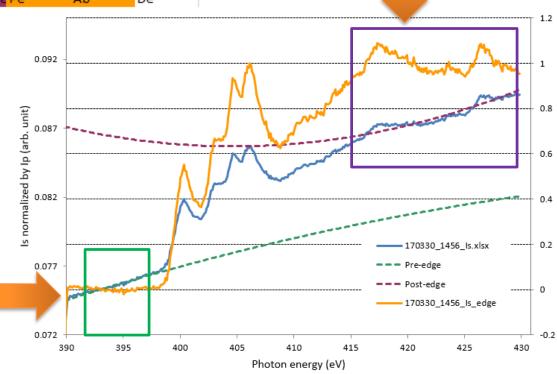
A1 cells: "edge"

New *Linear combination*

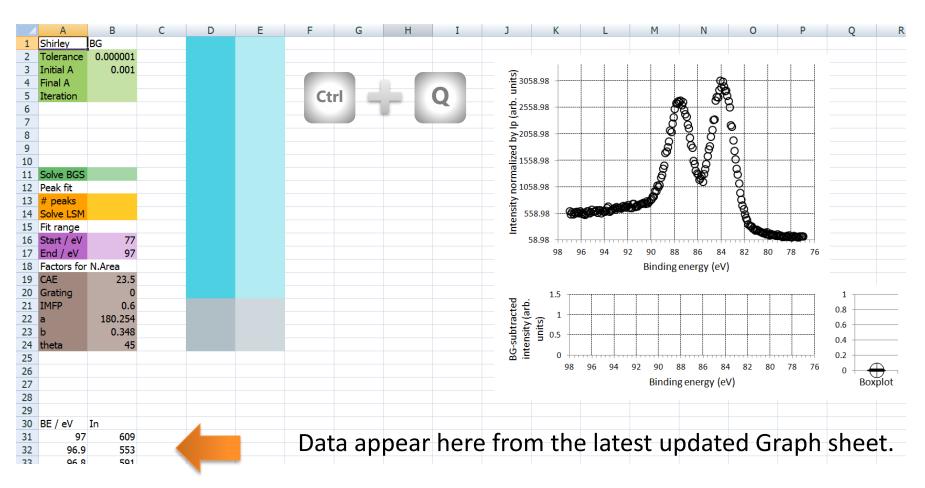
A1 cells: "lcmb"

After two references added.

Pre-edge regions to be 0



Fit sheet



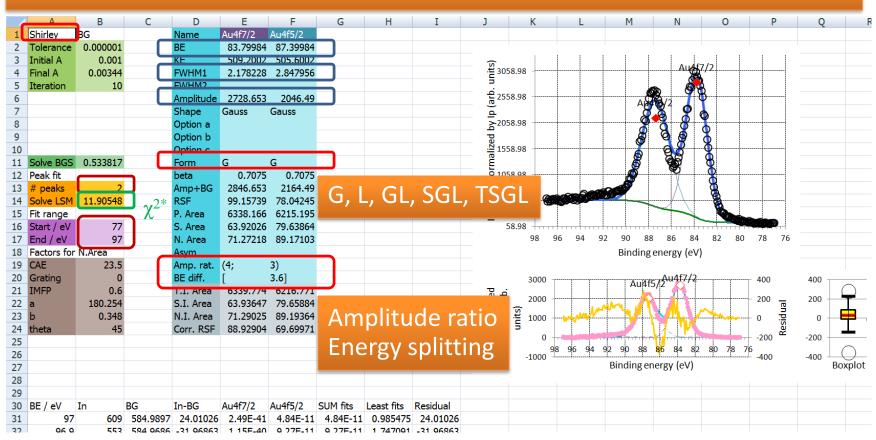


Switch the worksheet

Note that offset/multiple factors should be aligned for all spectra in Graph sheet to evaluate the peak area in Fit sheet.

Fitting functions and parameters

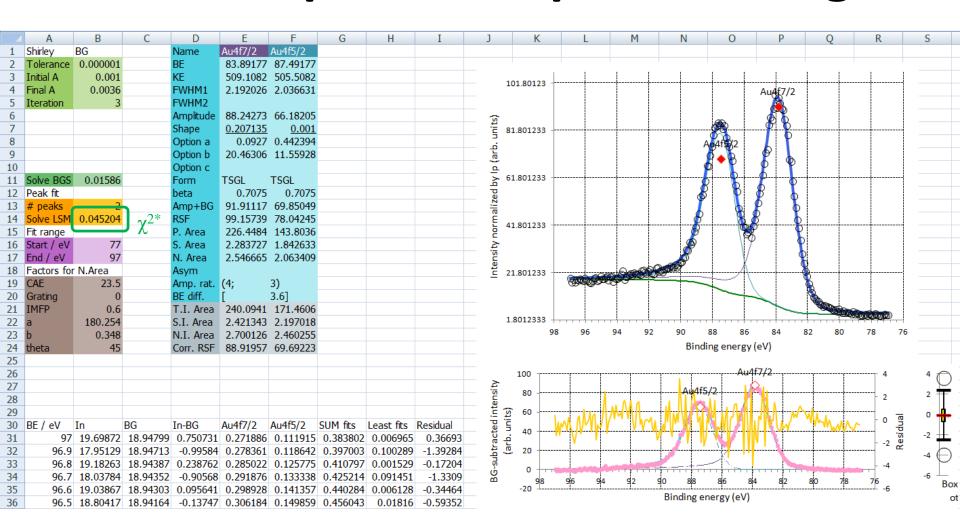
Shirley, Tougaard, Polynomial, Poly Shirley, Poly Tougaard, Victoreen, Arctan



Update results by setup constraints with bold font style, various BGs or Forms

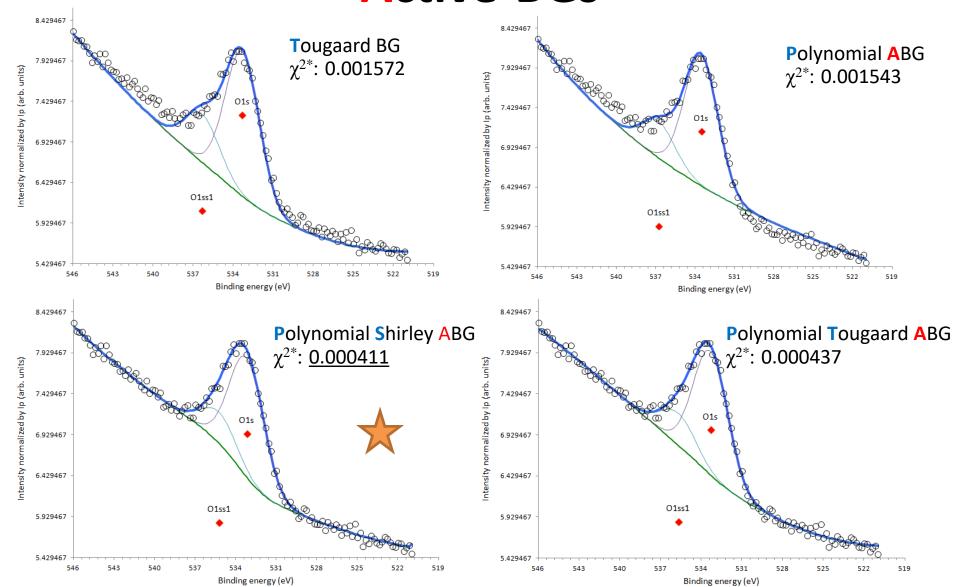


Peak shape: Tailed pseudo-Voigt

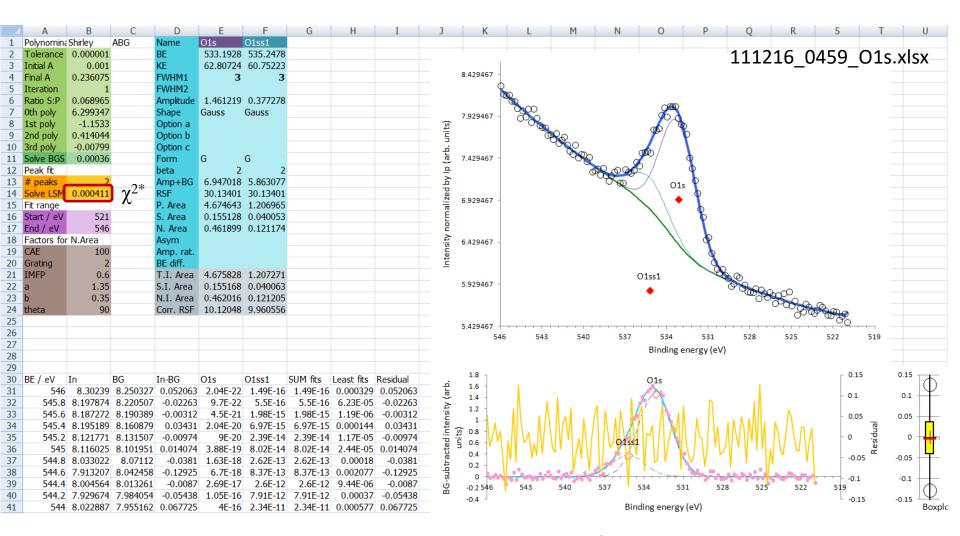


Shirley's BG, Tailed Voigt, Doublet 4f (4:3), ΔSO Au4f: 3.6 eV are used for above fitting. User-defined peak and BG functions can be easily implemented in your Visual Basic code.

Polynomial (Shirley or Tougaard) Active-BGs



Poly Shirley ABG



3-eV FWHM Gaussian fit

Lists of fitting functions

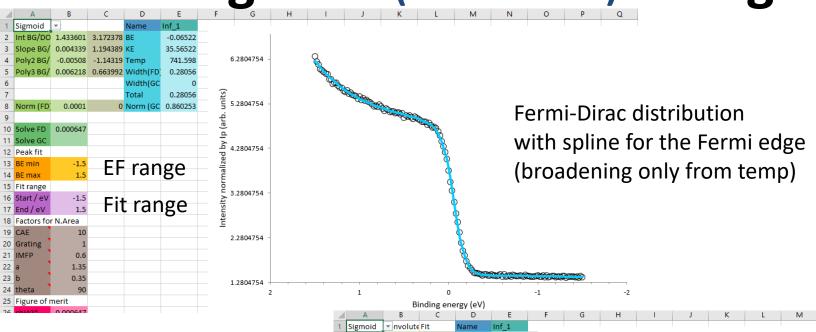
Peak shapes (total 5 forms)

Code (form)	Shape	Option a	Option b
G (0)	Gauss		
L (1)	Lorentz		
GL (0 < shape < 1)	G + L with the same FWHM1		
SGL (0 < shape < 1)	G (FWHM1) + L (FWHM2)		
<u>TSGL</u>	Exponential blend GL (FWHM1)	Tail scale	Tail length at half max

Backgrounds (total 9 BGs)

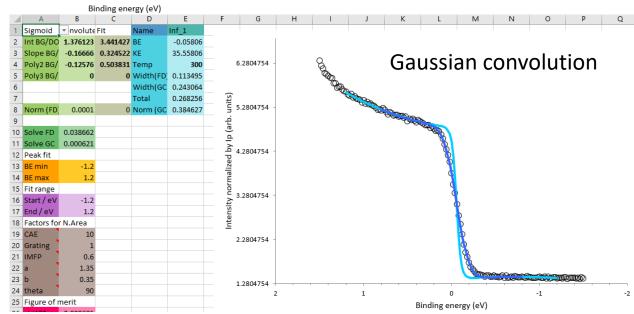
BG (A1)	BG (B1)	BG (C1)	Technique	Optimization
Shirley	BG/ABG		XPS	Static/Active
Tougaard	BG		XPS	Static
Polynomial	BG/ABG		XPS/XAS	Static/Active
Polynomial	Shirley	ABG	XPS	Active
Polynomial	Tougaard	ABG	XPS	Active
Victoreen	BG		XAS	Static
ArcTangent	BG		XAS	Active

Sigmoid (convoluted) fitting



Constraint temperature at 300 K, and polynomial parameters as well prior to convolution. Sample temp. leads to ΔE =113 meV, and instrumental resolution (BL & analyzer) to 243 meV. Total resolution becomes ΔE =268 meV.

Secondary electron cutoff can be analyzed in the same way.



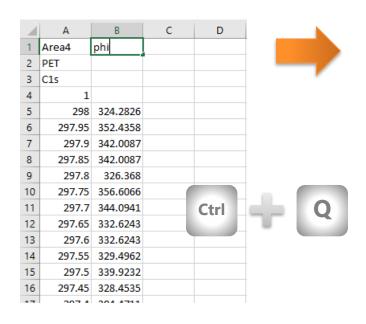
ULVAC-PHI

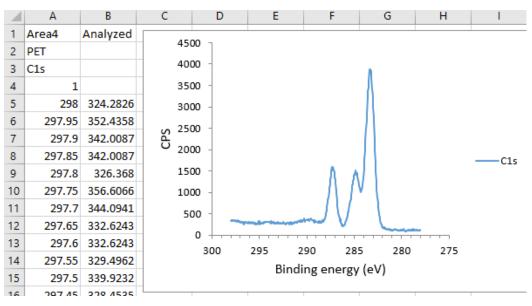
Multipak csv data analysis

Open Multipak-exported csv file with Excel, and type "phi" on B1 cell.

PET.csv	2/19/2016 2:40 PM	Microsoft Excel C	8 KB
PET.SPE	1/7/2015 4:45 PM	SPE File	7 KB

Run the macro, then new workbook appears with chart of spectrum in the csv file.





Open text file with Excel.

PET.csv	2/19/2016 2:40 PM	Microsoft Excel C	8 KB
PET.SPE	1/7/2015 4:45 PM	SPE File	7 KB
PET.xlsx	5/8/2019 2:28 PM	Microsoft Excel W	25 KB
PET_C1s.txt	5/8/2019 2:28 PM	TXT File	7 KB

Open Excel macro-exported txt file with Excel.

PET.csv	2/19/2016 2:40 PM	Microsoft Excel C	8 KB
PET.SPE	1/7/2015 4:45 PM	SPE File	7 KB
PET.xlsx	5/8/2019 2:28 PM	Microsoft Excel W	25 KB
PET_C1s.txt	5/8/2019 2:28 PM	TXT File	7 KB

5/8/2019 2:28 PM

5/8/2019 2:28 PM

5/8/2019 2:36 PM

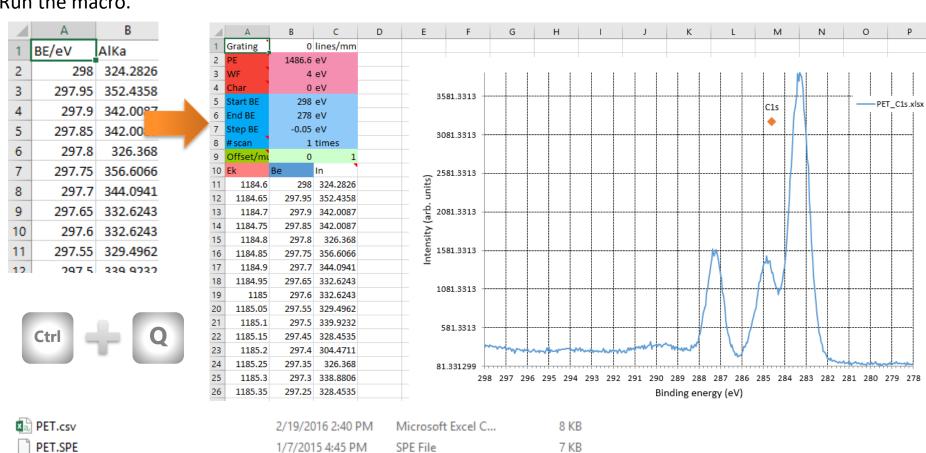
continued

Run the macro.

PET.xlsx

PET_C1s.txt

PET_C1s.xlsx



Microsoft Excel W...

Microsoft Excel W...

TXT File

25 KB

7 KB

91 KB

Multiple data-fit

	Α	В	С	D	Е	F
1	Shirley	BG		debug	Au4f7/2	Au4f5/2
2	Tolerance	0.000001		BE	83.89131	87.49132
3	Initial A	0.001		KE	509.1087	505.5087
4	Final A	0.00344		FWHM1	2.196802	2.052602
5	Iteration	6		FWHM2		

"debug"/"debuga" in D1 cell

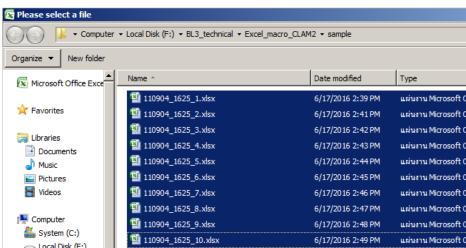
- debug: apply the original fit parameters
- debuga: apply those used just before



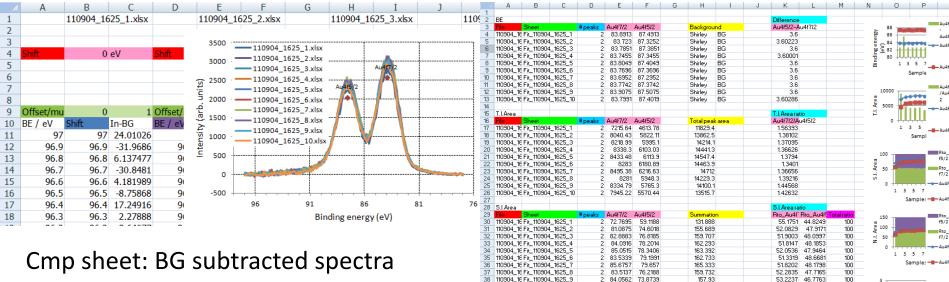
Once fitting process above done, "ana" in D1 cell to compare those fitted xlsx files.



	Α	В	С	D	E	F
1	Shirley	BG		ana	Au4f7/2	Au4f5/2
2	Tolerance	0.000001		BE	83.89131	87.49132
3	Initial A	0.001		KE	509.1087	505.5087
4	Final A	0.00344		FWHM1	2.196802	2.052602
5	Iteration	6		FWHM2		



Summary of fitting results



110904 16 Fir 110904 1625 9

110904_16 Fit_110904_1625_1

110904_16 Fit_110904_1625_2

110904 16 Fir 110904 1625 6

110904_16 Fit_110904_1625_1

110904 16 Fit 110904 1625 8

110904 16 Fit 110904 1625 9

110904_16 Fit_110904_1625_1

Ready

Ana sheet: BE, Peak areas, FWHM Rto sheet can be generated from this sheet if all elemental peak fitting results prepared for atomic ratio analysis.



T.I. area: Trapezoidal integration of peak area

173.93

178,432

181.314

182 553

181 814

184,699

178,456

176,463

53 2237

51.7958 48.2042

51.7101 48.2899

51.9491 48.0509

512273 48 7727

51,7157 48.2843

53.1194

55 0715 44 9285

51.9783 48.0217

47.821

100

100

100

100

S.I. area: T.I. area normalized with RSF

73 8739

86.0118

87.5562

87.7182

88 6757

85,3392

Au4f7/2

81 1479 66 202

90.4059 83.5241

92,4203

93.7576

94.8344

93 1385

95,5183 89,1807

93.7358

93.1164

N.I. area: T.I. area normalized with corr. RSF





Depth-profile XPS: RSC Adv. 6, 94905 (2016).

