

# Data analysis: Ctrl+Q

- efficient and effective analysis ever experienced -

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## Preface

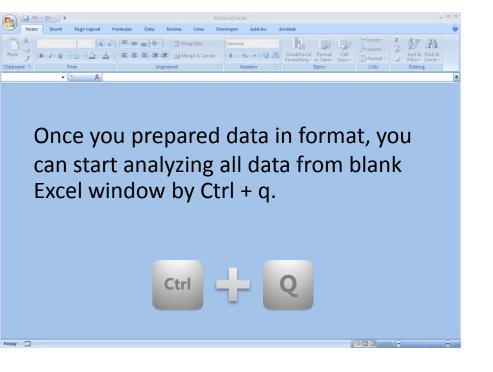
- Ctrl+Q is the first comprehensive data-analysis code for SR-XPS and XAS working on Windows Office Excel 2007 or later.
- Ctrl+Q will analyze the data by just pressing the shortcut key: Ctrl + q.
- It works great for the number of publications.
  - Carbon, JPAP, ASS, ACS AMI, SAB etc. at BL3.2Ua in 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 E	10 22572

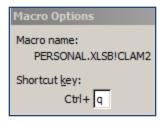
- Two-column data in the Tab-delimited format texts or spreadsheets are analyzed in the code.
- Sheet is named after the workbook name.
- Syntax in A1 cell has to be chosen as follows.



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

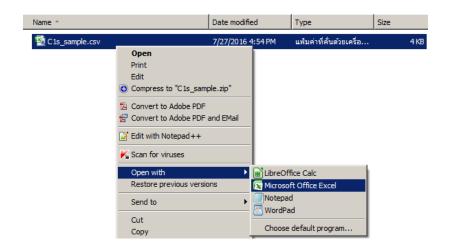
XPS sheet

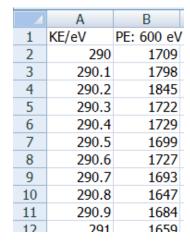
<b>A</b>	Α	В	C	D
1	Element	Orbit	BE(eV)	ASF
2	C	1s	284.6	1
3	0	1s	532	2.93

**AFS** sheet

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

# C1s sample csv







600 eV of photon energy

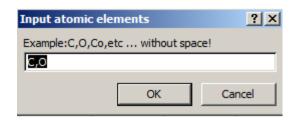


Input a photon energy [eV] or cancel to switch AES mode

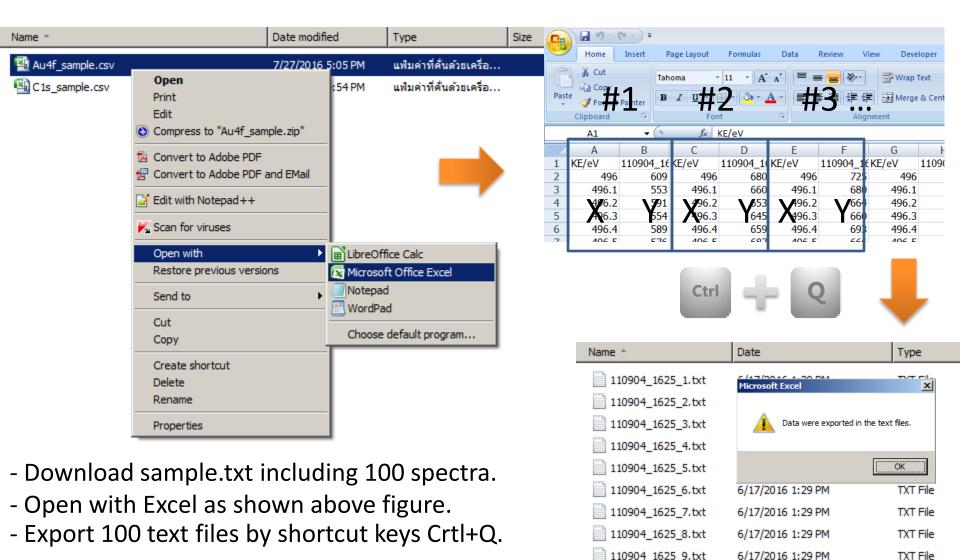
OK Cancel

XPS: Carbon 1s spectrum





# Sample file with multiple scans



110904\_1625\_10.txt

110904 1625 11.tvt

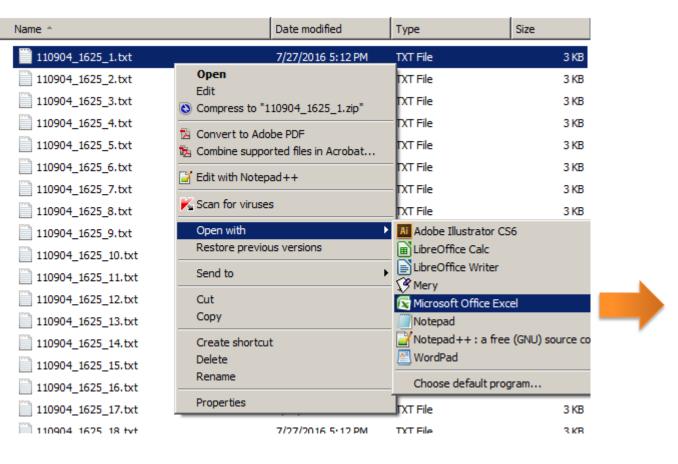
6/17/2016 1:29 PM

6/17/2016 1:29 PM

TXT File

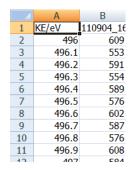
TXT File

# Sample file with single scan



	Α	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

# Data to plot in Graph sheet

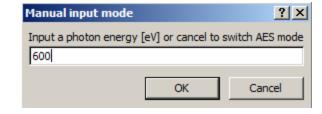


KE/eV represents a kinetic energy.



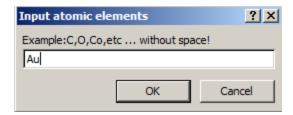
600 eV of photon energy





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 sheet on UD.xlsx



	Α	В	C	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

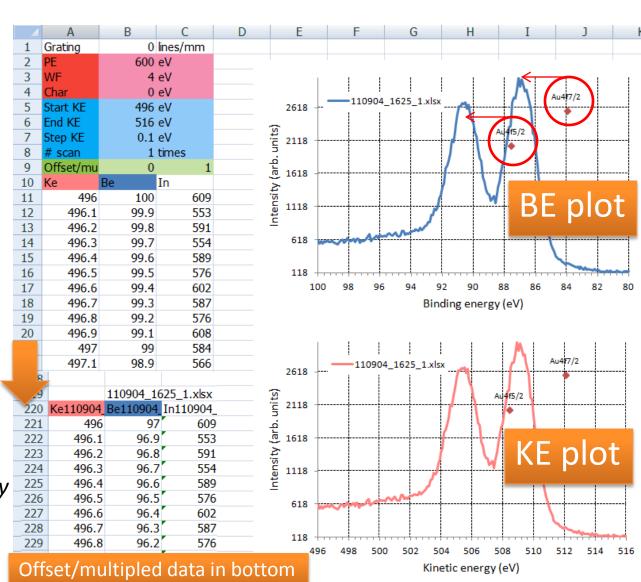
base line (0: default)

- Multiple: multiple

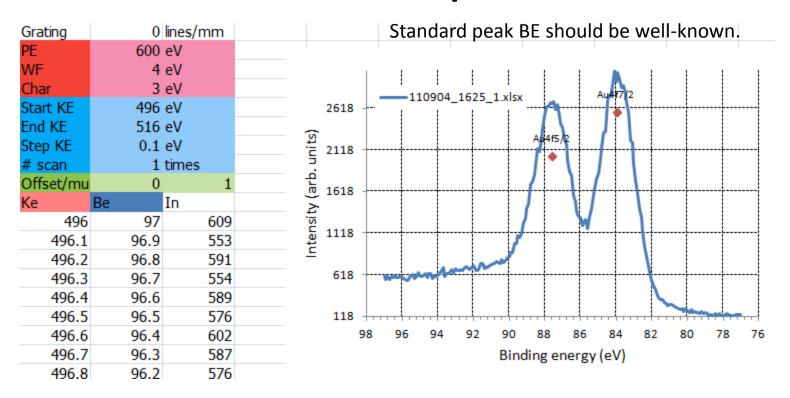
normalized factor (1)

Change Char to 3, and update by





# Reference spectrum

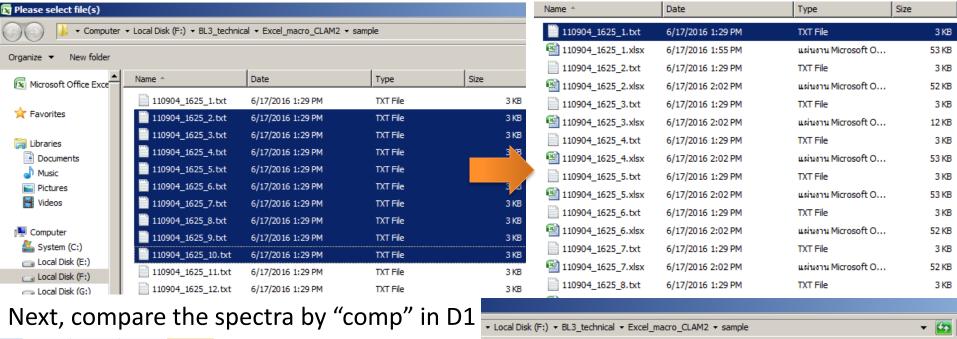


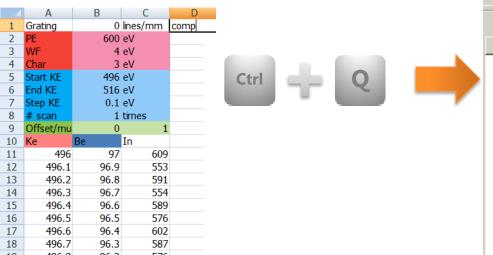
Next, apply these factors to another by "debug" in A1 cell then



	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

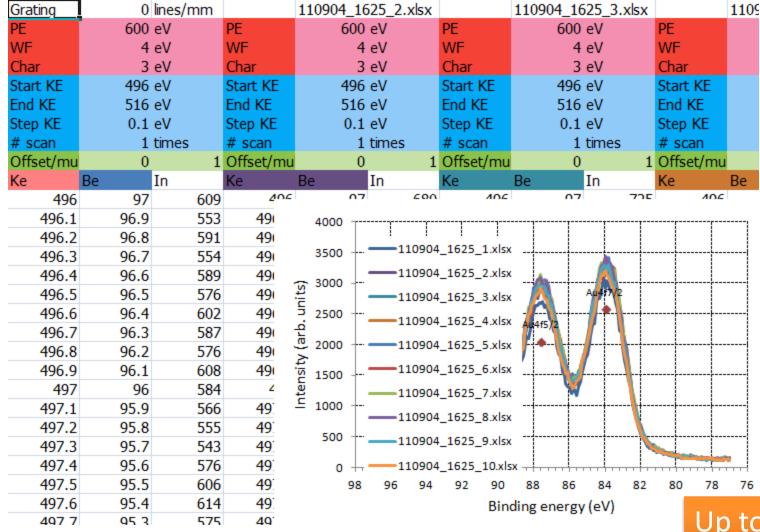
# Multiple-file selection





Name ^	Date modified	Туре	Size
110904_1625_1.xlsx	6/17/2016 2:06 PM	แผ่นงาน Microsoft O	177 KB
110904_1625_2.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	53 KB
110904_1625_3.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	53 KB
110904_1625_4.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	53 KB
110904_1625_5.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	53 KB
110904_1625_6.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	52 KB
110904_1625_7.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	52 KB
110904_1625_8.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	52 KB
110904_1625_9.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	52 KB
110904_1625_10.xlsx	6/17/2016 2:02 PM	แผ่นงาน Microsoft O	53 KB
sample.xlsx	6/17/2016 1:55 PM	แผ่นงาน Microsoft O	181 KB

# Compared spectra



110904 1625 1

Graph 110904 1625 1

Up to 100 spectra for BE&KE plots



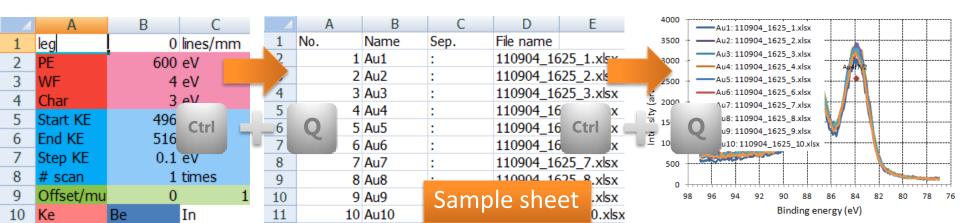
Fit 110904 1625 1

# Spectral comparison functions

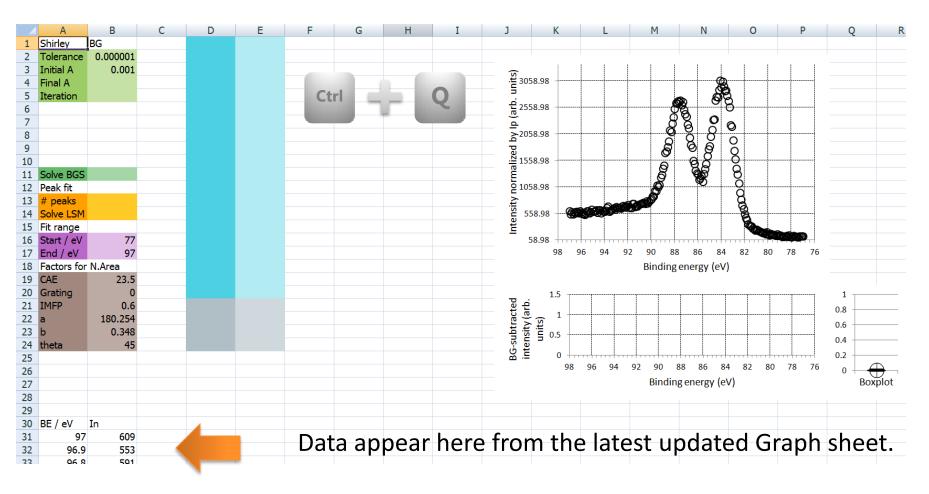
- Automatically scaled spectra for comparison
  - "auto" in A1 cell to scale at the both ends
  - "auto[x1:x2,x3:x4]" in A1 cell to scale in the ranges of (x1, x2) for offset (0) and (x3, x4) for multiple (1).

Every single step needs

- 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.



## Fit sheet

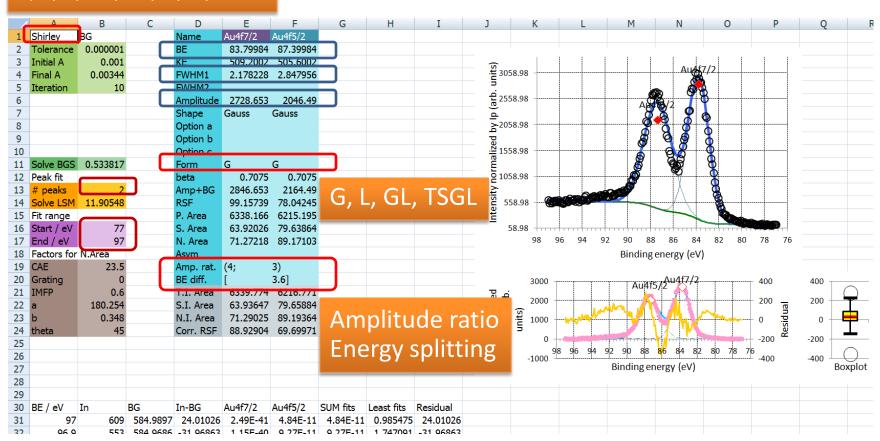




Switch the sheet

# Fitting parameters

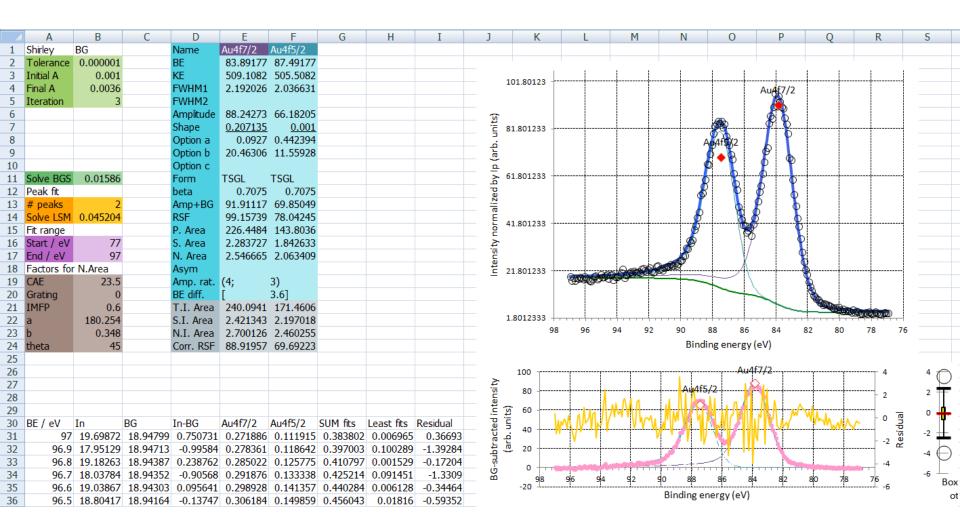
S, T, P, PS, PT, V, A, etc.



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

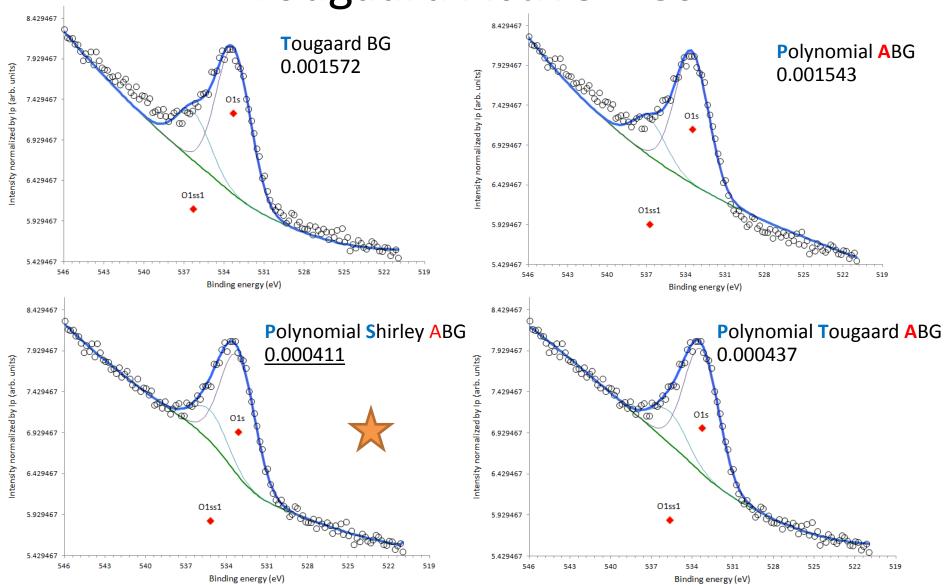


# Fitting result: tailed pseudo-Voigt

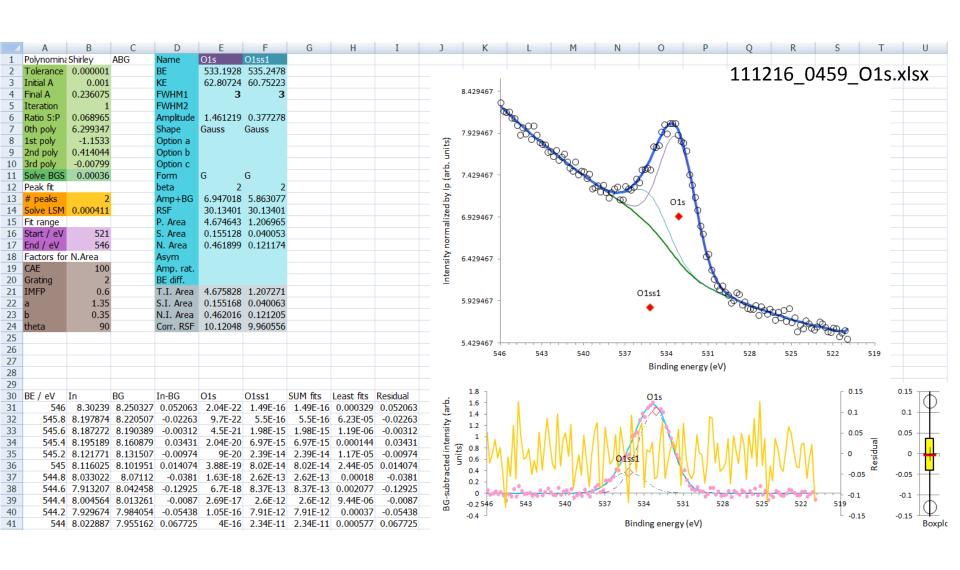


Shirley's BG, Tailed Voigt, Doublet 4f (4:3), ΔSO Au4f: 3.6 eV are used for above fitting. User-defined peak and BG functions can be easily implemented in your Visual Basic code.

# Fitting result: Polynomial Shirley & Tougaard Active-BGs



# Fitting result: Poly Shirley ABG



# Peak shapes (total 4 forms)

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

# **Backgrounds (total 8 BGs)**

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

# 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

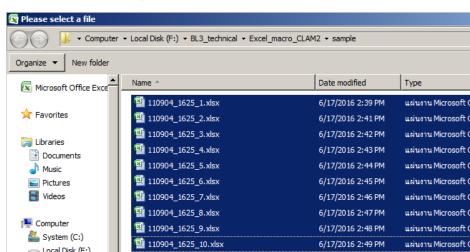
- debug: apply the original fit parameters
- debuga: apply those used just before



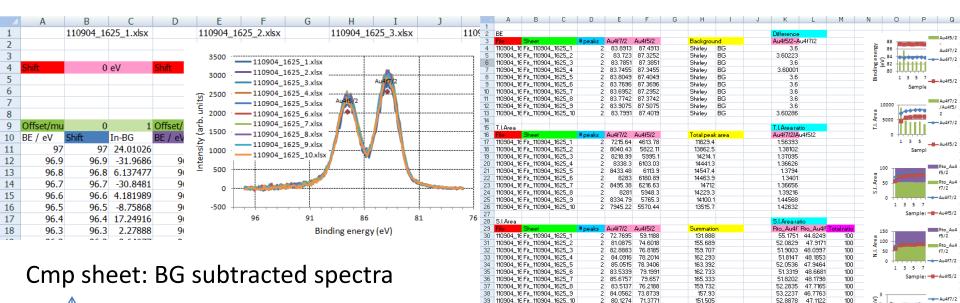
Once fitting process above done, "ana" in D1 cell to compare those fitted xlsx files.



	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



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

55 0715 44 9285

51.9783 48.0217

47.821

51.7958 48.2042

51.7101 48.2899

51.9491 48.0509

51 2273 48 7727

51,7157 48,2843

53.1194

100

100

100

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

Au4f7/2

81 1479 66 2021

90,4059 83,5241

92,4203

93.7576

94.8344

93 1385

95,5183 89,1807

93.7358

93.1164

86.0118

87.5562

87.7182

88 6757

85,3392

110904\_16 Fit\_110904\_1625\_1

110904\_16 Fit\_110904\_1625\_2

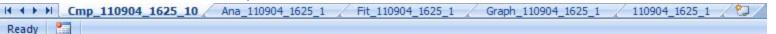
110904 16 Fir 110904 1625 6

110904\_16 Fit\_110904\_1625\_1

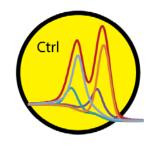
110904 16 Fit 110904 1625 8

110904\_16 Fit\_110904\_1625\_1

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



# Enjoy!



## Depth-profile XPS: under review in RSC Adv.

