

CloudCal

Lee Drake

August 2, 2017

Paleoresearch Institute

CloudCal Spectrum Counts Add Concentrations Cal Curves Apply Calibration

X-Ray Fluorescence Calibration

Calibration Name

Process Data Plot Spectrum Plot

Choose Spectra
 No file selected

Spectra
 Net

Element:

Load Cal File
 No file selected

Use Cal File

X-Ray Fluorescence Calibration

Calibration Name

Obsidian

Give your
calibration a name

Process Data

Plot Spectrum

Plot

Choose Spectra

Browse...

No file selected

When you are done there,
you can load spectra

Spectra

Net

Element:

(Fe) Iron

Load Cal File

Browse...

No file selected

Use Cal File

CloudCal 127.0.0.1:3315 Lee

CloudCal Spectrum Counts Add Concentrations Cal Curves Apply Cal

X-Ray Fluorescence Calibration

Calibration Name: Obsidian

Process Data Plot Spectrum Plot

Choose Spectra: No file selected

Spectra Net

Element: (Fe) Iron

Load Cal File: No file selected

Use Cal File

Select your spectra in the pop-up window

OB40Archibarca35.CSV
OB40Archibarca35.pdz
OB40Basalti...lateau20.CSV
OB40Basaltic_Plateau20.pdz
OB40Big_So...Butte06.CSV
OB40Big_So...Butte06.pdz
OB40Blue_Mountain04.CSV
OB40Blue_Mountain04.pdz
OB40Burns_Green15.CSV
OB40Burns_Green15.pdz
OB40Cannonball1_22.CSV
OB40Cannonball1_22.pdz
OB40Casa_Diablo10.CSV

Format: Custom Files

Options Cancel Open

X-Ray Fluorescence Calibration

Calibration Name
Obsidian

Process Data Plot Spectrum 

Choose Spectra
Browse... 40 files Upload complete

Spectra Net

Element:
(Fe) Iron

Load Cal File
Browse... No file selected

Use Cal File

When spectra are done uploading,
you will be notified here

X-Ray Fluorescence Calibration

Calibration Name
Obsidian

Process Data

Choose Spectra
Browse... 40 files
Upload complete

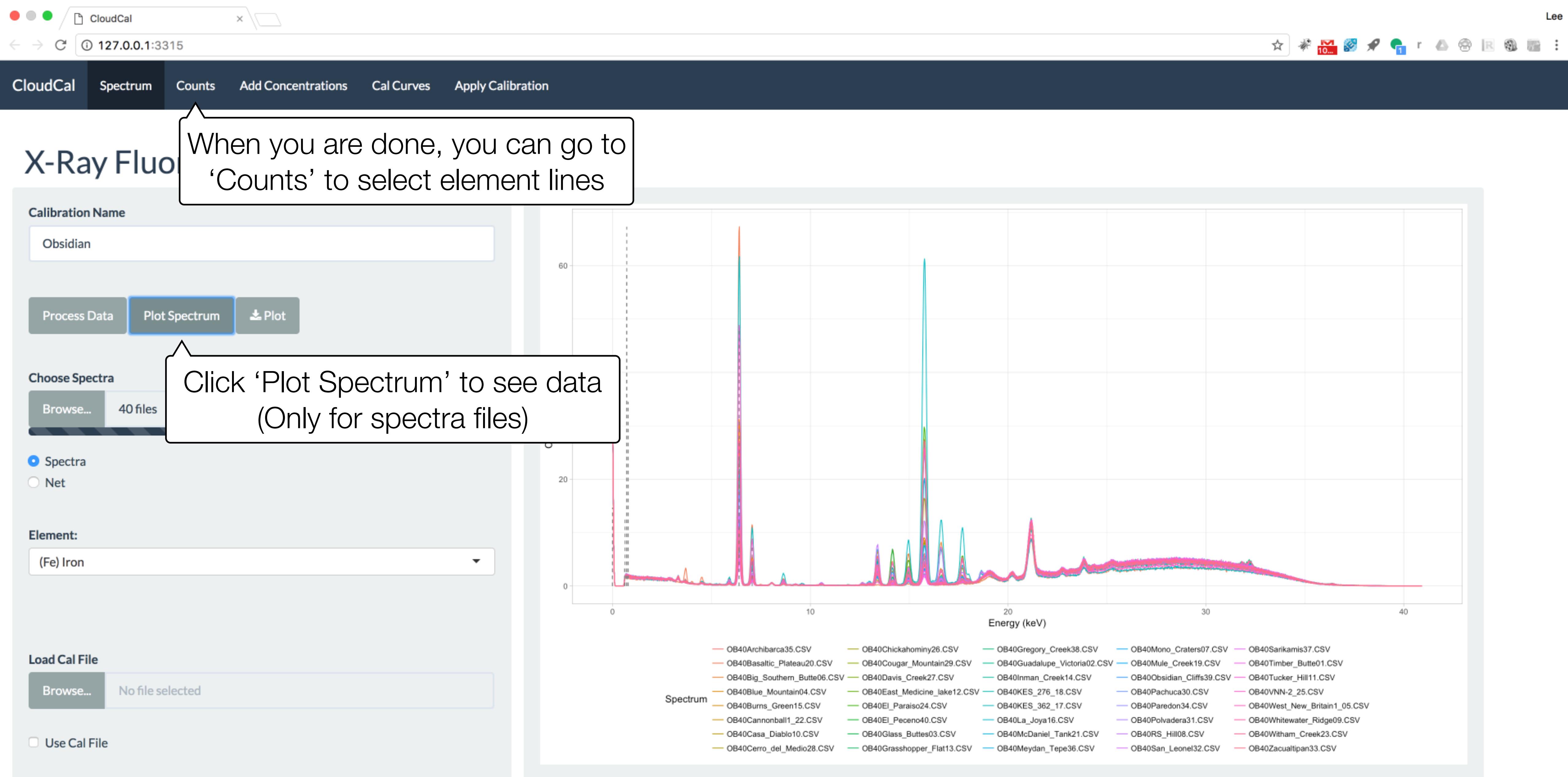
Spectra
 Net

Element:
(Fe) Iron

Load Cal File
Browse... No file selected

Use Cal File

Next, click 'Process Data'



[Confirm Elements](#)
 [Table](#)
Elemental lines to show:

- Ne.K.alpha
- Ne.K.beta
- Na.K.alpha
- Na.K.beta
- Mg.K.alpha
- Mg.K.beta
- Al.K.alpha
- Al.K.beta
- Si.K.alpha
- Si.K.beta
- P.K.alpha
- P.K.beta
- S.K.alpha
- S.K.beta
- Cl.K.alpha
- Cl.K.beta
- Ar.K.alpha
- Ar.K.beta
- K.K.alpha
- K.K.beta
- Ca.K.alpha
- Ca.K.beta
- Sc.K.alpha
- Sc.K.beta
- Ti.K.alpha
- Ti.K.beta
- V.K.alpha
- V.K.beta
- Cr.K.alpha
- Cr.K.beta
- Mn.K.alpha

Choose your spectra in the panel to the left

Spectral Lines

Show 10 entries

	Ca.K.alpha	Ti.K.alpha	Fe.K.alpha	Cu.K.alpha	Zn.K.alpha	Pb.L.alpha
1	3	1	27	2	1	2
2	6	3	194	2	2	1
3	2	1	38	2	4	3
4	1	1	90	2	3	1
5	1	1	57	2	2	2
6	1	1	75	2	3	2
7	2	1	29	2	1	2
8	1	1	23	2	1	2
9	2	1	37	2	1	2
10	2	1	26	2	2	1

Showing 1 to 10 of 40 entries

Previous 1 2 3 4 Next

Counts per second for element lines will appear here

127.0.0.1:3315



- Cr.K.alpha
- Cr.K.beta
- Mn.K.alpha
- Mn.K.beta
- Fe.K.alpha
- Fe.K.beta
- Co.K.alpha
- Co.K.beta
- Ni.K.alpha
- Ni.K.beta
- Cu.K.alpha
- Cu.K.beta
- Zn.K.alpha
- Zn.K.beta
- Ga.K.alpha
- Ga.K.beta
- Ge.K.alpha
- Ge.K.beta
- As.K.alpha
- As.K.beta
- Se.K.alpha
- Se.K.beta
- Br.K.alpha
- Br.K.beta
- Kr.K.alpha
- Kr.K.beta
- Rb.K.alpha
- Rb.K.beta
- Sr.K.alpha
- Sr.K.beta
- Y.K.beta
- Y.K.alpha
- Zr.K.alpha
- Zr.K.beta
- Nb.K.alpha
- Nb.K.beta
- Mo.K.alpha
- Mo.K.beta
- Mo.L.alpha

Here, you don't need to have concentrations for all elements - you can check some if you only want to use them for corrections (e.g. correct As K-alpha by Pb L-beta)

You will need to scroll down - all possible lines are listed here

[Confirm Elements](#) [Table](#)

When you are done, click
'Confirm Elements'

- Ele
- Na.K.beta
- Mg.K.alpha
- Mg.K.beta
- Al.K.alpha
- Al.K.beta
- Si.K.alpha
- Si.K.beta
- P.K.alpha
- P.K.beta
- S.K.alpha
- S.K.beta
- Cl.K.alpha
- Cl.K.beta
- Ar.K.alpha
- Ar.K.beta
- K.K.alpha
- K.K.beta
- Ca.K.alpha
- Ca.K.beta
- Sc.K.alpha
- Sc.K.beta
- Ti.K.alpha
- Ti.K.beta
- V.K.alpha
- V.K.beta
- Cr.K.alpha
- Cr.K.beta
- Mn.K.alpha

Spectral Lines

Show 10 entries

Search:

	K.K.alpha	Ca.K.alpha	Ti.K.alpha	Mn.K.alpha	Fe.K.alpha	Cu.K.alpha	Zn.K.alpha	Ga.K.alpha	As.K.alpha	Rb.K.alpha	Sr.K.alpha	Y.K.b
1	3	3	1	2	27	2	1	1	1	11	38	
2	2	6	3	3	194	2	2	1	0	1	25	
3	4	2	1	1	38	2	4	1	2	26	1	
4	3	1	1	4	90	2	3	1	1	5	1	
5	3	1	1	2	57	2	2	1	1	10	1	
6	3	1	1	1	75	2	3	1	2	31	1	
7	4	2	1	1	29	2	1	1	1	14	13	
8	3	1	1	1	23	2	1	1	1	15	1	
9	3	2	1	1	37	2	1	1	2	10	3	
10	3	2	1	1	26	2	2	1	1	9	4	

Showing 1 to 10 of 40 entries

Previous [1](#) [2](#) [3](#) [4](#) Next

Enter Values

Navigate to 'Add Concentrations'

Concentration data
will begin empty,
you will need to
provide these values

Enter Concentrations

Enter Values

Enter Concentrations

	Spectrum	K.K.alpha	Ca.K.alpha	Ti.K.alpha	Mn.K.alpha	Fe.K.alpha	Co.K.alpha	Cu.K.alpha	Zn.K.alpha	Ga.K.alpha	As.K.alpha	Rb.K.alpha	Sr.K.alpha	Y.K.
1	OB40Archibarca35.CSV	3.17	1.08	0.08	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
2	OB40Basaltic_Plateau20.CSV	0.20	4.99	0.78	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
3	OB40Big_Southern_Butte06.CSV	4.06	0.32	0.05	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
4	OB40Blue_Mountain04.CSV	2.77	0.10	0.12	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
5	OB40Burns_Green15.CSV	3.50	0.13	0.10										
6	OB40Cannonball1_22.CSV	3.68	0.18	0.10										
7	OB40Casa_Diablo10.CSV	3.98	0.61	0.11										
8	OB40Cerro_del_Medio28.CSV	3.64	0.25	0.05										
9	OB40Chickahominy26.CSV	3.52	0.42	0.11										
10	OB40Cougar_Mountain29.CSV	3.22	0.47	0.03										
11	OB40Davis_Creek27.CSV	4.01	0.58	0.05										
12	OB40East_Medicine_lake12.CSV	3.66	0.63	0.14										
13	OB40El_Paraiso24.CSV	3.76	0.11	0.08	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
14	OB40El_Peceno40.CSV	3.85	0.81	0.05	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
15	OB40Glass_Buttes03.CSV	3.60	0.52	0.06	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
16	OB40Grasshopper_Flat13.CSV	3.76	0.58	0.13	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
17	OB40Gregory_Creek38.CSV	3.60	0.82	0.02	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
18	OB40Guadalupe_Victoria02.CSV	3.40	0.34	0.06	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
19	OB40Inman_Creek14.CSV	2.46	0.71	0.05	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
20	OB40KES_276_18.CSV	4.14	0.57	0.31	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
21	OB40KES_362_17.CSV	3.31	0.15	0.16	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
22	OB40La_Joya16.CSV	3.48	0.12	0.09	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
23	OB40McDaniel_Tank21.CSV	3.90	0.69	0.16	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
24	OB40Meydan_Tepe36.CSV	3.46	0.29	0.05	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
25	OB40Mono_Craters07.CSV	3.53	0.38	0.04	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
26	OB40Mule_Creek19.CSV	3.57	0.39	0.04	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
27	OB40Obsidian_Cliffs39.CSV	2.83	0.59	0.06	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
28	OB40Pachuca30.CSV	3.29	0.08	0.11	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
29	OB40Paredon34.CSV	3.91	0.26	0.08	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
30	OB40Polvadera31.CSV	3.98	0.31	0.04	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
31	OB40RS_Hill08.CSV	3.35	0.25	0.02	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
32	OB40San_Leonel32.CSV	3.56	0.17	0.06	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
33	OB40Sarikamis37.CSV	3.92	0.32	0.05	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
34	OB40Timber_Butte01.CSV	3.65	0.49	0.03	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
35	OB40Tucker_Hill11.CSV	3.50	0.54	0.03	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
36	OB40VNN-2_25.CSV	3.02	0.10	0.07	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
37	OB40West_New_Britain1_05.CSV	1.71	0.86	0.11	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00

You can simply paste values from another spreadsheet here

CloudCal Spectrum Counts Add Concentrations

Cal Curves Apply Calibration

Enter Values

Enter Concentrations

	pha	Ti.K.alpha	Mn.K.alpha	Fe.K.alpha	Co.K.alpha	Cu.K.alpha	Zn.K.alpha	Ga.K.alpha	As.K.alpha	Rb.K.alpha	Sr.K.alpha	Y.K.alpha	Zr.K.alpha	Nb.K.alpha	Pb.L.alpha	Th.L.alpha
1	1.08	0.08	0.05	0.87	0.00	0.00	0.00	0.00	0.00	0.01	0.03	0.00	0.01	0.00	0.00	0.00
2	1.99	0.78	0.11	6.85	0.00	0.00	0.01	0.00	0.00	0.00	0.03	0.00	0.01	0.00	0.00	0.00
3	0.32	0.05	0.03	1.17	0.00	0.00	0.03	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
4	0.10	0.12	0.16	2.74	0.00	0.00	0.02	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
5	0.13	0.10	0.05	1.72	0.00	0.00	0.01	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
6	0.18	0.10	0.05	2.33	0.00	0.00	0.02	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
7	0.61	0.11	0.03	0.93	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
8	0.25	0.05	0.04	0.71	0.00	0.00	0.01	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
9	0.42	0.11	0.04	1.18	0.00	0.00	0.01	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
10	0.47	0.03	0.03	0.81	0.00	0.00	0.01	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
11	0.58	0.05	0.04	0.53	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
12	0.63	0.14	0.03	1.04	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.02	0.00	0.00	0.00
13	0.11	0.08	0.02	1.94	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.02	0.00	0.11	0.01	0.00
14	0.81	0.05	0.09	0.61	0.00	0.00	0.01	0.00	0.00	0.02	0.03	0.00	0.01	0.00	0.00	0.00
15	0.52	0.06	0.03	0.62	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
16	0.58	0.13	0.03	0.92	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.02	0.00	0.00	0.00
17	0.82	0.02	0.07	0.65	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
18	0.34	0.06	0.05	0.43	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
19	0.71	0.05	0.05	1.08	0.00	0.00	0.01	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
20	0.57	0.31	0.11	2.34	0.00	0.00	0.01	0.00	0.00	0.02	0.01	0.01	0.10	0.03	0.00	0.00
21	0.15	0.16	0.18	5.37	0.00	0.00	0.06	0.00	0.00	0.04	0.00	0.04	0.31	0.06	0.00	0.01
22	0.12	0.09	0.06	1.89	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.07	0.01	0.00	0.00
23	0.69	0.16	0.06	1.00	0.00	0.00	0.01	0.00	0.00	0.02	0.02	0.00	0.02	0.00	0.00	0.00
24	0.29	0.05	0.05	0.93	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.03	0.00	0.00	0.00
25	0.38	0.04	0.04	0.78	0.00	0.00	0.00	0.00	0.00	0.02	0.00	0.01	0.00	0.00	0.00	0.00
26	0.39	0.04	0.04	0.66	0.00	0.00	0.00	0.00	0.00	0.02	0.00	0.00	0.01	0.00	0.00	0.00
27	0.59	0.06	0.03	0.73	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
28	0.08	0.11	0.11	1.63	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.01	0.09	0.01	0.00	0.00
29	0.26	0.08	0.04	0.85	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.00	0.00
30	0.31	0.04	0.04	0.38	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
31	0.25	0.02	0.04	0.72	0.00	0.00	0.01	0.00	0.00	0.04	0.00	0.01	0.02	0.02	0.00	0.00
32	0.17	0.06	0.03	1.24	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.01	0.04	0.00	0.00	0.00
33	0.32	0.05	0.04	0.55	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
34	0.49	0.03	0.08	0.37	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.00	0.01	0.00	0.00	0.00
35	0.54	0.03	0.05	0.47	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
36	0.10	0.07	0.12	4.38	0.00	0.00	0.03	0.00	0.00	0.02	0.00	0.01	0.11	0.01	0.00	0.00
37	0.86	0.11	0.06	0.86	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.00	0.00	0.00	0.00

Even though data rounds to two decimal places here, the values you paste are captured

CloudCal Spectrum Counts Add Concentrations

Cal Curves Apply Calibration

Enter Values

When the table is complete, you can click 'Enter Values'

Enter Concentrations

	pha	Ti.K.alpha	Mn.K.alpha	Fe.K.alpha	Co.K.alpha	Cu.K.alpha	Zn.K.alpha	Ga.K.alpha	As.K.alpha	Rb.K.alpha	Sr.K.alpha	Y.K.alpha	Zr.K.alpha	Nb.K.alpha	Pb.L.alpha	Th.L.alpha
1	1.08	0.08	0.05	0.87	0.00	0.00	0.00	0.00	0.00	0.01	0.03	0.00	0.01	0.00	0.00	0.00
2	1.99	0.78	0.11	6.85	0.00	0.00	0.01	0.00	0.00	0.00	0.03	0.00	0.01	0.00	0.00	0.00
3	0.32	0.05	0.03	1.17	0.00	0.00	0.03	0.00	0.00	0.03	0.00	0.02	0.03	0.03	0.00	0.00
4	0.10	0.12	0.16	2.74	0.00	0.00	0.02	0.00	0.00	0.01	0.00	0.01	0.04	0.00	0.00	0.00
5	0.13	0.10	0.05	1.72	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.01	0.06	0.00	0.00	0.00
6	0.18	0.10	0.05	2.33	0.00	0.00	0.02	0.00	0.00	0.03	0.00	0.01	0.11	0.01	0.00	0.00
7	0.61	0.11	0.03	0.93	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.02	0.00	0.00	0.00
8	0.25	0.05	0.04	0.71	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.00	0.00
9	0.42	0.11	0.04	1.18	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.01	0.03	0.00	0.00	0.00
10	0.47	0.03	0.03	0.81	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.01	0.01	0.00	0.00	0.00
11	0.58	0.05	0.04	0.53	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
12	0.63	0.14	0.03	1.04	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.02	0.00	0.00	0.00
13	0.11	0.08	0.02	1.94	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.02	0.11	0.01	0.00	0.00
14	0.81	0.05	0.09	0.61	0.00	0.00	0.01	0.00	0.00	0.02	0.03	0.00	0.01	0.00	0.00	0.00
15	0.52	0.06	0.03	0.62	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
16	0.58	0.13	0.03	0.92	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.02	0.00	0.00	0.00
17	0.82	0.02	0.07	0.65	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
18	0.34	0.06	0.05	0.43	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
19	0.71	0.05	0.05	1.08	0.00	0.00	0.01	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
20	0.57	0.31	0.11	2.34	0.00	0.00	0.01	0.00	0.00	0.02	0.01	0.01	0.10	0.03	0.00	0.00
21	0.15	0.16	0.18	5.37	0.00	0.00	0.06	0.00	0.00	0.04	0.00	0.04	0.31	0.06	0.00	0.01
22	0.12	0.09	0.06	1.89	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.07	0.01	0.00	0.00
23	0.69	0.16	0.06	1.00	0.00	0.00	0.01	0.00	0.00	0.02	0.02	0.00	0.02	0.00	0.00	0.00
24	0.29	0.05	0.05	0.93	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.03	0.00	0.00	0.00
25	0.38	0.04	0.04	0.78	0.00	0.00	0.00	0.00	0.00	0.02	0.00	0.01	0.00	0.00	0.00	0.00
26	0.39	0.04	0.04	0.66	0.00	0.00	0.00	0.00	0.00	0.02	0.00	0.00	0.01	0.00	0.00	0.00
27	0.59	0.06	0.03	0.73	0.00	0.00	0.00	0.00	0.00	0.01	0.01	0.00	0.01	0.00	0.00	0.00
28	0.08	0.11	0.11	1.63	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.01	0.09	0.01	0.00	0.00
29	0.26	0.08	0.04	0.85	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.00	0.02	0.00	0.00	0.00
30	0.31	0.04	0.04	0.38	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
31	0.25	0.02	0.04	0.72	0.00	0.00	0.01	0.00	0.00	0.04	0.00	0.01	0.02	0.02	0.00	0.00
32	0.17	0.06	0.03	1.24	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.01	0.04	0.00	0.00	0.00
33	0.32	0.05	0.04	0.55	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
34	0.49	0.03	0.08	0.37	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.00	0.01	0.00	0.00	0.00
35	0.54	0.03	0.05	0.47	0.00	0.00	0.00	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00	0.00
36	0.10	0.07	0.12	4.38	0.00	0.00	0.03	0.00	0.00	0.02	0.00	0.01	0.11	0.01	0.00	0.00
37	0.86	0.11	0.06	0.86	0.00	0.00	0.01	0.00	0.00	0.02	0.00	0.01	0.00	0.00	0.00	0.00

Plot Update Save 

Element
K.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept

K.Kalpha
 Ca.Kalpha
 Ti.Kalpha
 Mn.Kalpha
 Fe.Kalpha
 Co.Kalpha
 Cu.Kalpha
 Zn.Kalpha
 Ga.Kalpha
 As.Kalpha

View Curves Diagnostics Standards

Error: arguments imply differing number of rows: 40, 0

You may see error when you begin - this is fine unless it persists after you begin calibration

When the table is complete, you can click 'Enter Values'

Element

K.Kalpha

- As.Kalpha
- Rb.Kalpha
- Sr.Kalpha
- Y.Kalpha
- Zr.Kalpha
- Nb.Kalpha
- Pb.L.alpha
- Th I alpha

Compton

Min

18.5

Max

19.5

Intercept

K.Kalpha

Ca.Kalpha

Ti.Kalpha

Mn.Kalpha

Fe.K.alpha

Co.Kalpha

Cu.Kalpha

Zn.Kalpha

Ga.Kalpha

Al.Kalpha

View Curves Diagnostics Standards

Error: arguments imply differing number of rows: 40, 0

Select the element line you would like to calibrate

Plot **Update** **Save** 

Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

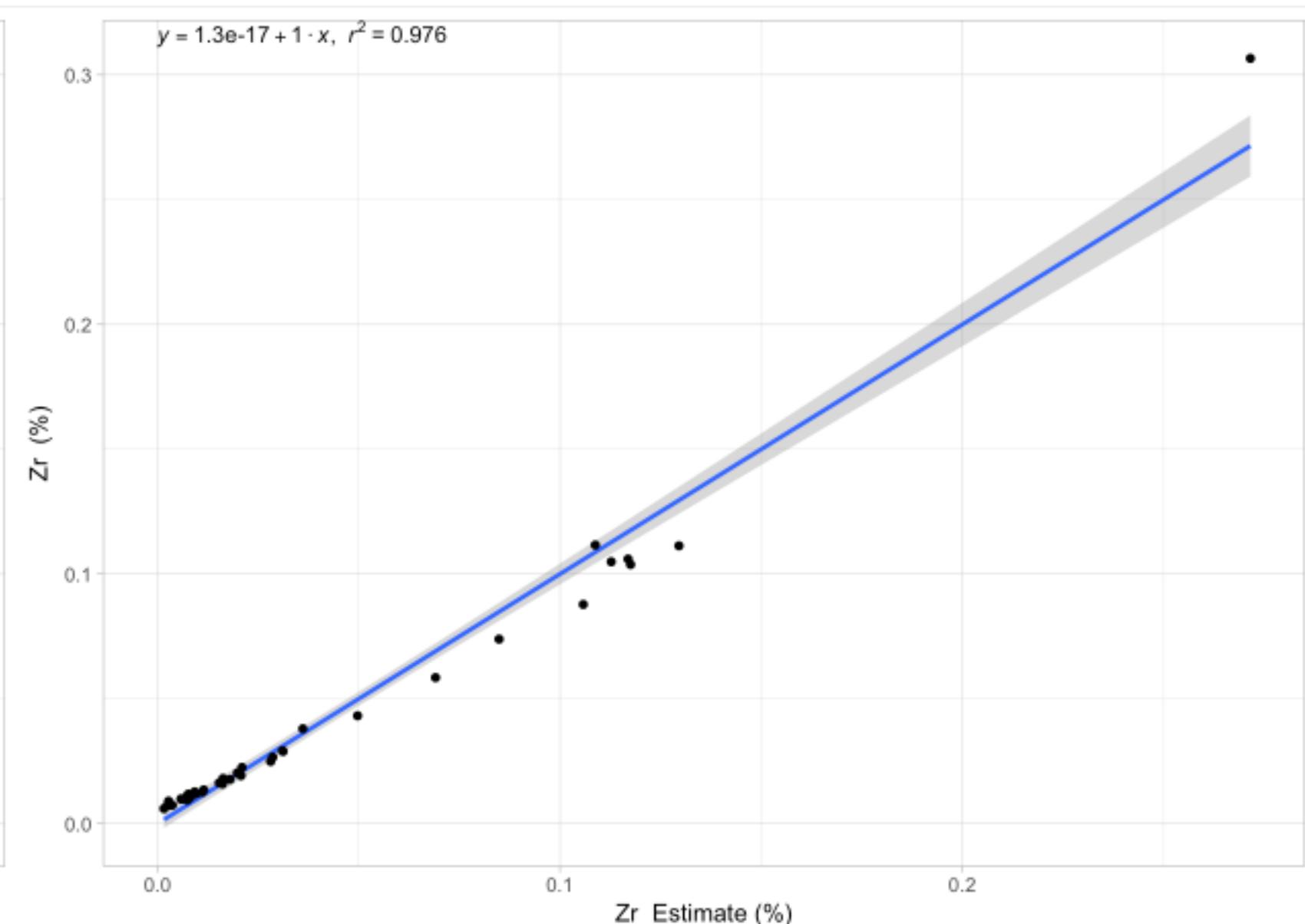
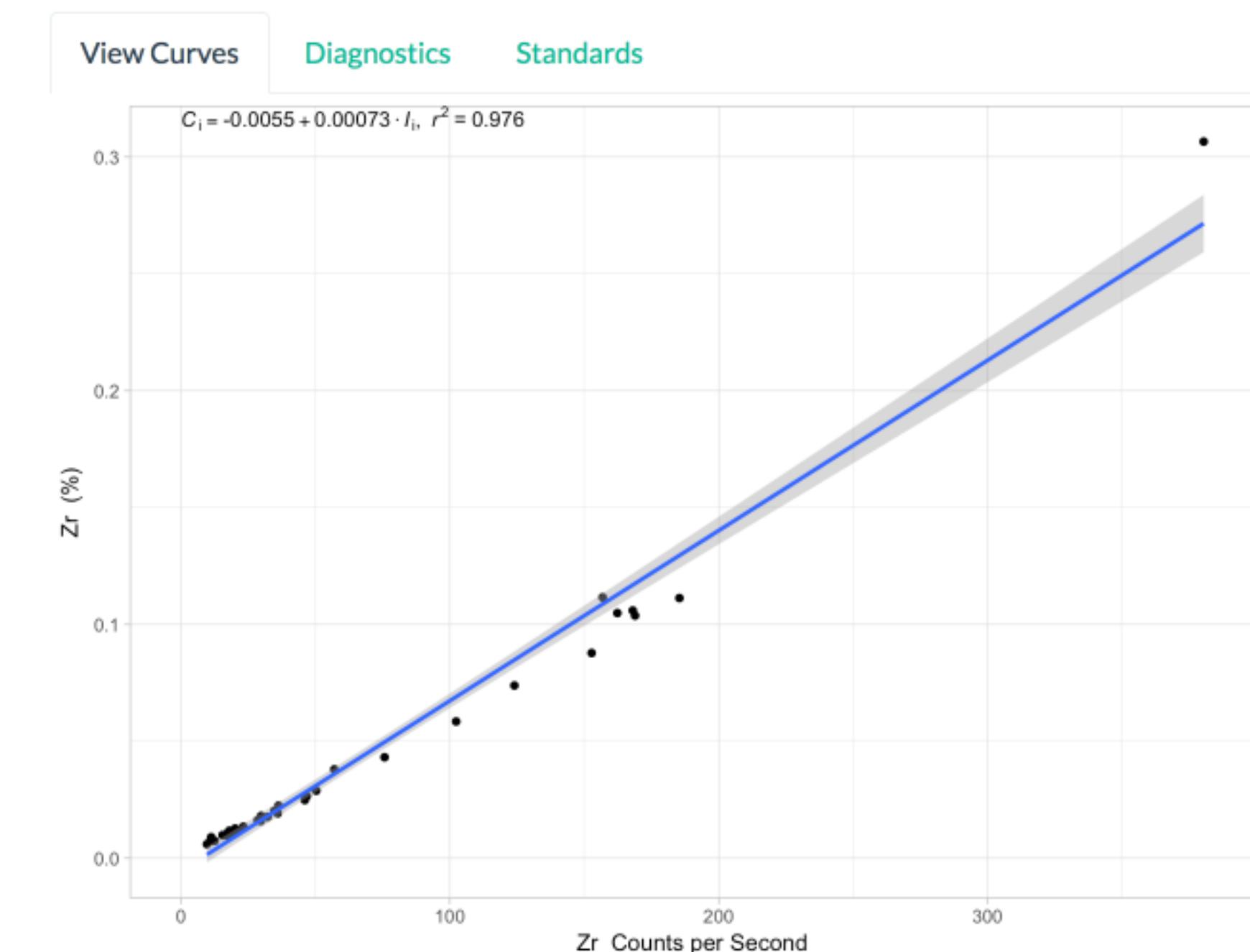
Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



The plot on the left is the calibration curve, with photon counts on the x-axis and known values on the y-axis

Plot **Update** **Save**

Element
Zr.Kalpha

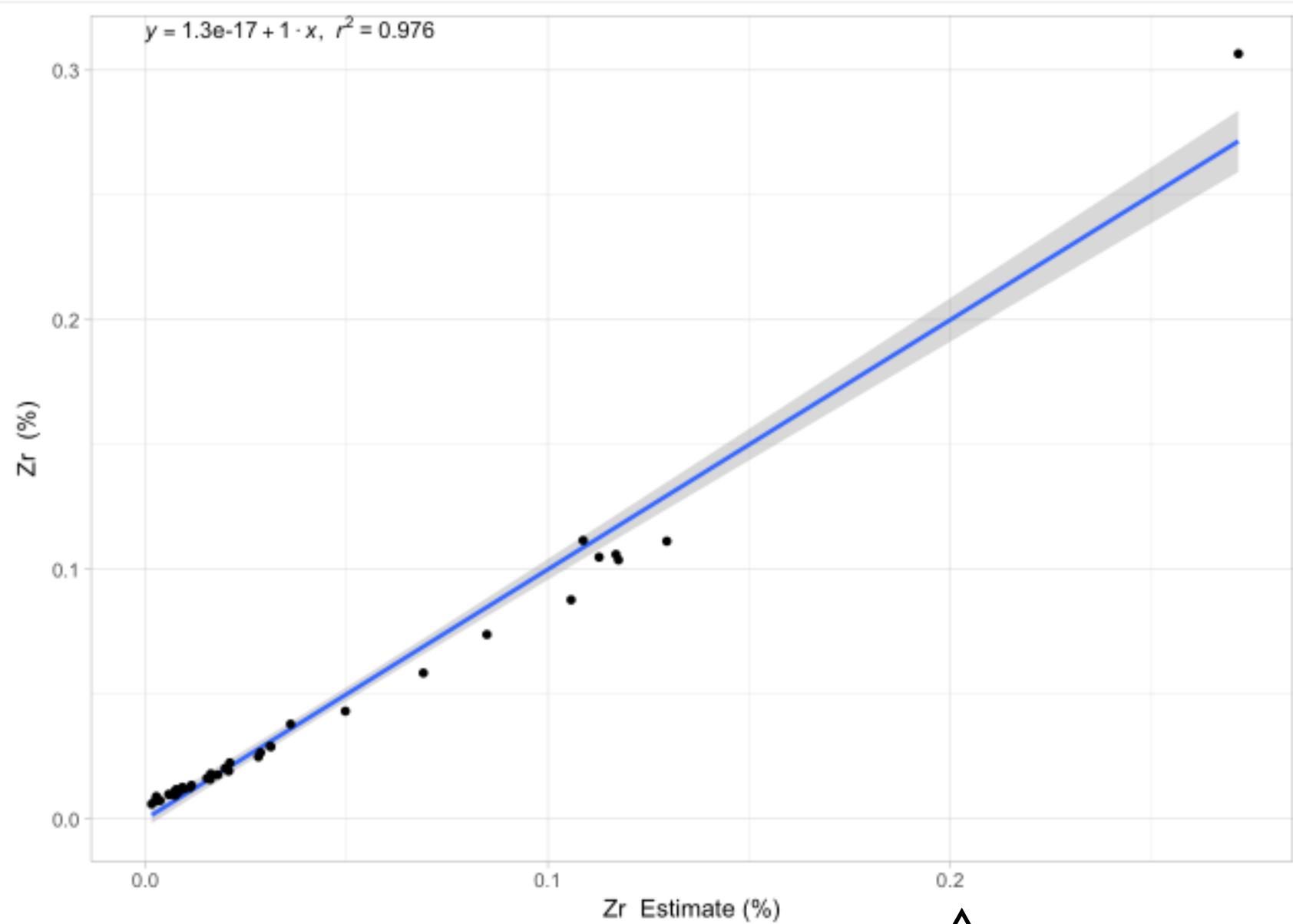
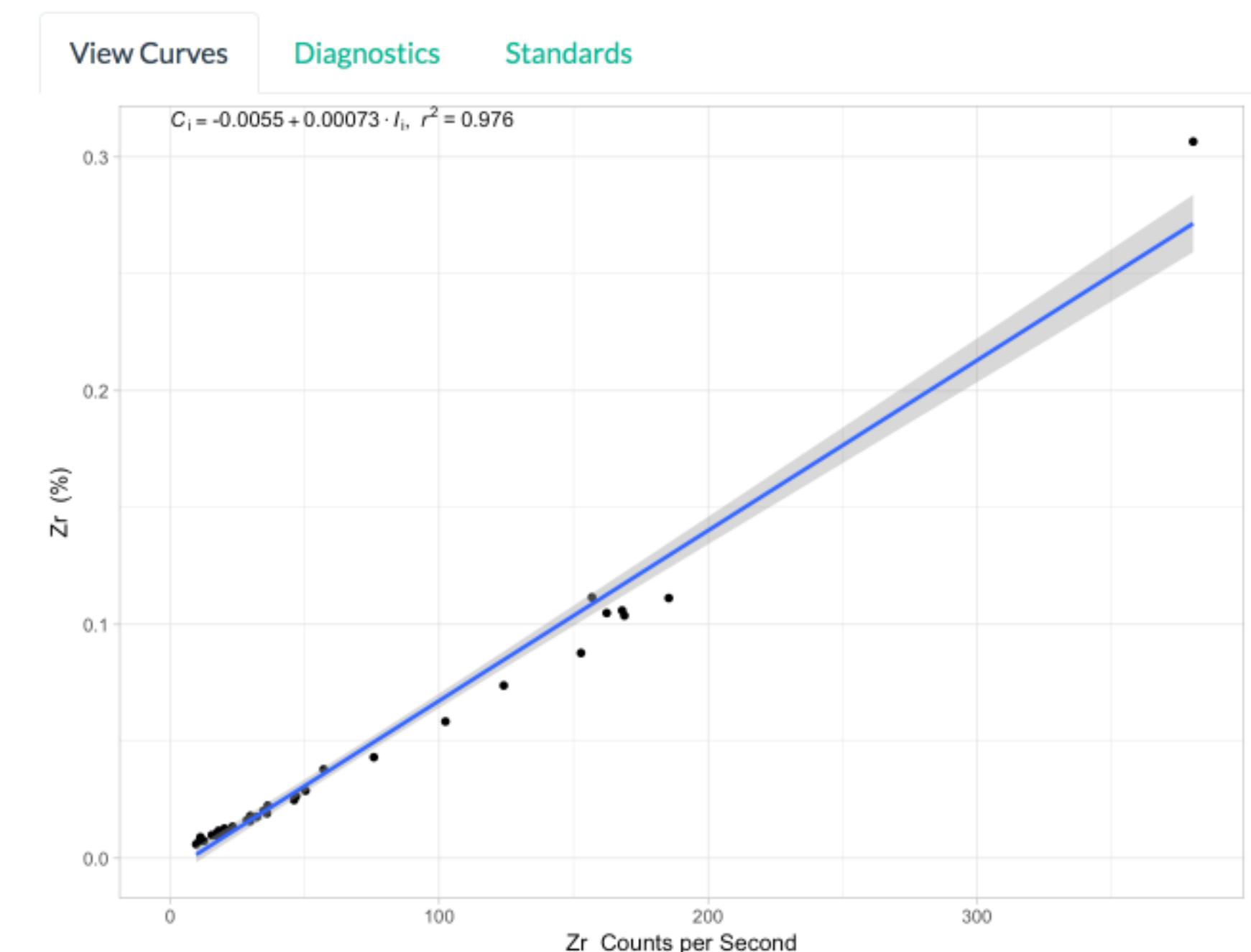
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



The plot on the right is the validation plot - with XRF estimates on the x-axis and known values on the y-axis

Plot **Update** **Save** 

Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

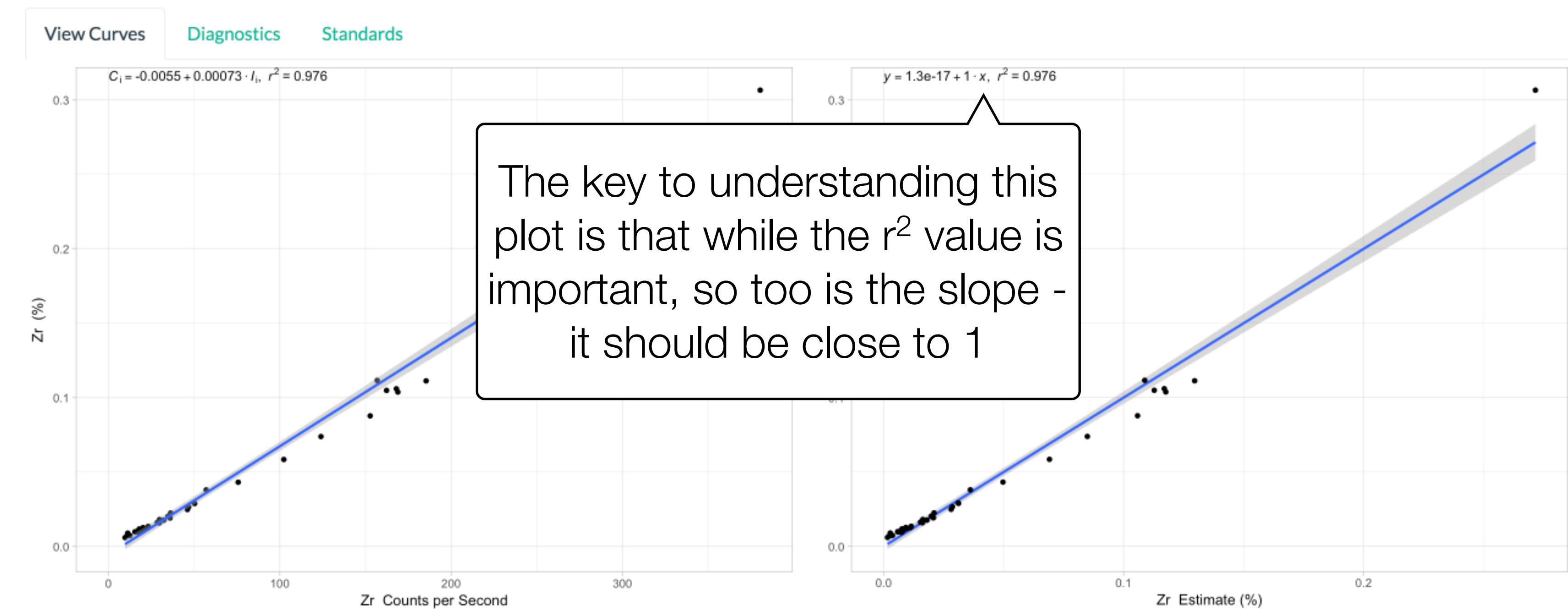
Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



CloudCal Spectrum Counts Add Concentrations Cal Curves Apply Calibration

Plot Update Save 

Element

Zr.Kalpha

Calibration Curve

Linear

Non-Linear

Lucas-Tooth

Normalization

Time

Total Counts

Compton

Min

18.5

Max

19.5

Intercept

K.K.alpha

Ca.K.alpha

Ti.K.alpha

Mn.K.alpha

Fe.K.alpha

Co.K.alpha

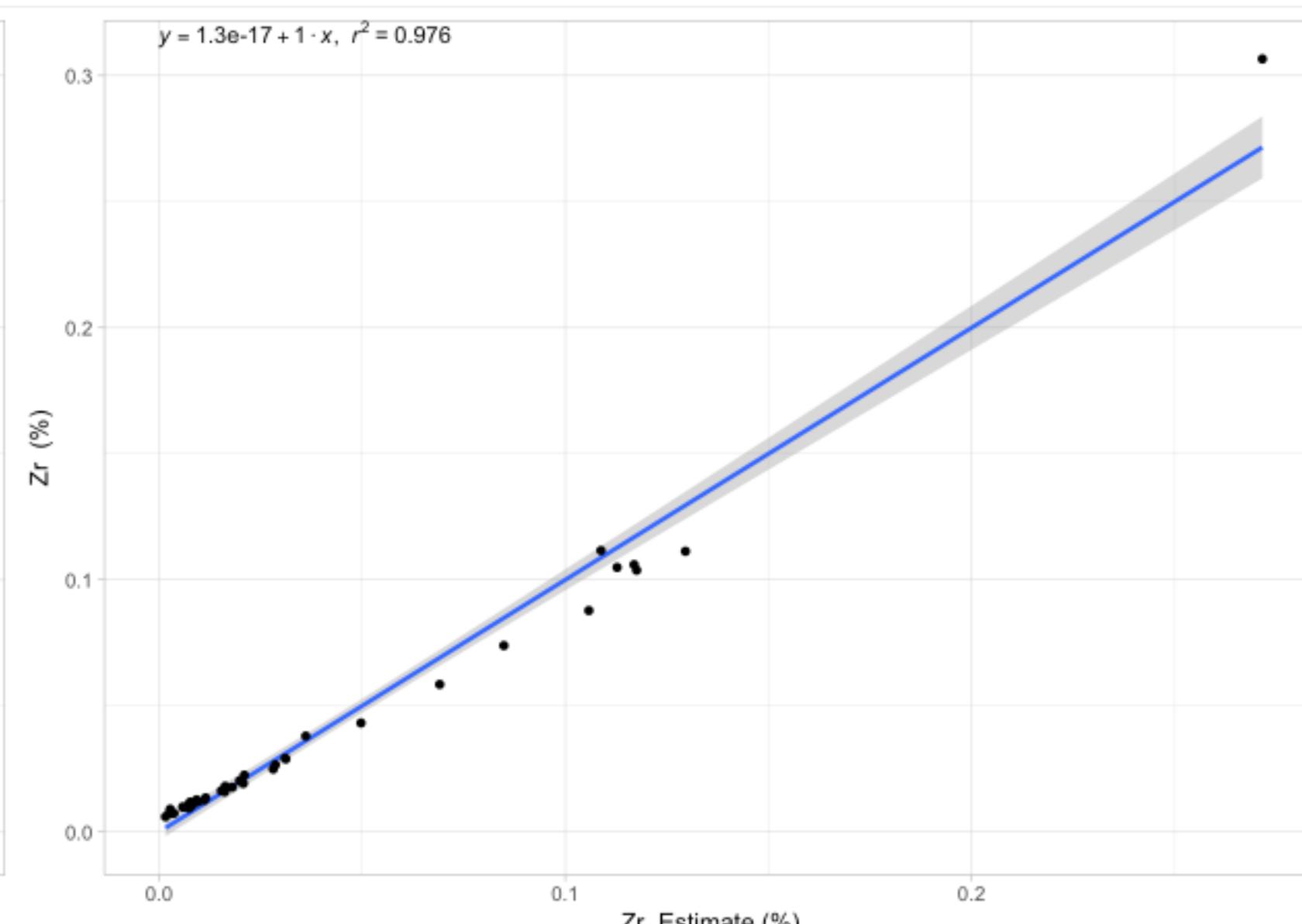
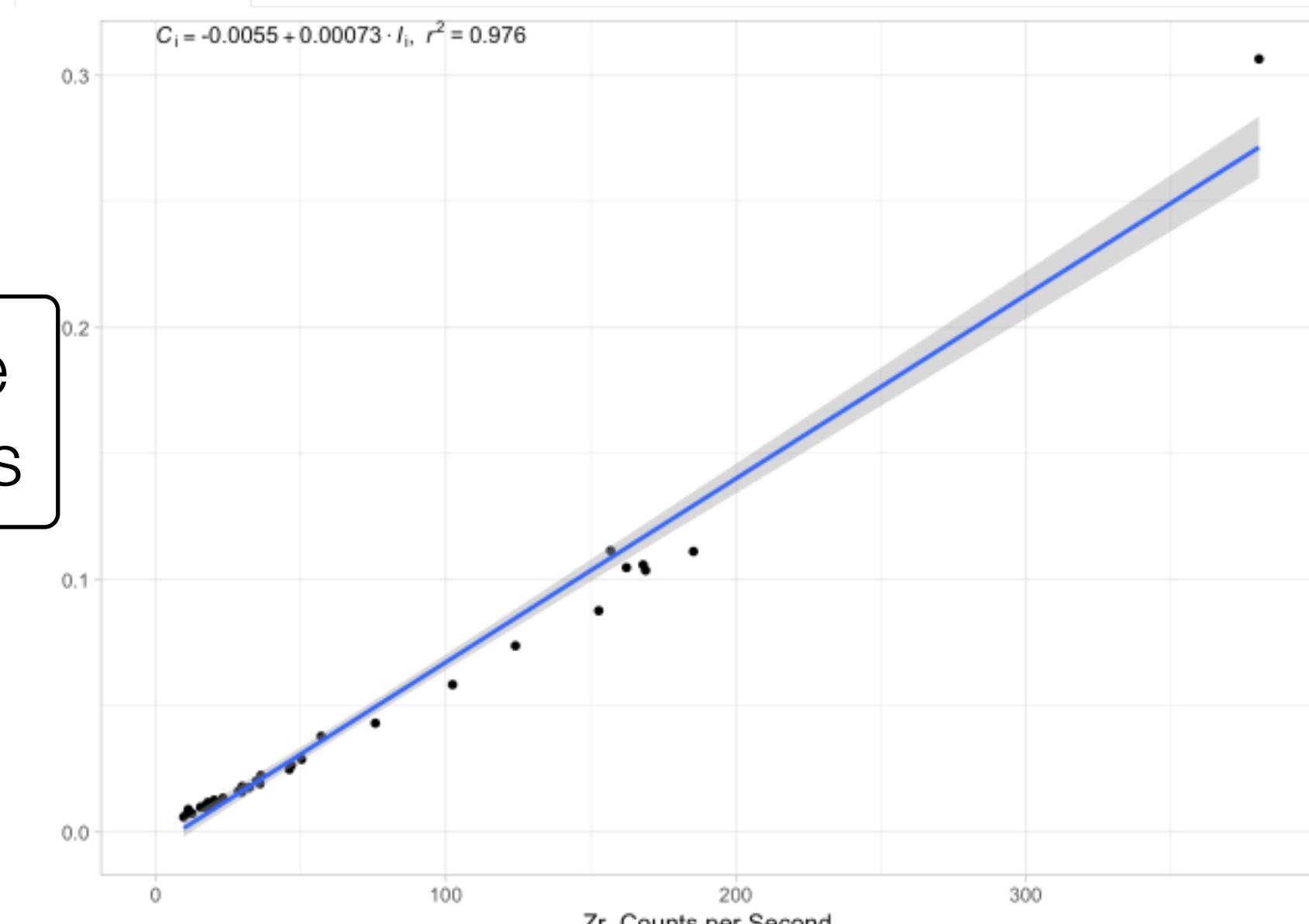
Cu.K.alpha

Zn.K.alpha

Ga.K.alpha

As.K.alpha

View Curves Diagnostics Standards



Currently, we are using a simple linear calibration - no corrections

We are also only normalizing data to time

Plot **Update** **Save**

Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

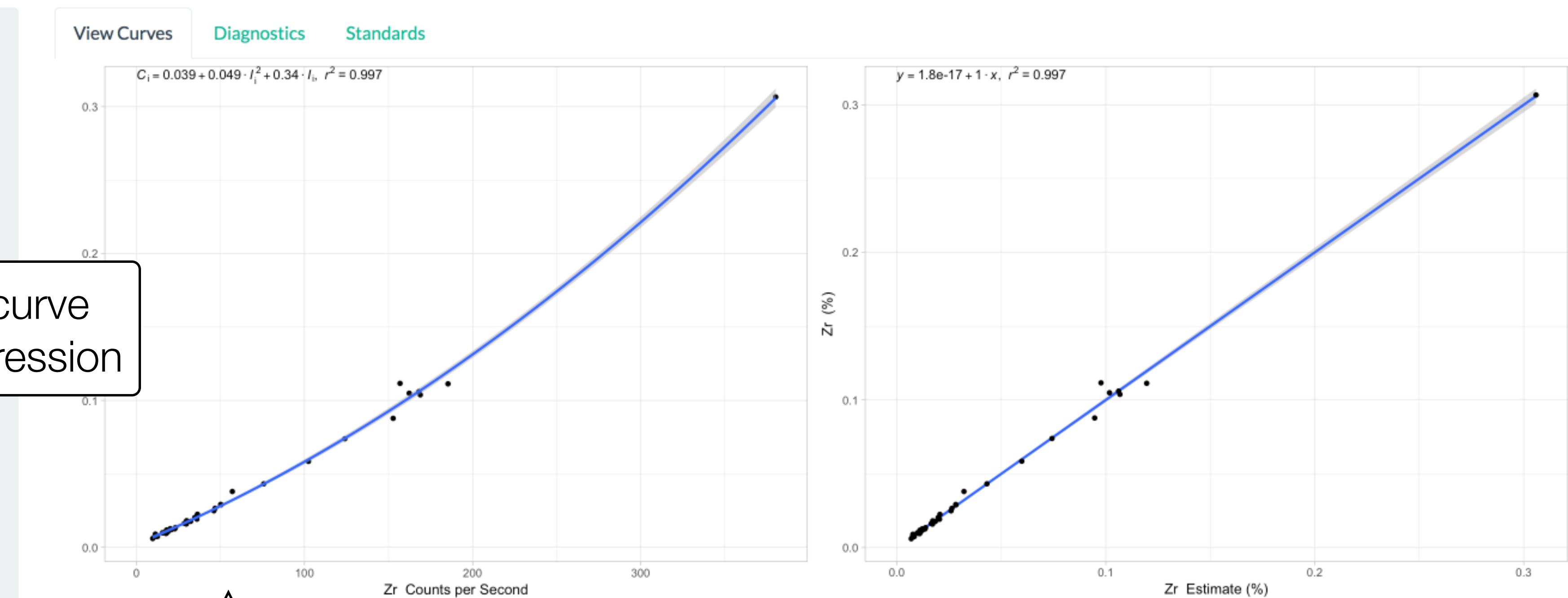
Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



Switching to a non-linear curve simply uses a quadratic regression

Note that while the cal curve plots is curvilinear, the validation plot remains linear

Plot Update Save 

Element

Zr.Kalpha

Calibration Curve

- Linear
- Non-Linear
- Lucas-Tooth

Normalization

- Time
- Total Counts
- Compton

Min

18.5

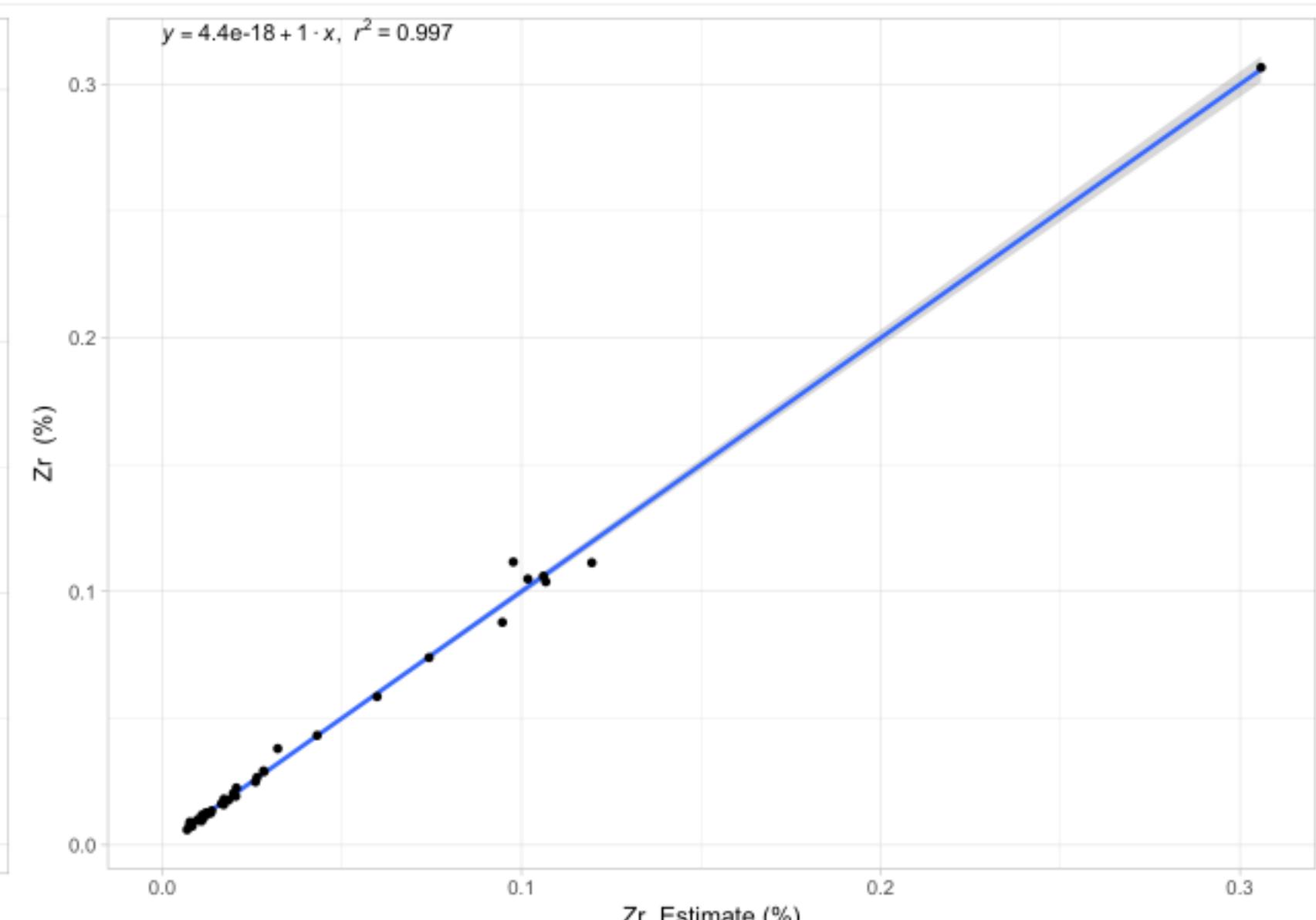
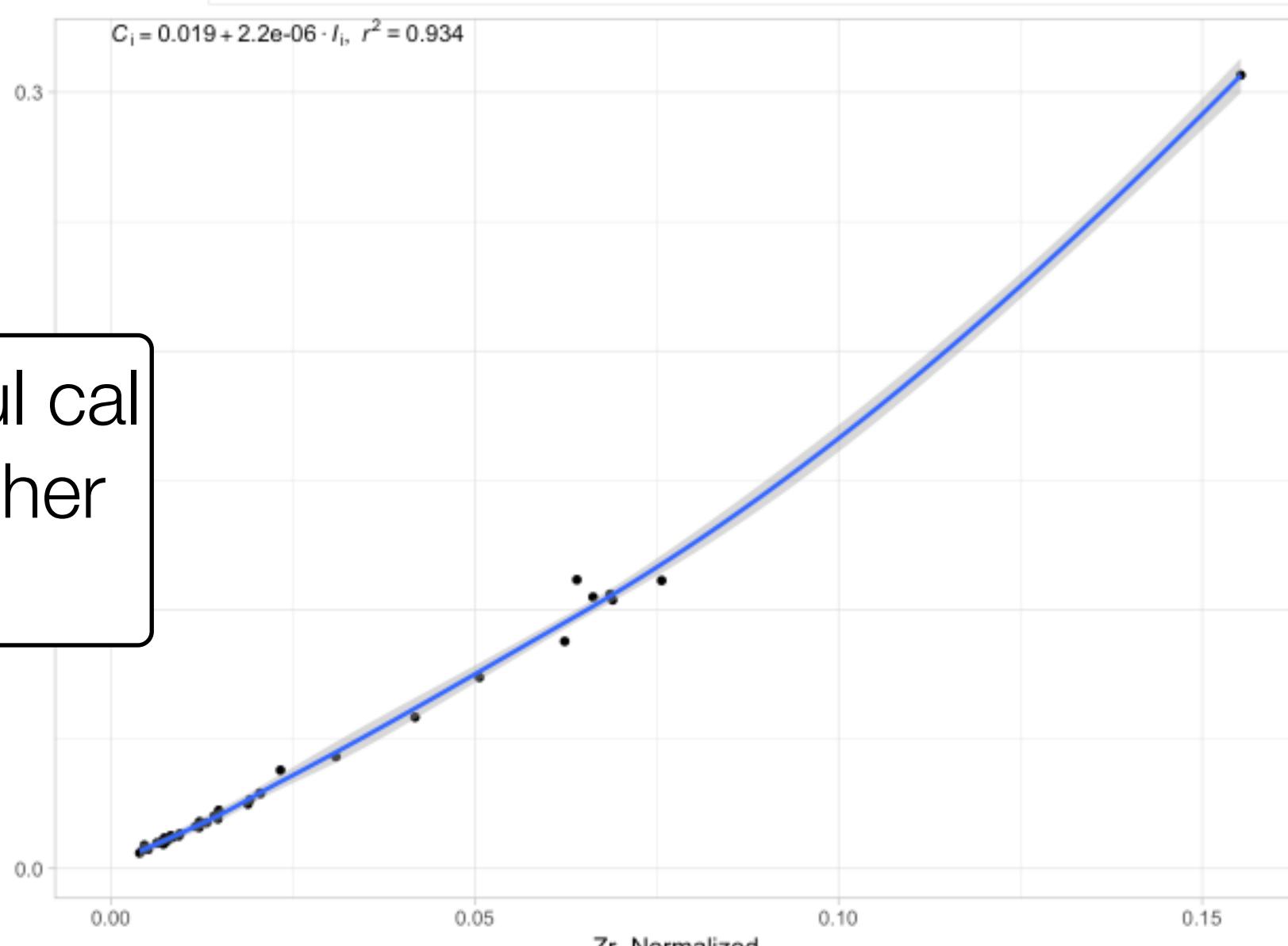
Max

19.5

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha

View Curves Diagnostics Standards



Lukas-Tooth is the most powerful cal engine - it enables the use of other elements as corrections

Plot **Update** **Save** 

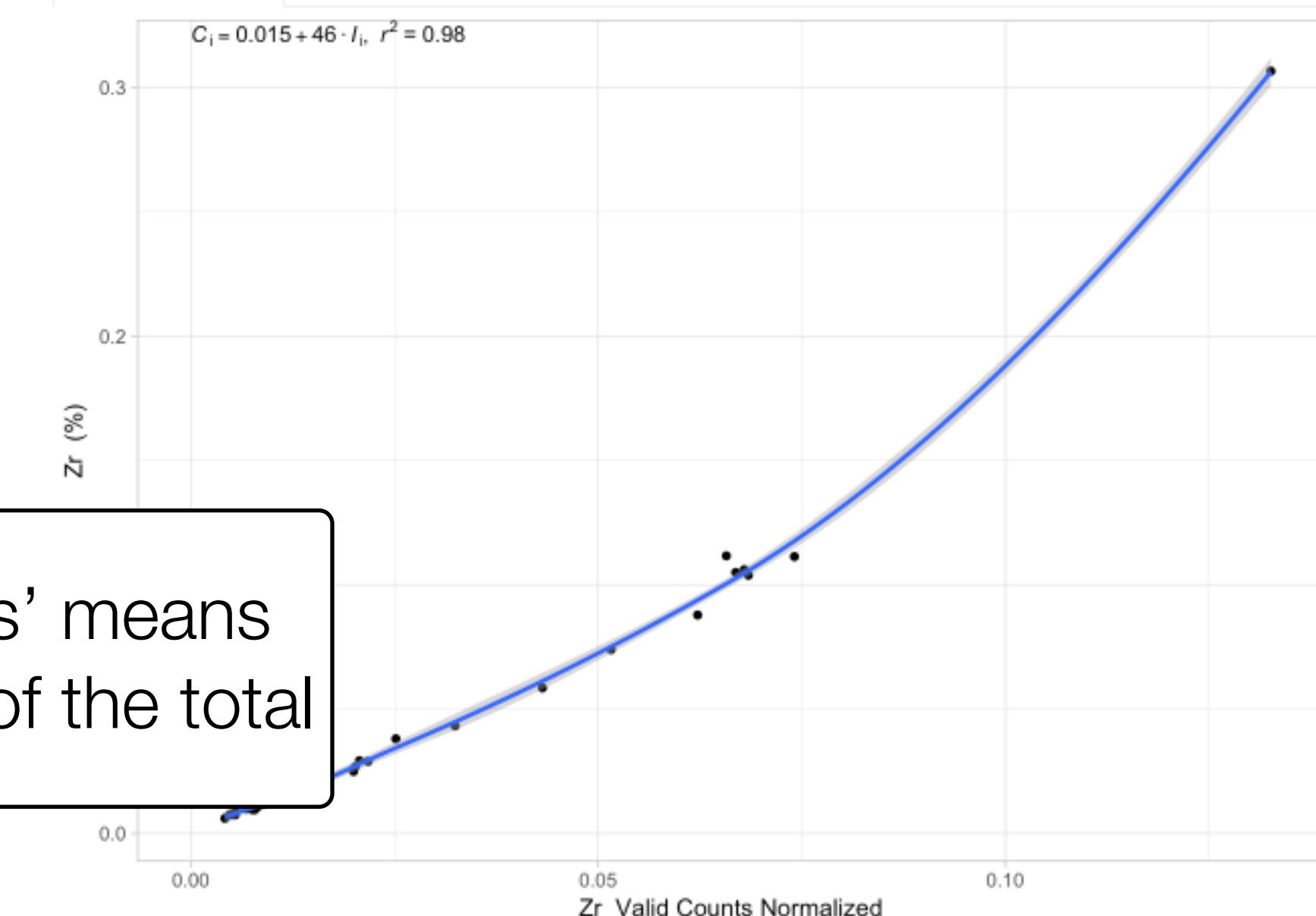
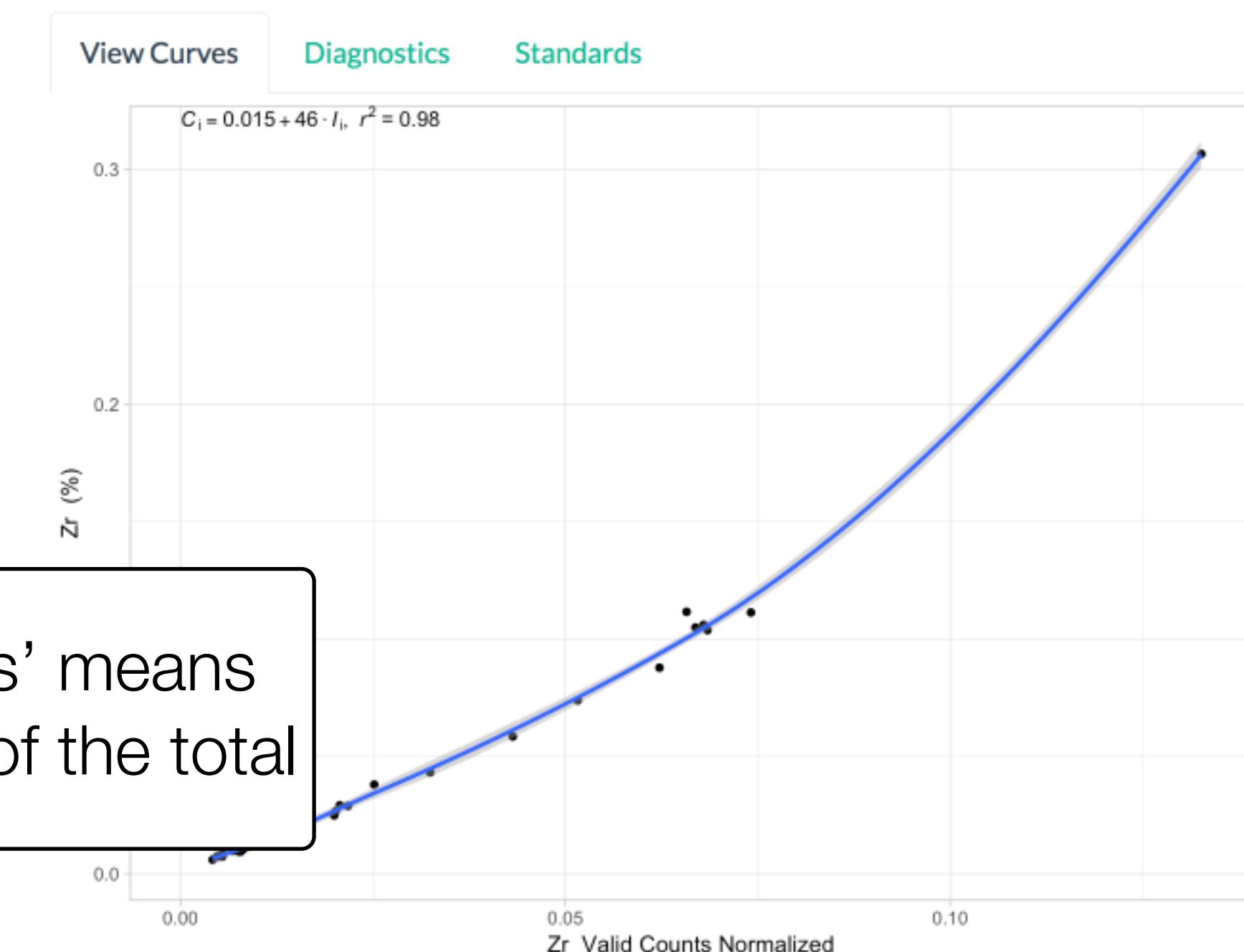
Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5



Normalizing to 'Total Counts' means we are using the peak as % of the total

Plot Update Save

Element
Zr.Kalpha

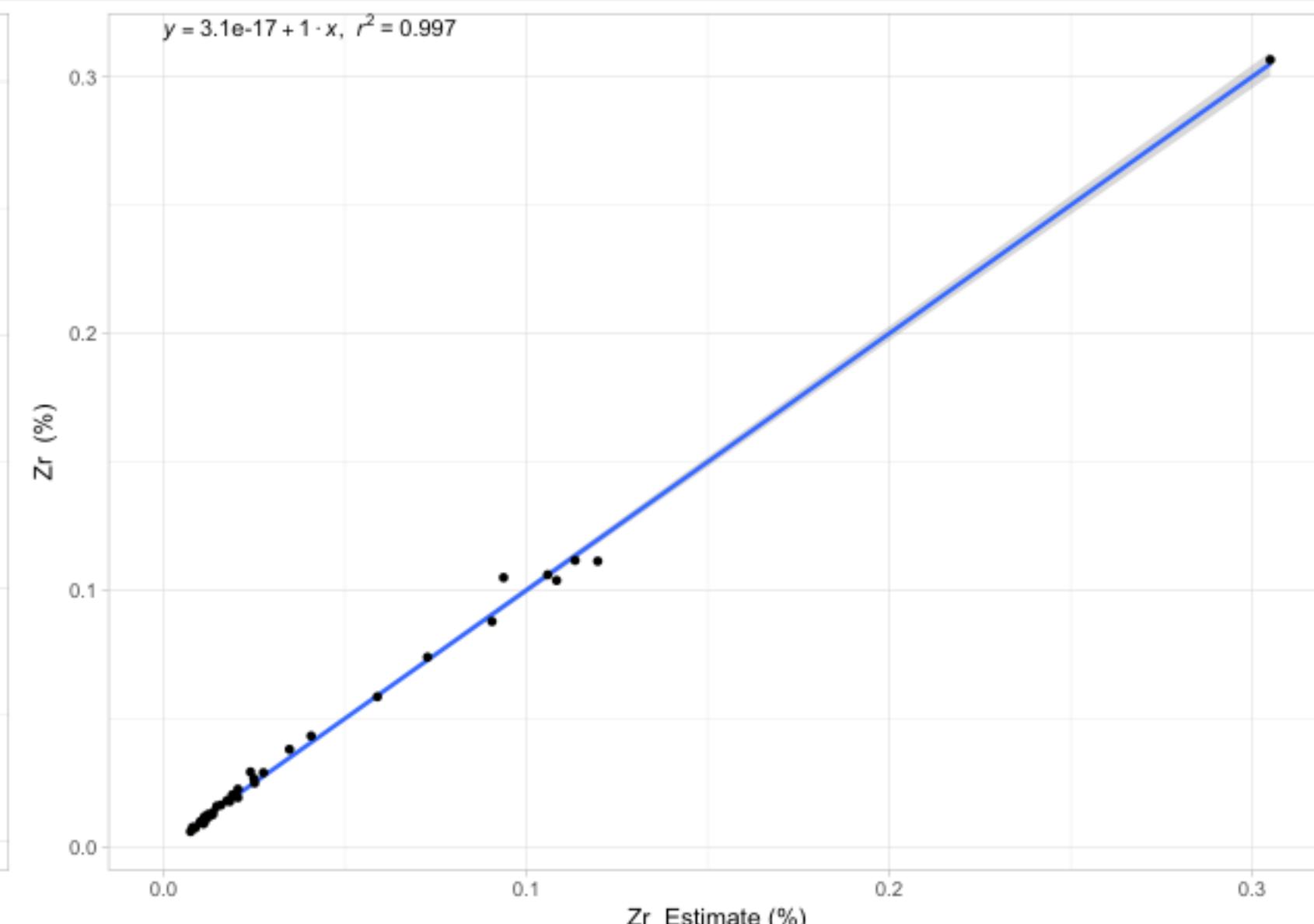
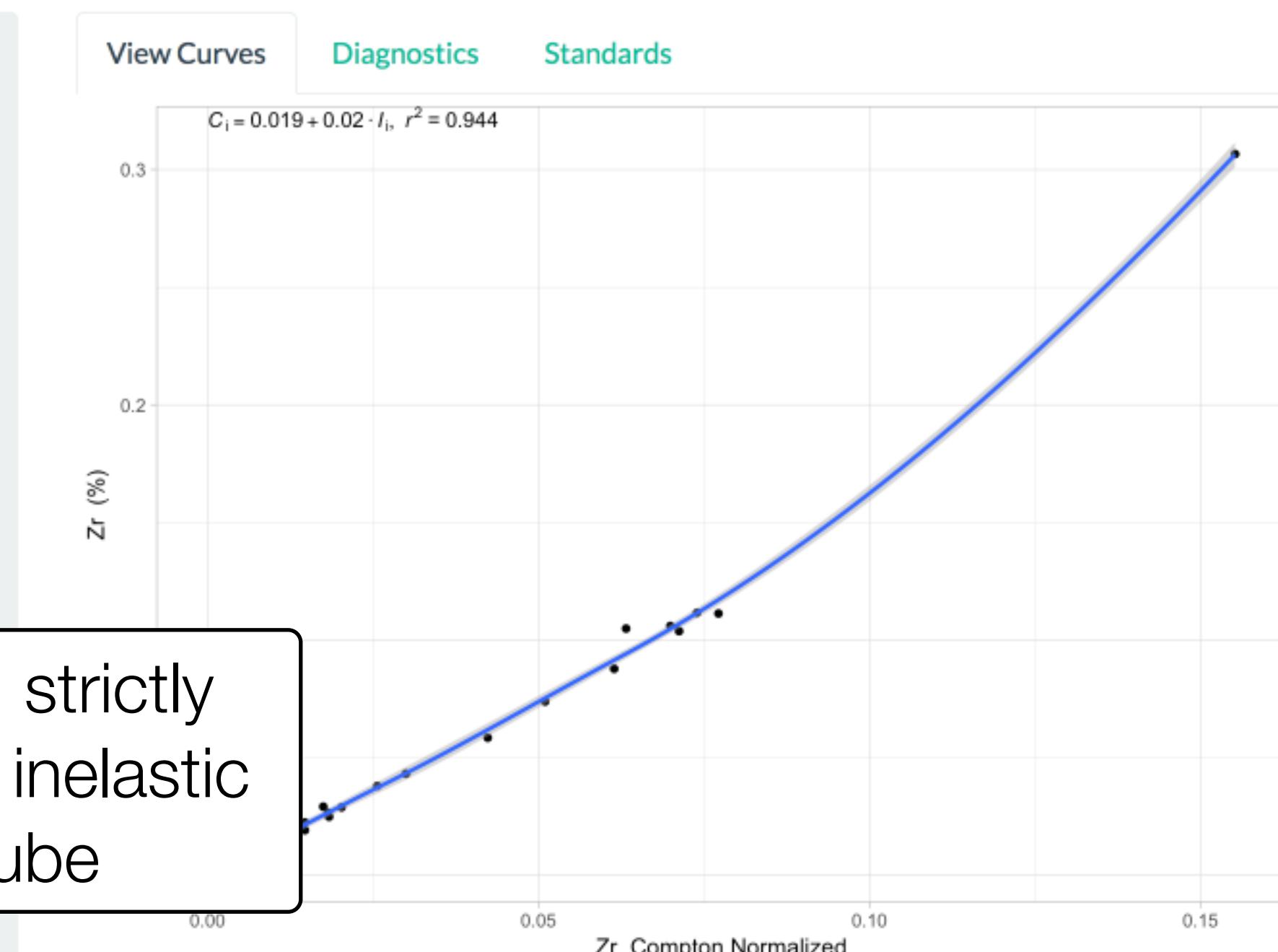
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



Normalizing to Compton is, strictly speaking, normalizing to the inelastic scatter from the x-ray tube

Plot Update Save

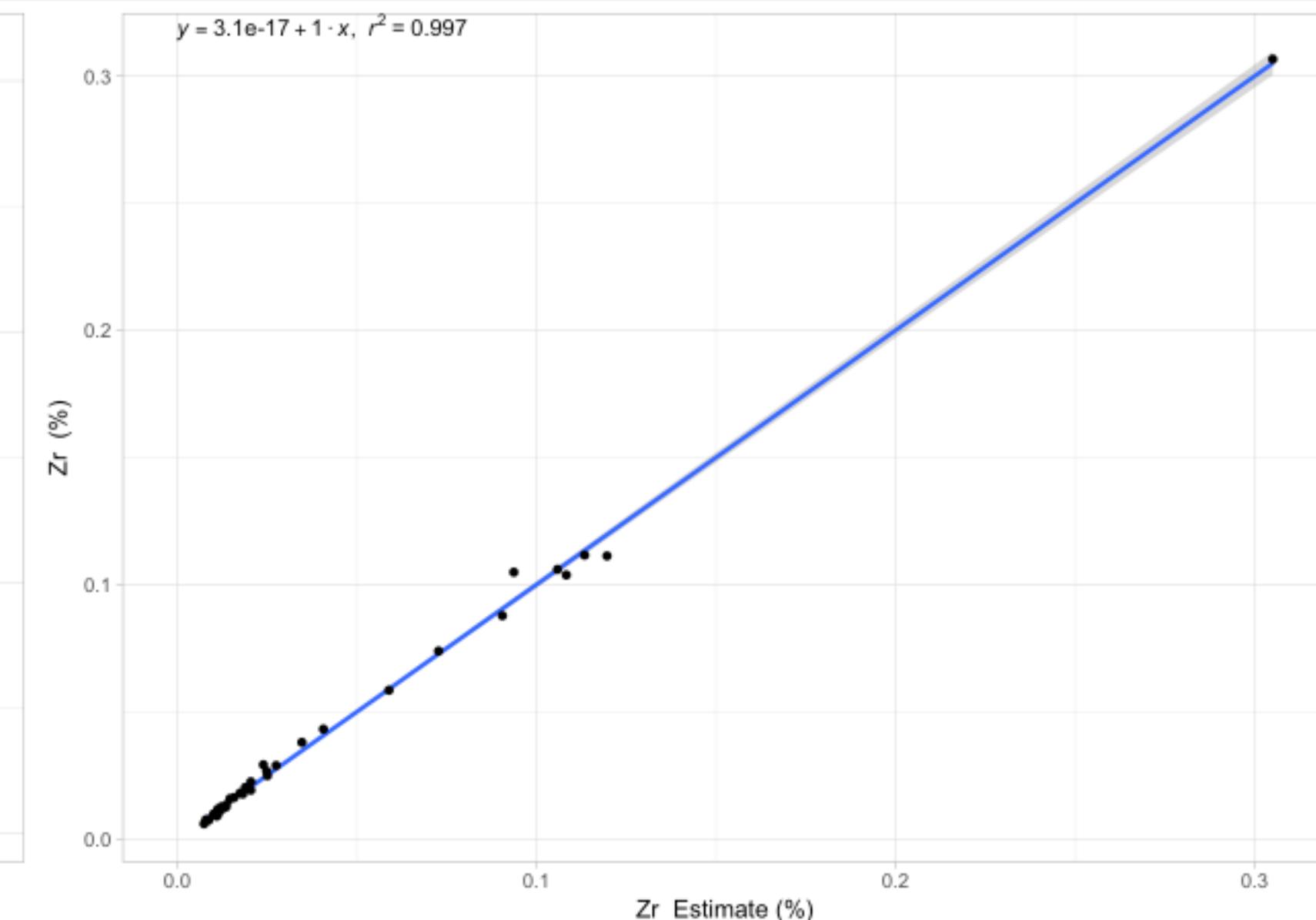
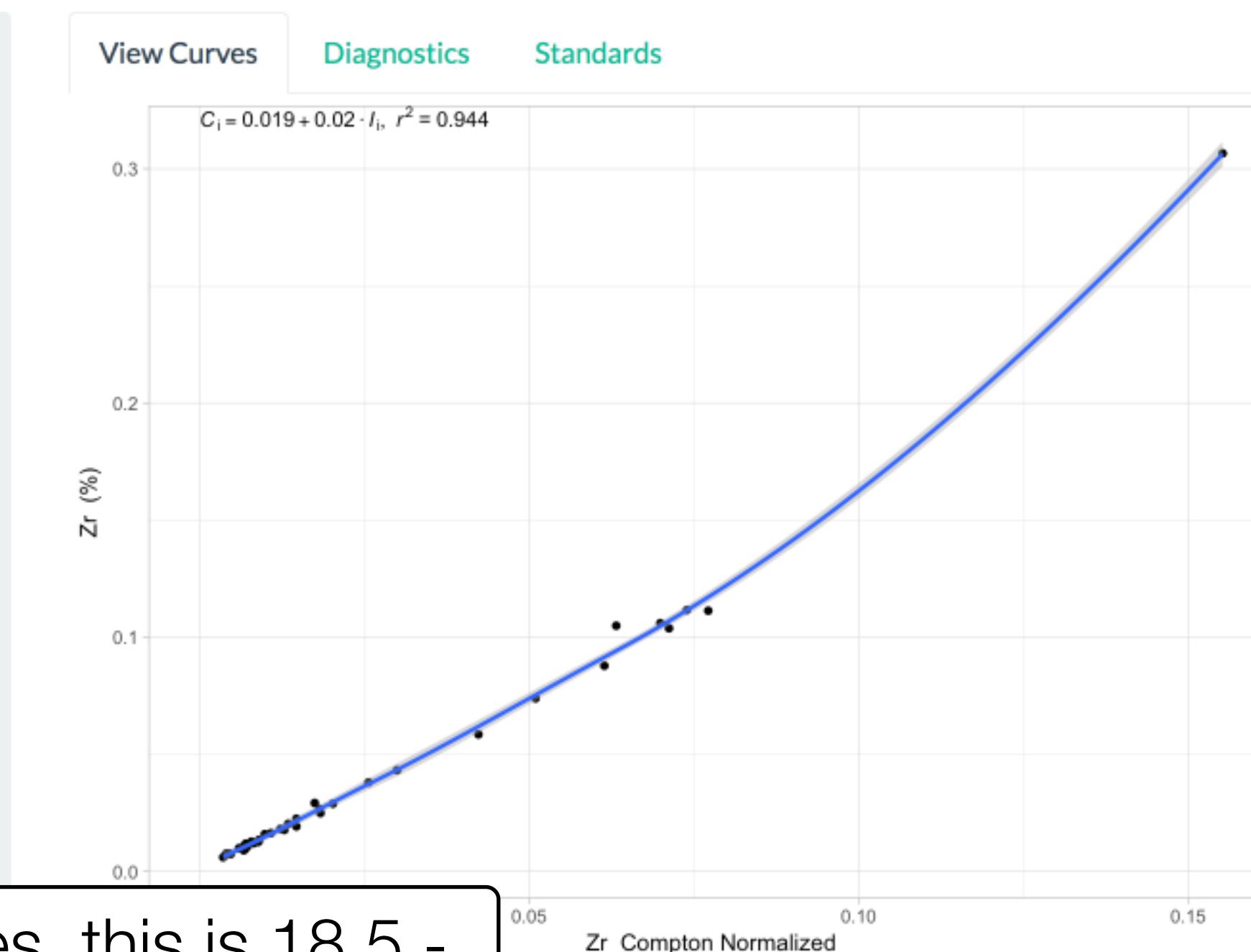
Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
18.5

Max
19.5



For Rhodium x-ray tubes, this is 18.5 - 19.5 keV, though this can shift depending on the density of your sample

...however, you can normalize to any range using this menu. You can even do this separately for each element

Plot **Update** **Save** 

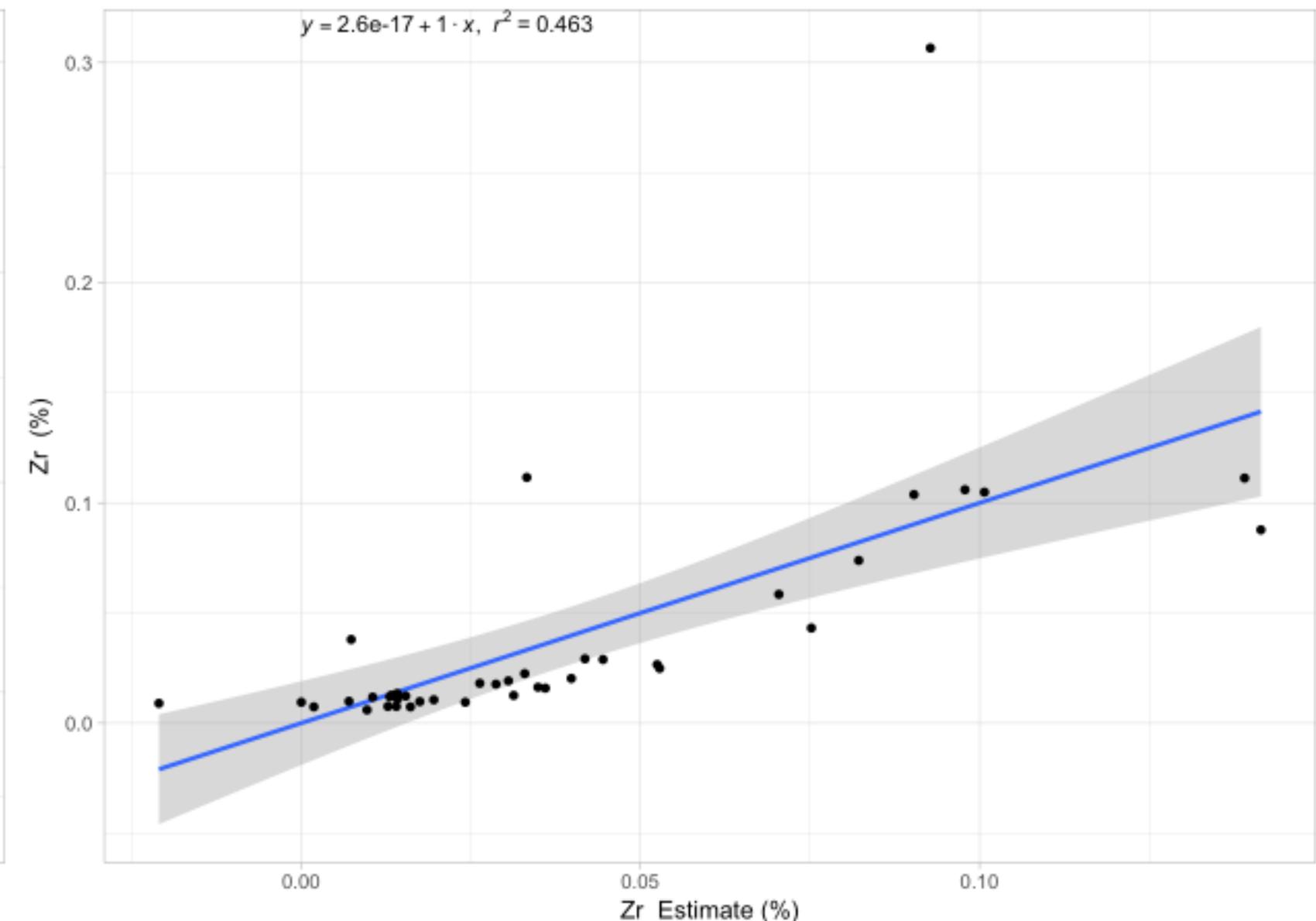
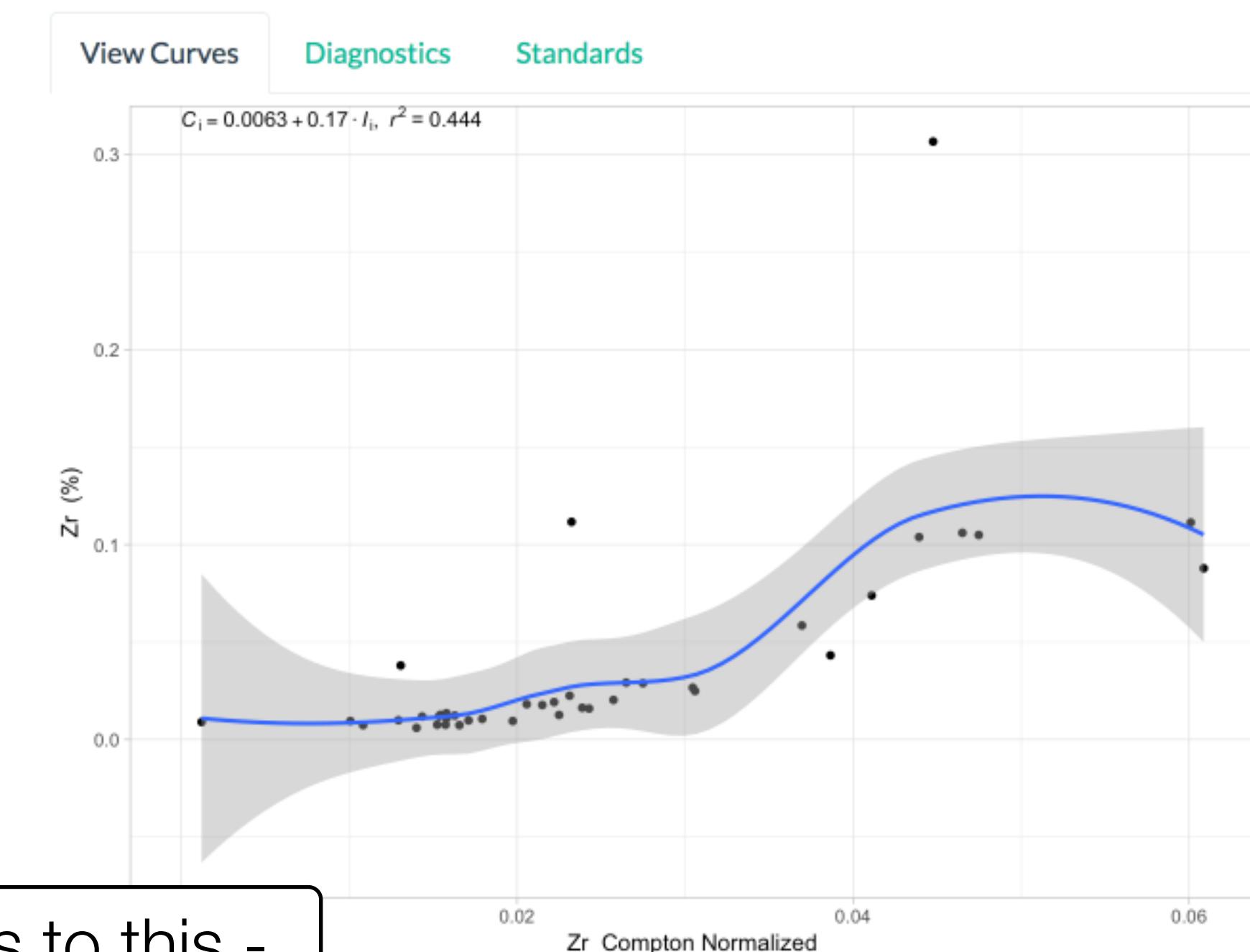
Element
Zr.Kalpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
6

Max
7



There are consequences to this -
choosing a bad range can disrupt
your calibration quality

Plot **Update** **Save** **Download**

Element
Zr.Kalpha

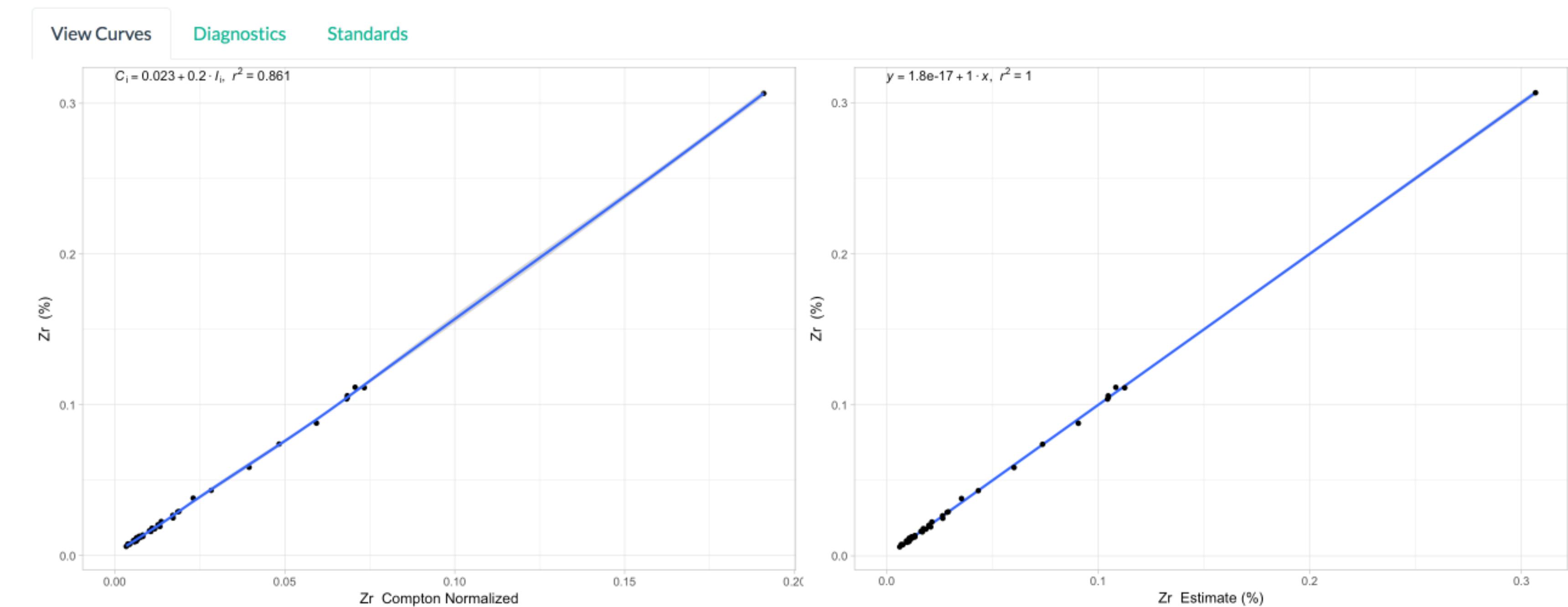
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
19.5

Max
22

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



For this element and this set, we have hit the theoretical best performance - an r^2 of 1, a slope of 1. This is not common

- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- As.K.alpha
- Rb.K.alpha
- Sr.K.alpha
- Y.K.alpha
- Zr.K.alpha
- Nb.K.alpha
- Pb.L.alpha
- Th.L.alpha
- U.L.alpha
- None

Slope

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- As.K.alpha
- Rb.K.alpha
- Sr.K.alpha
- Y.K.alpha
- Zr.K.alpha
- Nb.K.alpha
- Pb.L.alpha
- Th.L.alpha
- U.L.alpha
- None

Nonetheless, this element can still be influenced by an overlap with Sr - you can add that as a slope or intercept correction

Plot **Update** **Save** **Download**

To show the effects of your corrections, click on 'Plot'

Calibration Curve

- Linear
- Non-Linear
- Lucas-Tooth

Normalization

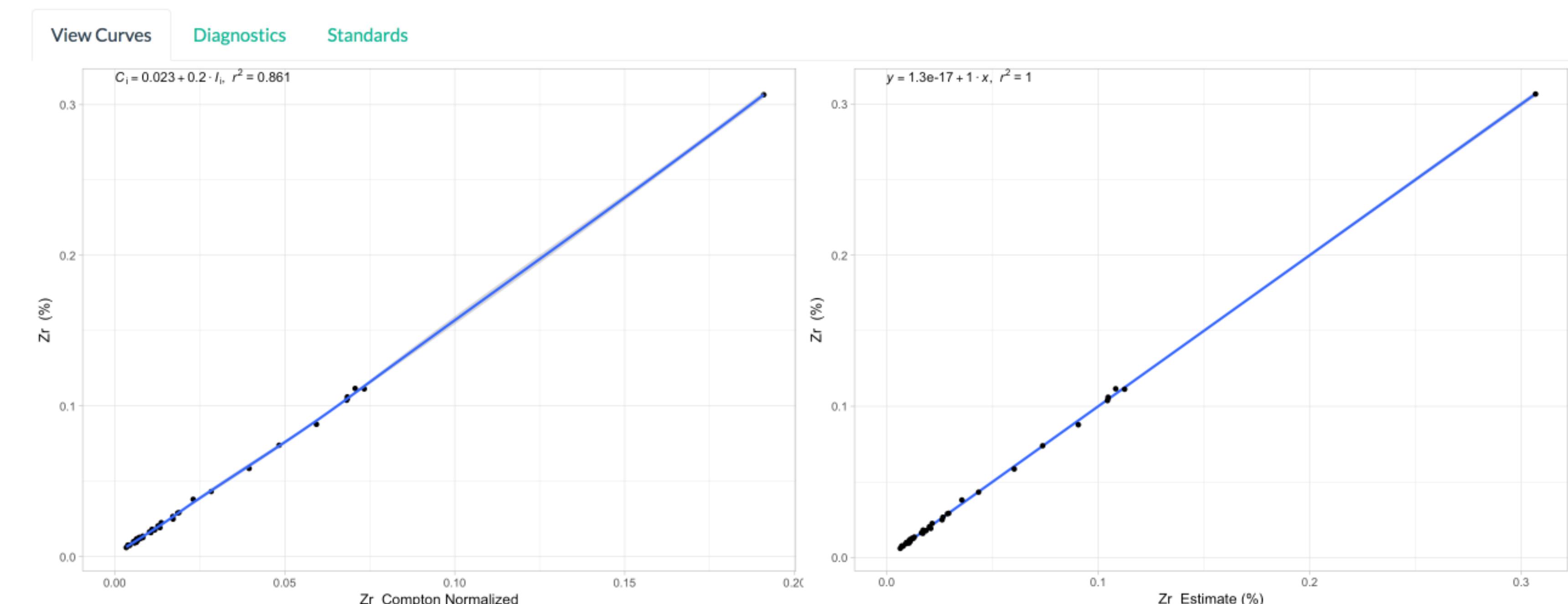
- Time
- Total Counts
- Compton

Min
19.5

Max
22

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



Plot Update Save

Element Zr.Kalph

Calibration Curve

- Linear
- Non-Linear
- Lucas-Tooth

Normalization

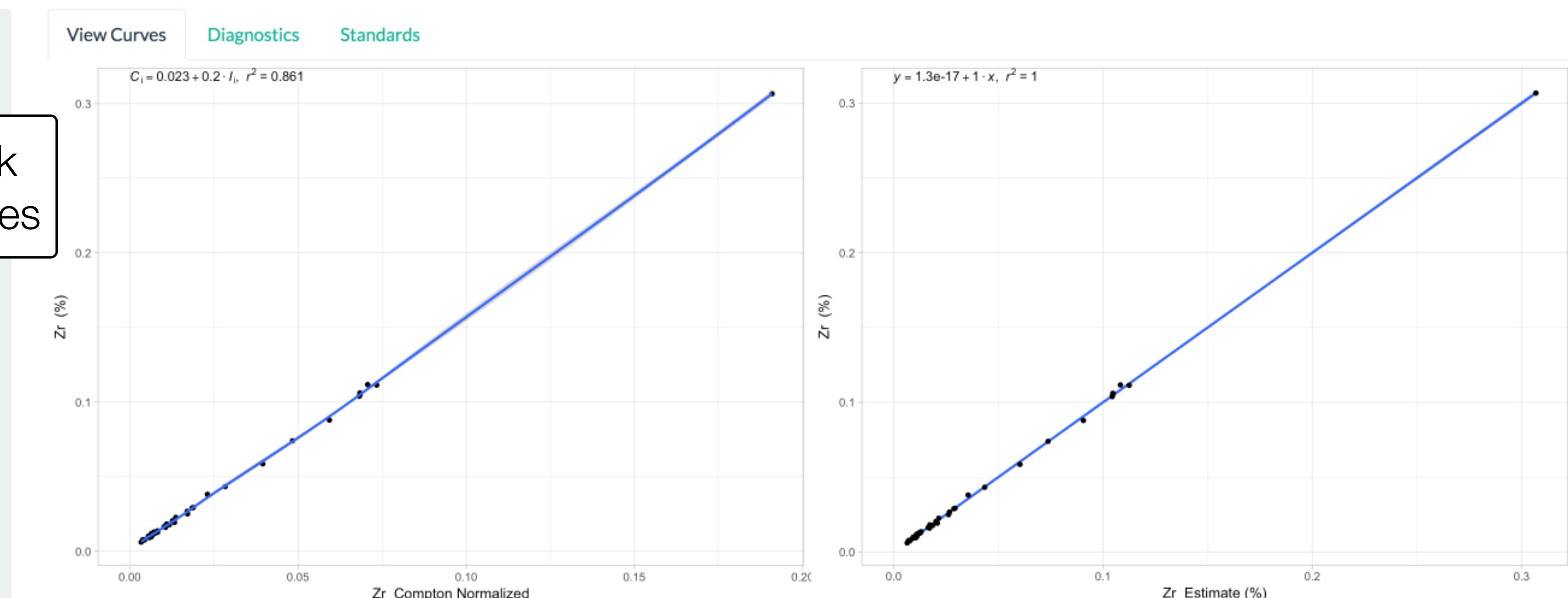
- Time
- Total Counts
- Compton

Min 19.5

Max 22

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



If you are satisfied, you can click
'Update' to commit these changes

Plot Update Save

Element
Rb.K.alpha

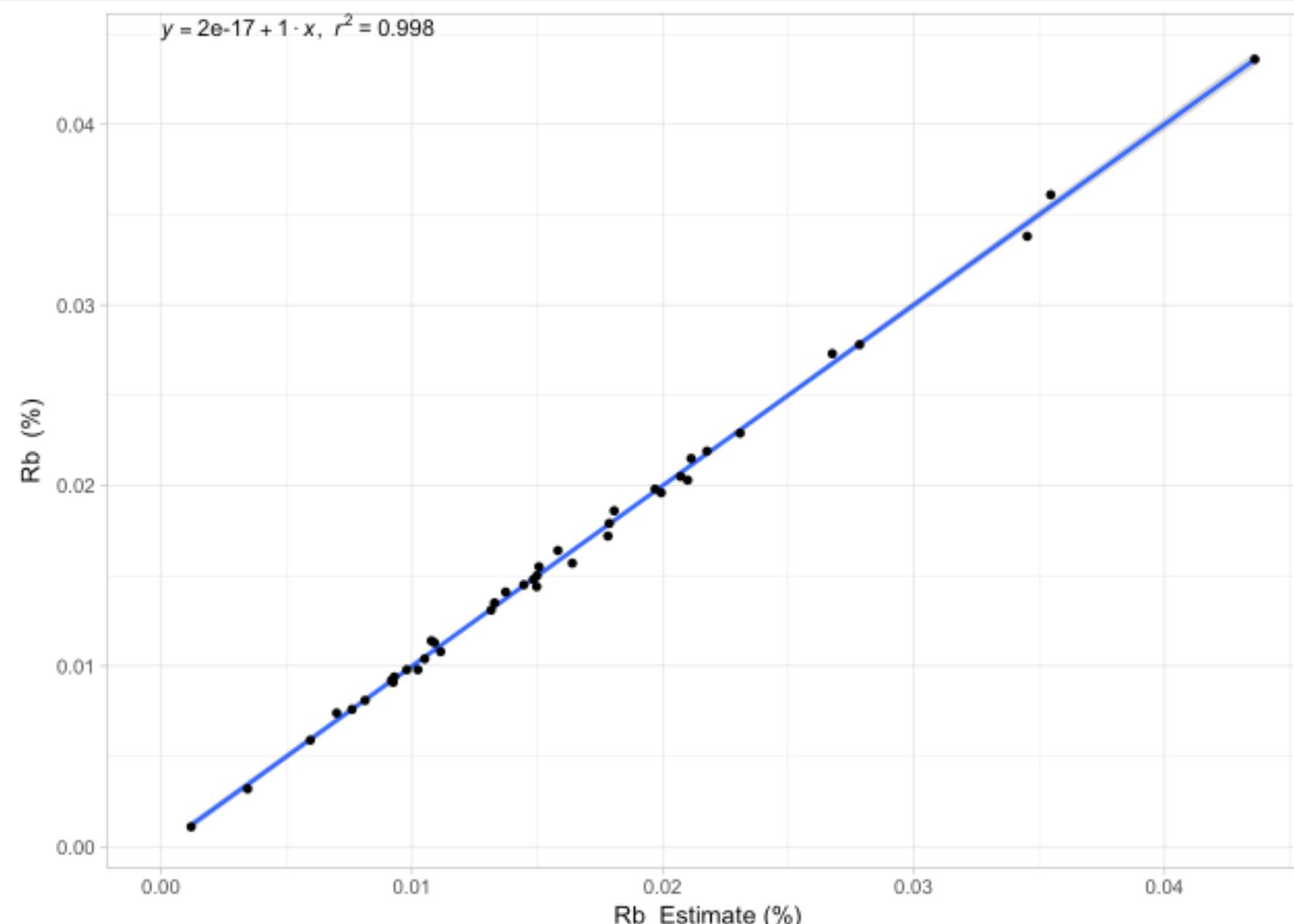
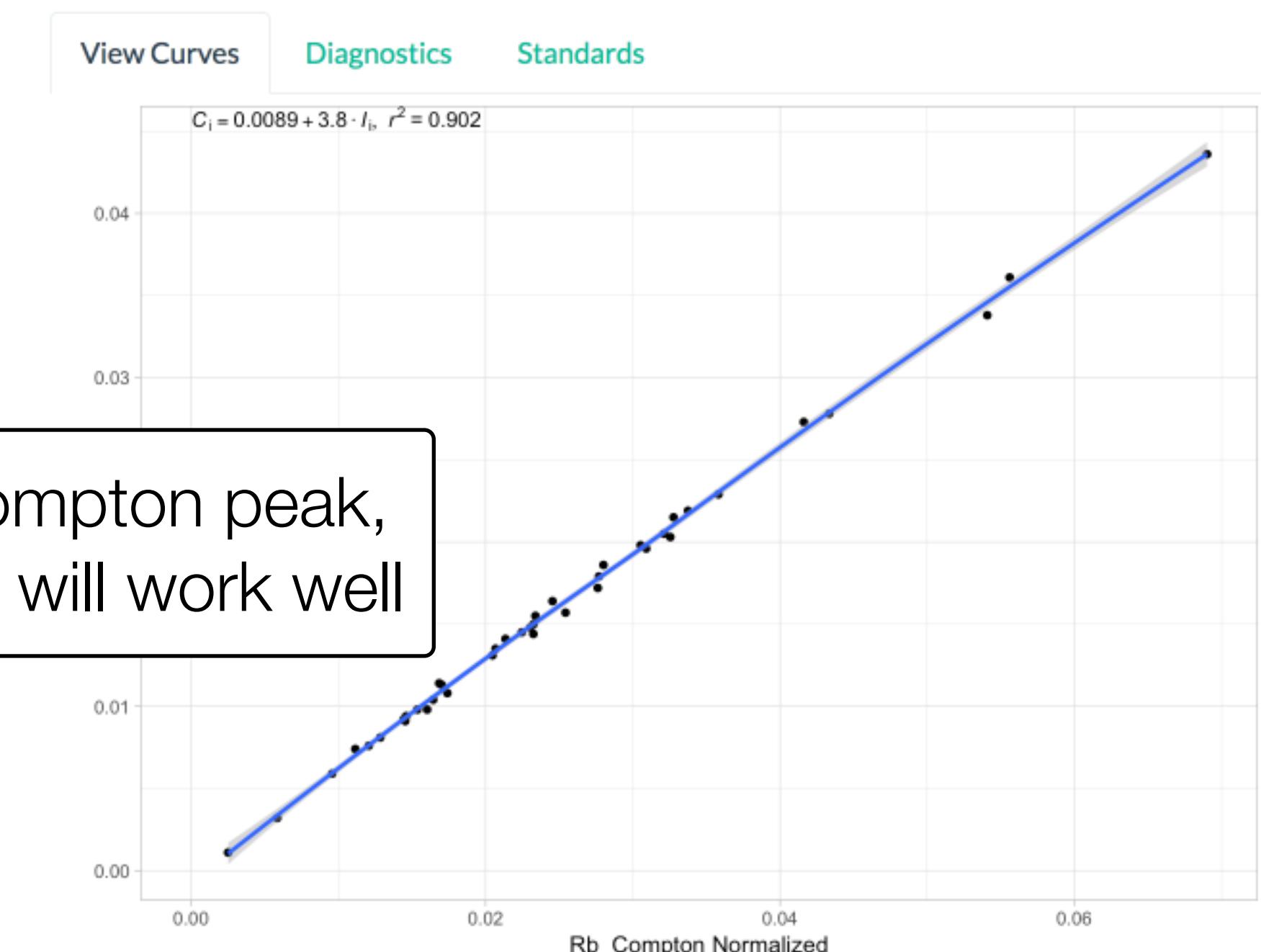
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
19.5

Max
22

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



For elements near the Compton peak,
that form of normalization will work well

Plot Update Save

Element
K.Kalpha

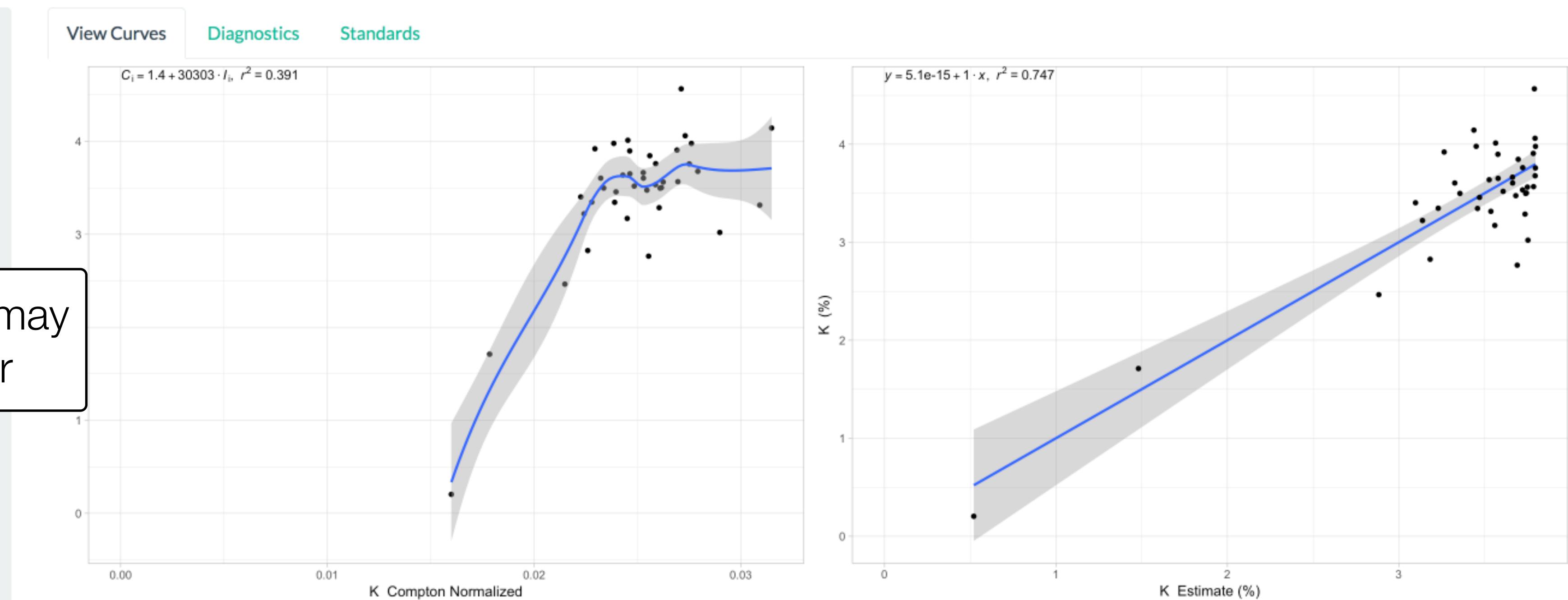
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
19.5

Max
22

Intercept
 K.Kalpha
 Ca.Kalpha
 Ti.Kalpha
 Mn.Kalpha
 Fe.K.alpha
 Co.Kalpha
 Cu.Kalpha
 Zn.Kalpha
 Ga.Kalpha
 Al.Kalpha



For other elements, this may not work well however

X-Ray Fluorescence Calibration

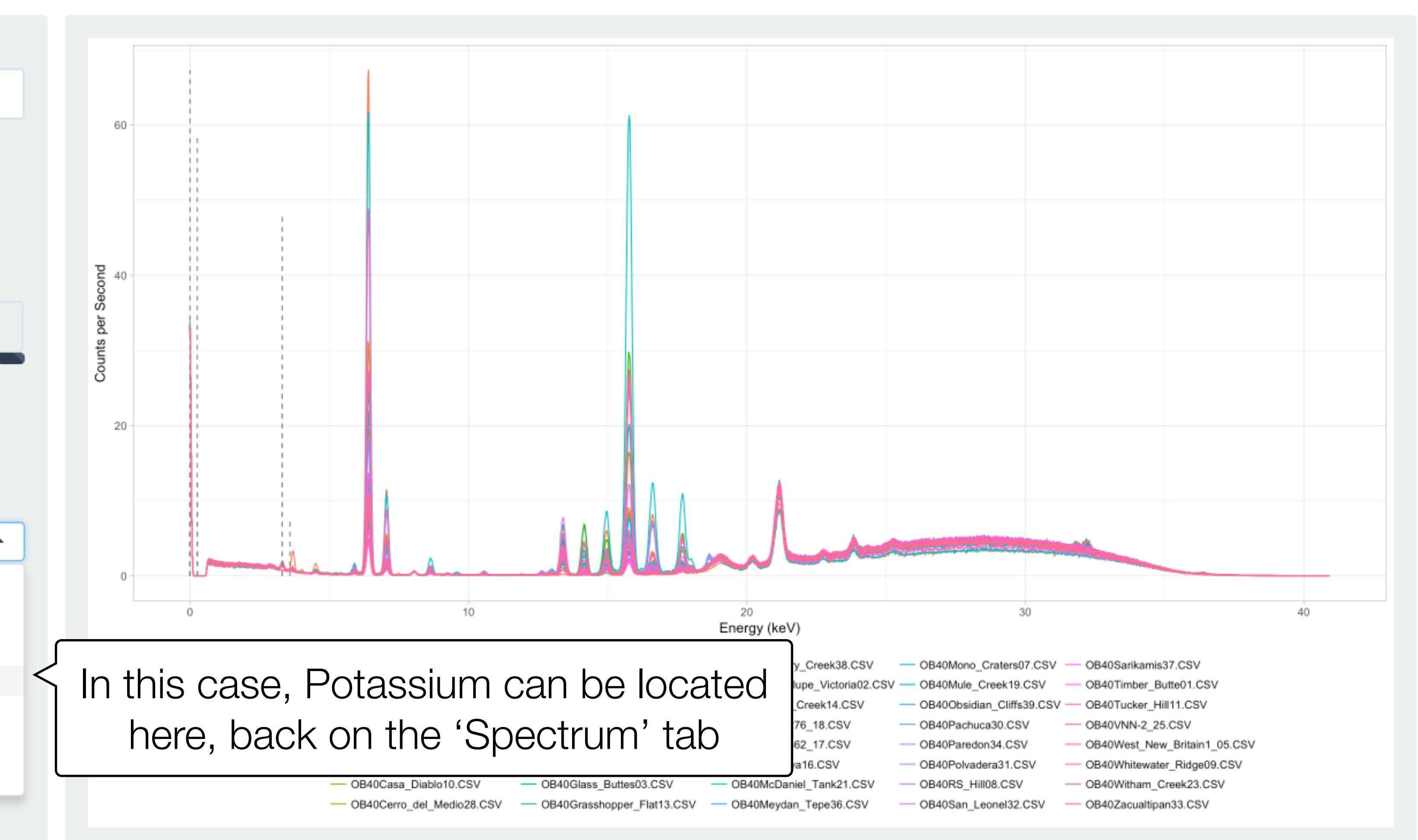
Calibration Name
Obsidian

Process Data Plot Spectrum 

Choose Spectra
Browse... 40 files Upload complete

Spectra Net

Element:
(K) Potassium (S) Sulfur (Cl) Chlorine (Ar) Argon
(K) Potassium (Ca) Calcium (Sc) Scandium (Ti) Titanium
Use Cal File



X-Ray Fluorescence Calibration

Calibration Name

Obsidian

Process Data **Plot Spectrum** **Plot**

Choose Spectra

Browse... 40 files **Upload complete**

Spectra Net

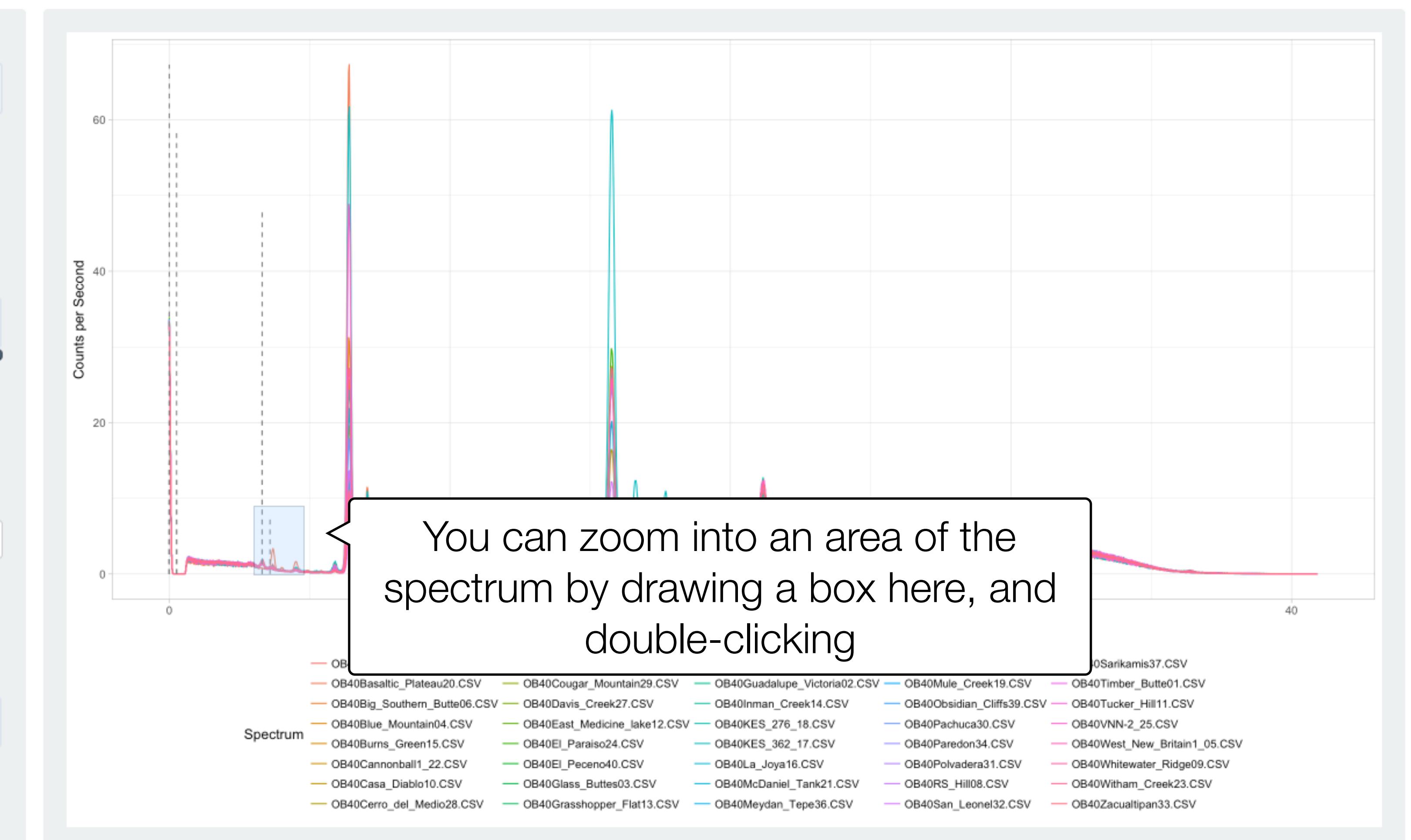
Element:

(K) Potassium

Load Cal File

Browse... No file selected

Use Cal File



X-Ray Fluorescence Calibration

Calibration Name
Obsidian

Process Data Plot Spectrum 

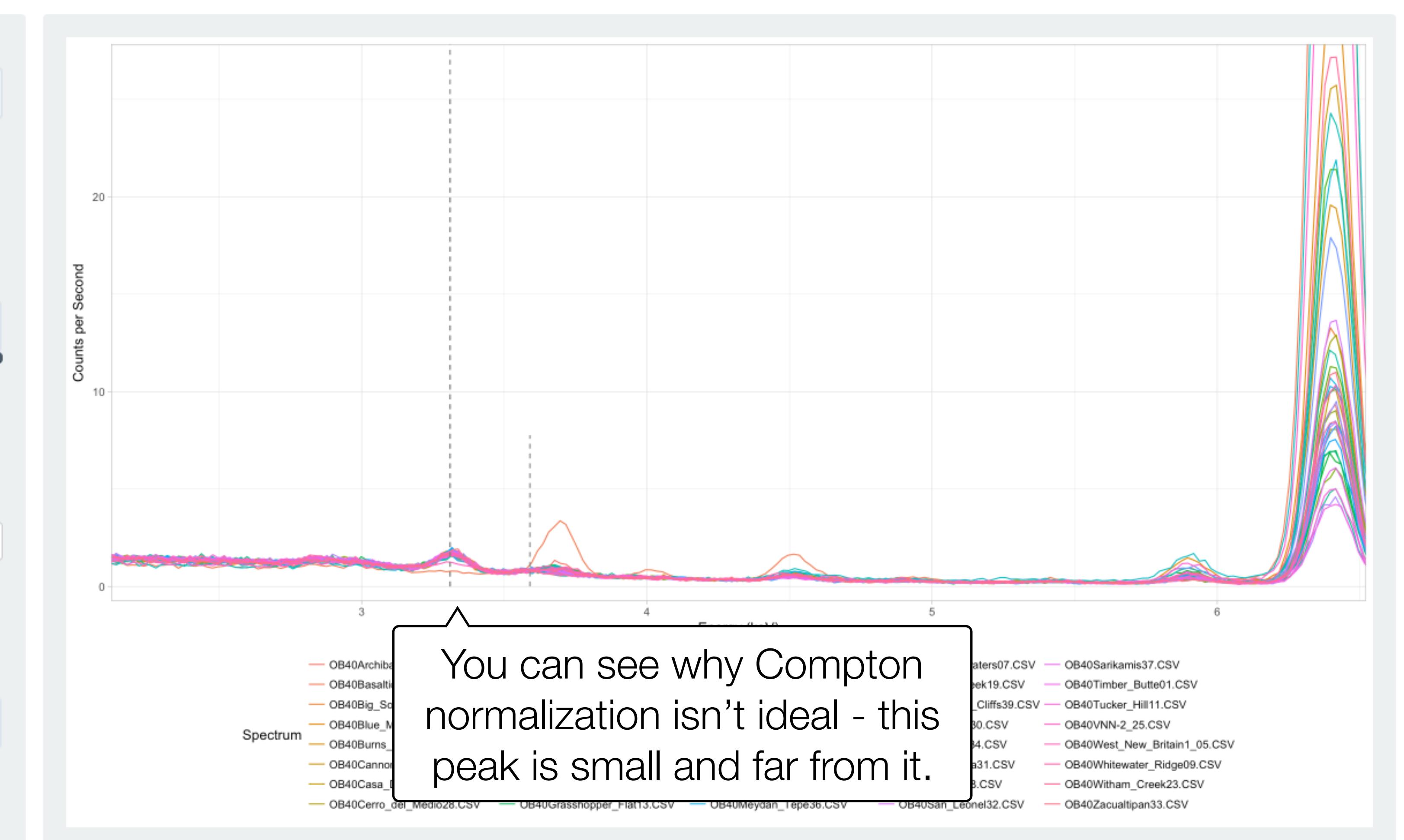
Choose Spectra
Browse... 40 files Upload complete

Spectra Net

Element: (K) Potassium

Load Cal File
Browse... No file selected

Use Cal File



Plot **Update** **Save**

Element
K.K.alpha

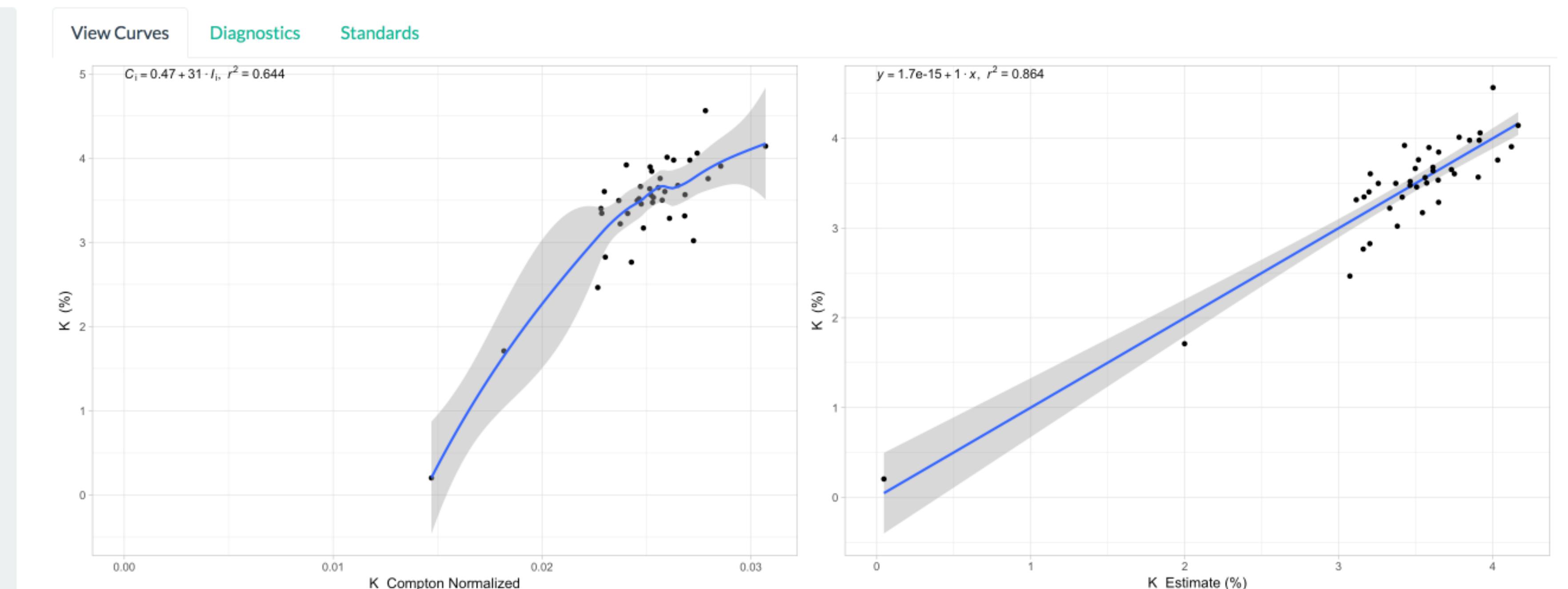
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
3

Max
3.2

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



Here, when the normalization is changed to something closer to the peak, the model improves

Plot **Update** **Save**

Element: K.Kalpha

Calibration Curve: Lucas-Tooth

Normalization: Compton

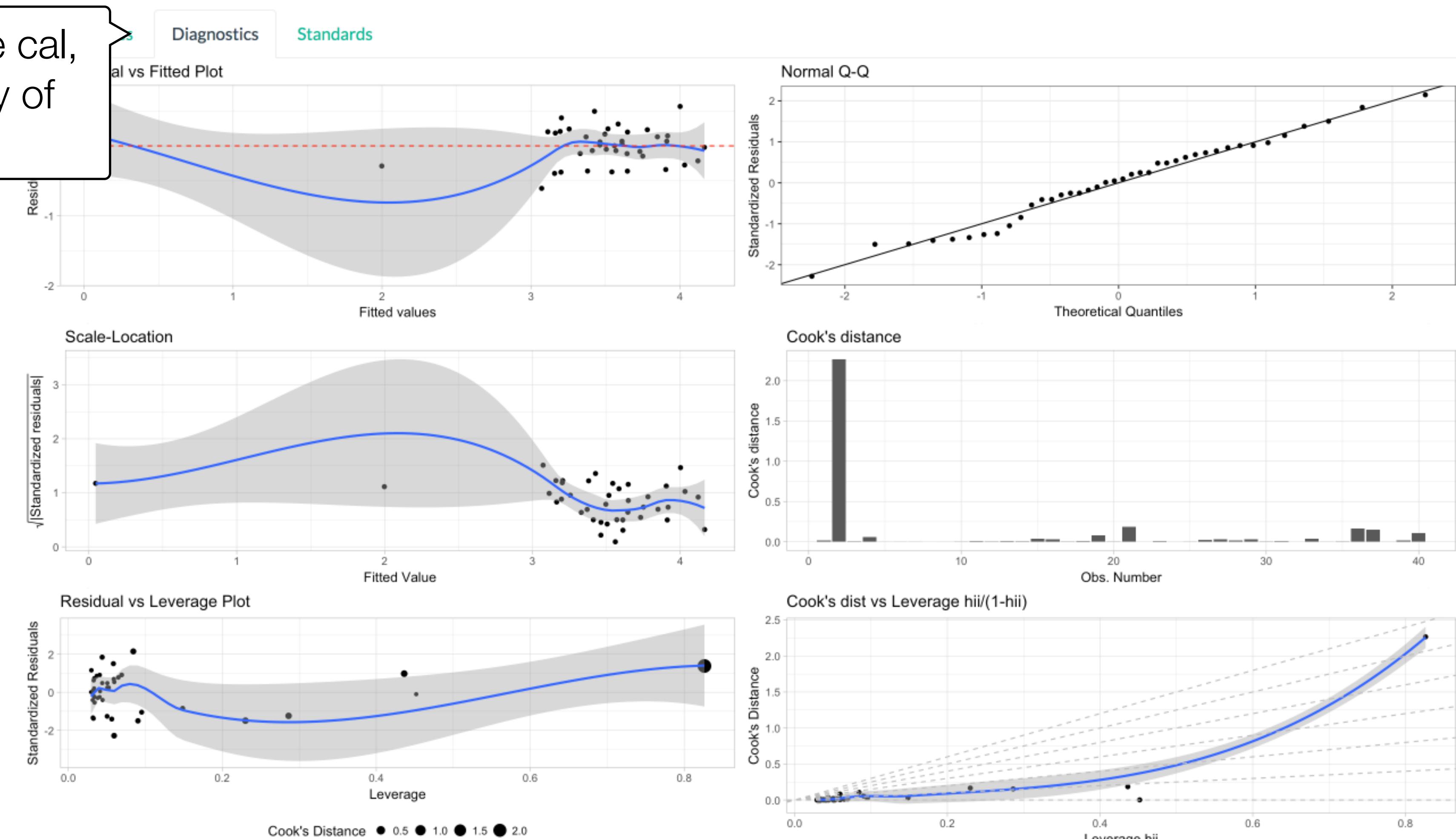
Min: 3

Max: 3.2

Intercept: K.Kalpha

- K.Kalpha
- Ca.Kalpha
- Ti.Kalpha
- Mn.Kalpha
- Fe.Kalpha
- Co.Kalpha
- Cu.Kalpha
- Zn.Kalpha
- Ga.Kalpha
- Al.Kalpha

To check the quality of the cal, we can evaluate a variety of diagnostic plots



[Plot](#)
[Update](#)
[Save](#)

We can also evaluate the product data of the calibration in absolute and relative terms

[Standards](#)
[Search:](#)

Element		Concentration	Prediction	Difference	Relative
K.Kalpha	5.CSV	3.1712	3.543092800145	-0.371892800145003	-0.117271947573475
	2 OB40Basaltic_Plateau20.CSV	0.2036	0.0457670629668963	0.157832937033104	0.775210889160627
Calibration Curve	3 OB40Big_Southern_Butte06.CSV	4.0605	3.91723052438532	0.143269475614676	0.0352837028973466
<input type="radio"/> Linear	4 OB40Blue_Mountain04.CSV	2.7656	3.15848555757976	-0.392885557579755	-0.142061598777754
<input type="radio"/> Non-Linear	5 OB40Burns_Green15.CSV	3.5022	3.57038916259993	-0.068189162599928	-0.0194703793615236
<input checked="" type="radio"/> Lucas-Tooth	6 OB40Cannonball1_22.CSV	3.6776	3.61095241204926	0.0666475879507411	0.0181225766670495
Normalization	7 OB40Casa_Diablo10.CSV	3.9778	3.91174367817721	0.0660563218227885	0.0166062451160914
<input type="radio"/> Time	8 OB40Cerro_del_Medio28.CSV	3.6373	3.61179432075415	0.0255056792458457	0.00701225613665237
<input type="radio"/> Total Counts	9 OB40Chickahominy26.CSV	3.519	3.46324429496596	0.0557557050340427	0.0158441901205009
<input checked="" type="radio"/> Compton	10 OB40Cougar_Mountain29.CSV	3.2214	3.33031419391722	-0.108914193917224	-0.0338095840060916

Showing 1 to 10 of 40 entries

 Previous [1](#) [2](#) [3](#) [4](#) Next

Min

3

Max

3.2

Intercept

- K.Kalpha
- Ca.Kalpha
- Ti.Kalpha
- Mn.Kalpha
- Fe.Kalpha
- Co.Kalpha
- Cu.Kalpha
- Zn.Kalpha
- Ga.Kalpha
- Al.Kalpha

Plot Update Save

Element
Ca.K.alpha

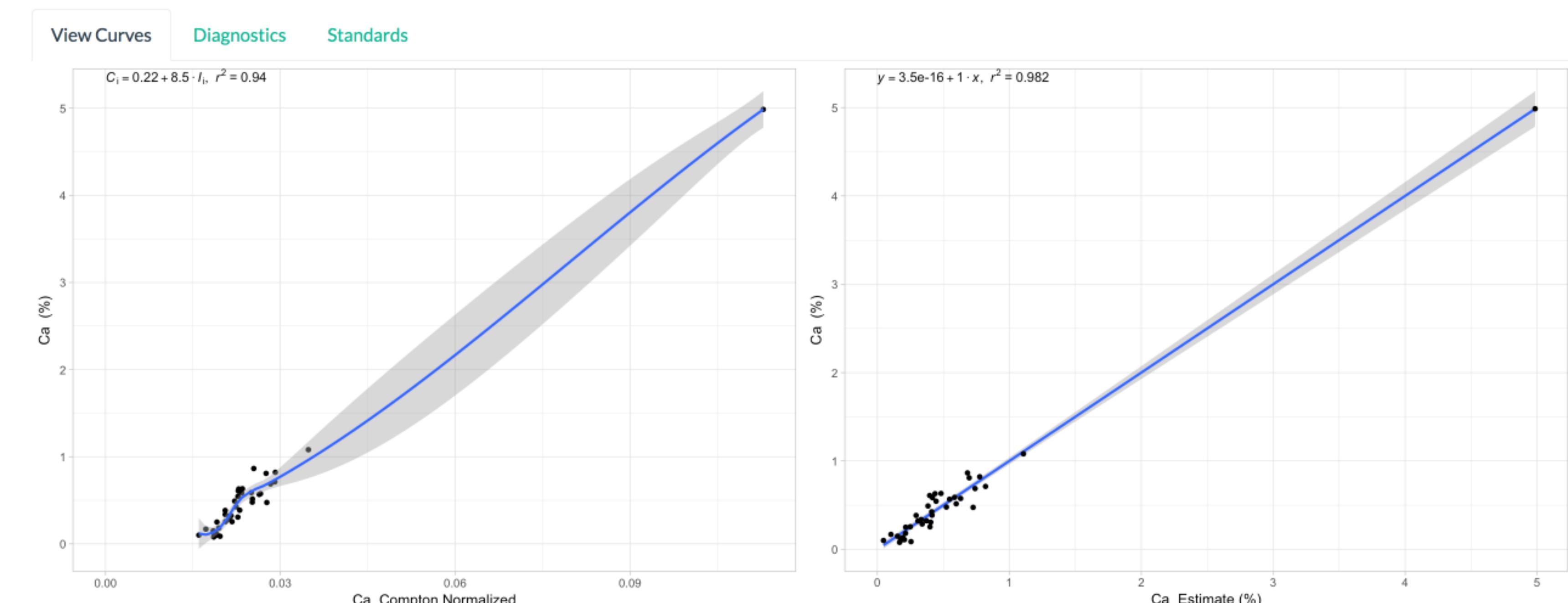
Calibration Curve
 No

 Total Counts
 Compton

Min
3

Max
3.2

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



To see why this is important, we can look at another element

One point has a concentration much higher than all the others, is this ok for the model?

Plot **Update** **Save**

Element
Ca.K.alpha

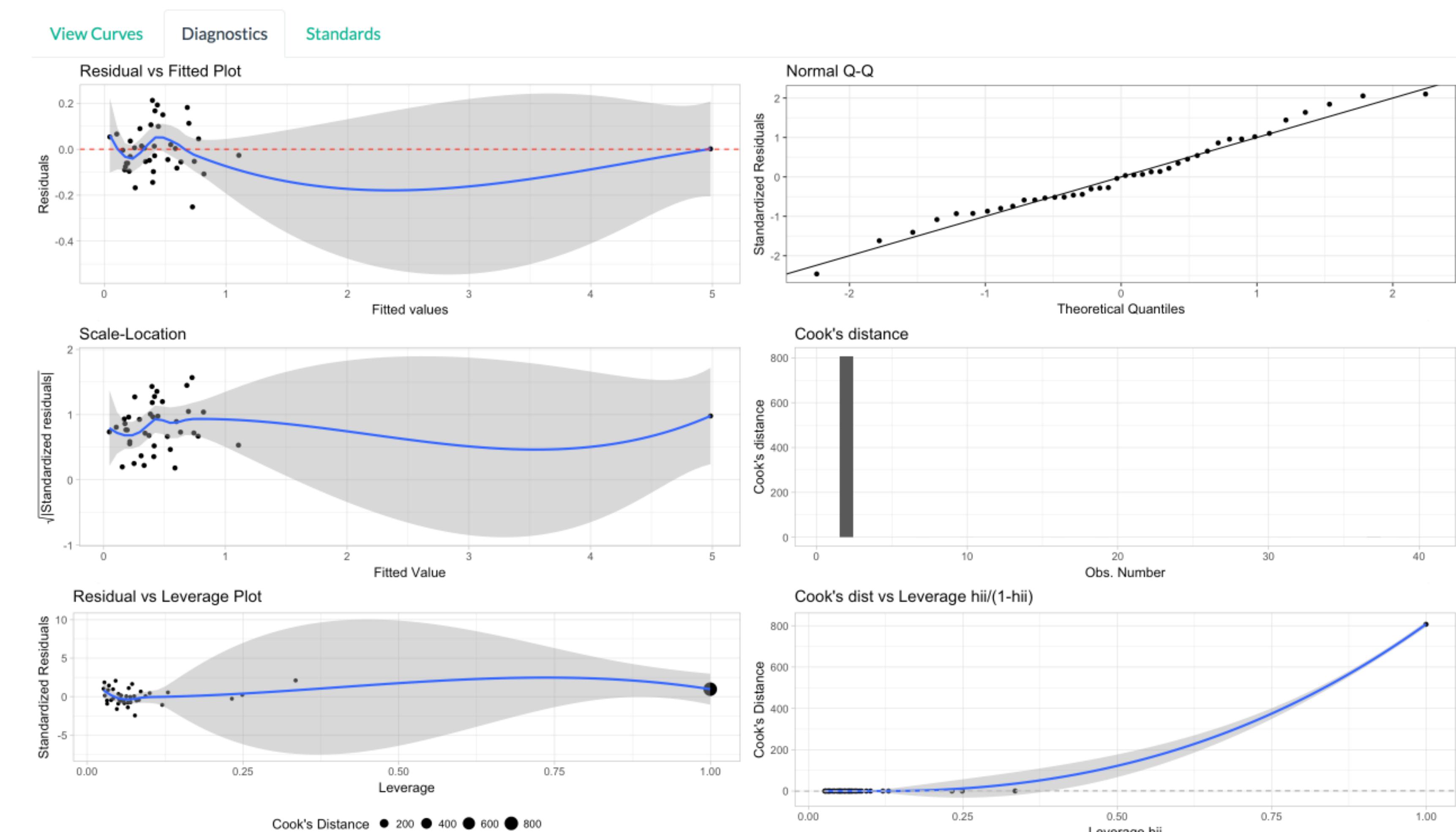
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
3

Max
3.2

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



Diagnostic plots indicate that while the Q-Q is good, that point has a large Cook's Distance

Enter Values

If we return to 'Add Concentrations', we can put in '999' in the cell with the high value to delete it. Then hit 'Enter Values'

Enter Concentrations

	Spectrum	K.K.alpha	Ca.K.alpha	Ti.K.alpha	Mn.K.alpha	Fe.K.alpha	Co.K.alpha	Cu.K.alpha	Zn.K.alpha	Ga.K.alpha	As.K.alpha	Rb.K.alpha	Sr.K.alpha	Y.K
1	OB40Archibarca35.CSV	3.17	1.08	0.08	0.05	0.87	0.00	0.00	0.00	0.00	0.00	0.01	0.03	
2	OB40Basaltic_Plateau20.CSV	0.20	999	0.78	0.11	6.85	0.00	0.00	0.01	0.00	0.00	0.00	0.03	0.03
3	OB40Big_Southern_Butte06.CSV	4.06	0.32	0.05	0.03	1.17	0.00	0.00	0.03	0.00	0.00	0.03	0.00	
4	OB40Blue_Mountain04.CSV	2.77	0.10	0.12	0.16	2.74	0.00	0.00	0.02	0.00	0.00	0.01	0.01	0.00
5	OB40Burns_Green15.CSV	3.50	0.13	0.10	0.05	1.72	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00
6	OB40Cannonball1_22.CSV	3.68	0.18	0.10	0.05	2.33	0.00	0.00	0.02	0.00	0.00	0.00	0.03	0.00
7	OB40Casa_Diablo10.CSV	3.98	0.61	0.11	0.03	0.93	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
8	OB40Cerro_del_Medio28.CSV	3.64	0.25	0.05	0.04	0.71	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.00
9	OB40Chickahominy26.CSV	3.52	0.42	0.11	0.04	1.18	0.00	0.00	0.01	0.00	0.00	0.01	0.00	0.00
10	OB40Cougar_Mountain29.CSV	3.22	0.47	0.03	0.03	0.81	0.00	0.00	0.01	0.00	0.00	0.01	0.01	0.00
11	OB40Davis_Creek27.CSV	4.01	0.58	0.05	0.04	0.53	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
12	OB40East_Medicine_lake12.CSV	3.66	0.63	0.14	0.03	1.04	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
13	OB40El_Paraiso24.CSV	3.76	0.11	0.08	0.02	1.94	0.00	0.00	0.02	0.00	0.00	0.00	0.02	0.00
14	OB40El_Peceno40.CSV	3.85	0.81	0.05	0.09	0.61	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.03
15	OB40Glass_Buttes03.CSV	3.60	0.52	0.06	0.03	0.62	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
16	OB40Grasshopper_Flat13.CSV	3.76	0.58	0.13	0.03	0.92	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
17	OB40Gregory_Creek38.CSV	3.60	0.82	0.02	0.07	0.65	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
18	OB40Guadalupe_Victoria02.CSV	3.40	0.34	0.06	0.05	0.43	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
19	OB40Inman_Creek14.CSV	2.46	0.71	0.05	0.05	1.08	0.00	0.00	0.01	0.00	0.00	0.01	0.01	0.01
20	OB40KES_276_18.CSV	4.14	0.57	0.31	0.11	2.34	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.01
21	OB40KES_362_17.CSV	3.31	0.15	0.16	0.18	5.37	0.00	0.00	0.06	0.00	0.00	0.00	0.04	0.00
22	OB40La_Joya16.CSV	3.48	0.12	0.09	0.06	1.89	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.00
23	OB40McDaniel_Tank21.CSV	3.90	0.69	0.16	0.06	1.00	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.02
24	OB40Meydan_Tepe36.CSV	3.46	0.29	0.05	0.05	0.93	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.00
25	OB40Mono_Craters07.CSV	3.53	0.38	0.04	0.04	0.78	0.00	0.00	0.00	0.00	0.00	0.00	0.02	0.00
26	OB40Mule_Creek19.CSV	3.57	0.39	0.04	0.04	0.66	0.00	0.00	0.00	0.00	0.00	0.00	0.02	0.00
27	OB40Obsidian_Cliffs39.CSV	2.83	0.59	0.06	0.03	0.73	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.01
28	OB40Pachuca30.CSV	3.29	0.08	0.11	0.11	1.63	0.00	0.00	0.02	0.00	0.00	0.00	0.02	0.00
29	OB40Paredon34.CSV	3.91	0.26	0.08	0.04	0.85	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.00
30	OB40Polvadera31.CSV	3.98	0.31	0.04	0.04	0.38	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.00
31	OB40RS_Hill08.CSV	3.35	0.25	0.02	0.04	0.72	0.00	0.00	0.01	0.00	0.00	0.00	0.04	0.00
32	OB40San_Leonel32.CSV	3.56	0.17	0.06	0.03	1.24	0.00	0.00	0.01	0.00	0.00	0.00	0.01	0.00
33	OB40Sarikamis37.CSV	3.92	0.32	0.05	0.04	0.55	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.00
34	OB40Timber_Butte01.CSV	3.65	0.49	0.03	0.08	0.37	0.00	0.00	0.01	0.00	0.00	0.00	0.02	0.00
35	OB40Tucker_Hill11.CSV	3.50	0.54	0.03	0.05	0.47	0.00	0.00	0.00	0.00	0.00	0.00	0.01	0.00
36	OB40VNN-2_25.CSV	3.02	0.10	0.07	0.12	4.38	0.00	0.00	0.03	0.00	0.00	0.00	0.02	0.00
37	OB40West_New_Britain1_05.CSV	1.71	0.86	0.11	0.06	0.86	0.00	0.00	0.01	0.00	0.00	0.00	0.02	

Plot Update Save

Element
Ca.K.alpha

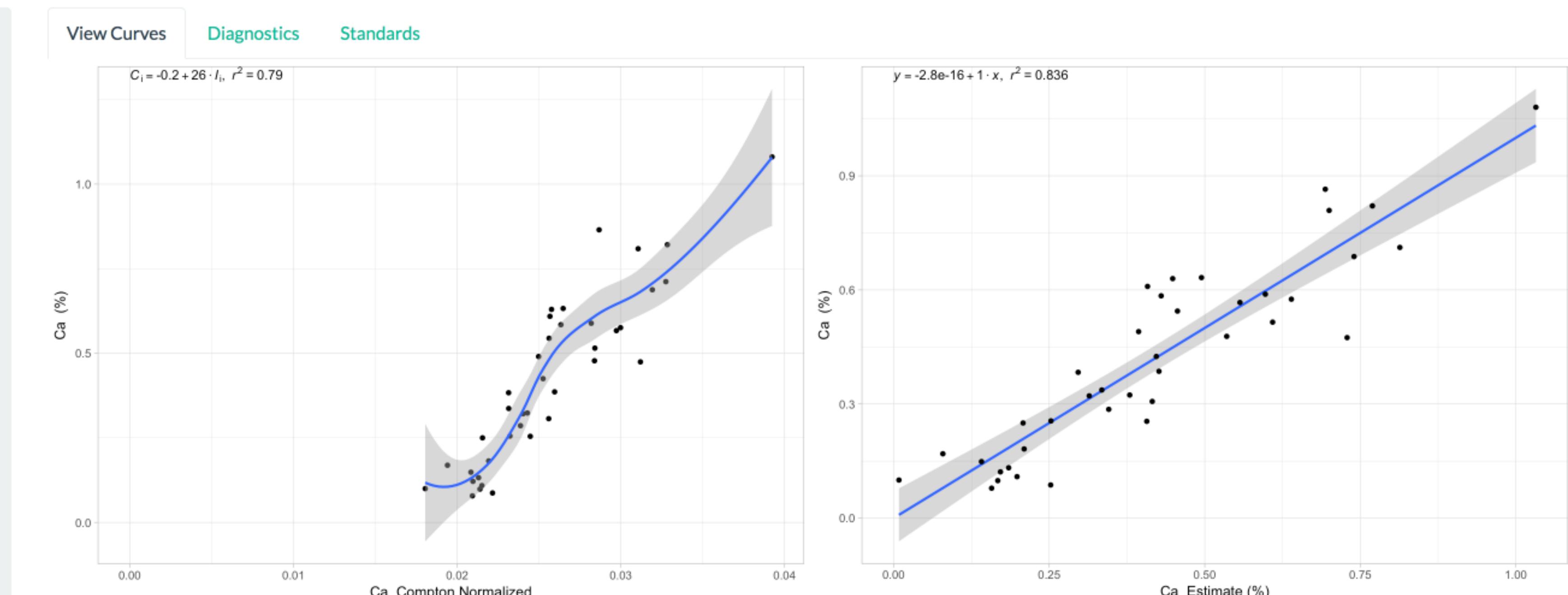
Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
3

Max
3.2

Intercept
 K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



Click 'Plot' in the 'Cal Curves' window, and you can now see the plot without the data point

[Plot](#)
[Update](#)
[Save](#)

Element

Calibration Curve

- Linear
- Non-Linear
- Lucas-Tooth

Normalization

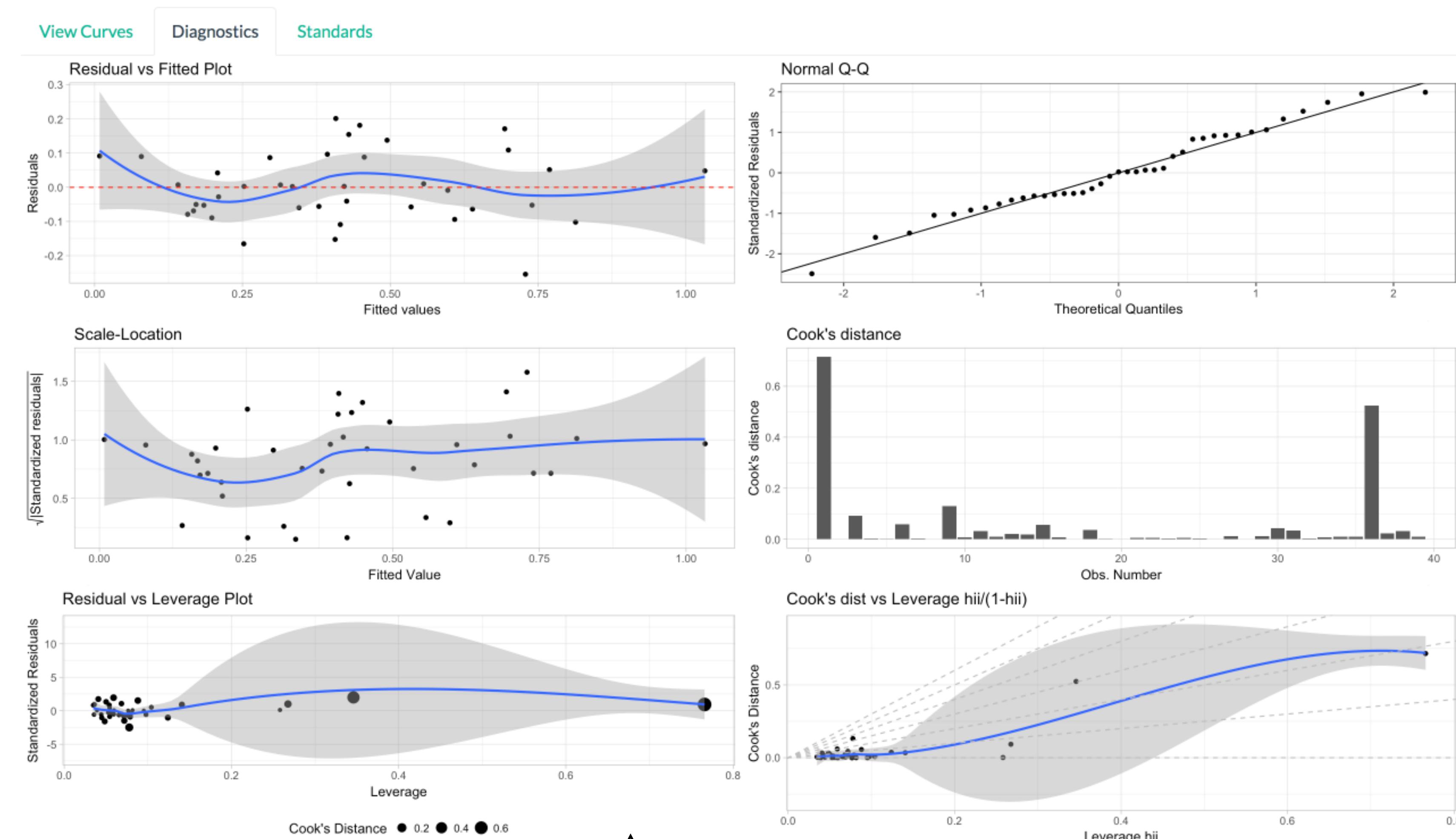
- Time
- Total Counts
- Compton

Min

Max

Intercept

- K.K.alpha
- Ca.K.alpha
- Ti.K.alpha
- Mn.K.alpha
- Fe.K.alpha
- Co.K.alpha
- Cu.K.alpha
- Zn.K.alpha
- Ga.K.alpha
- Al.K.alpha



Here, we see that there is a more even spread of data, creating a calibration less leveraged on one value

Plot Update Save 

Element
U.L.alpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

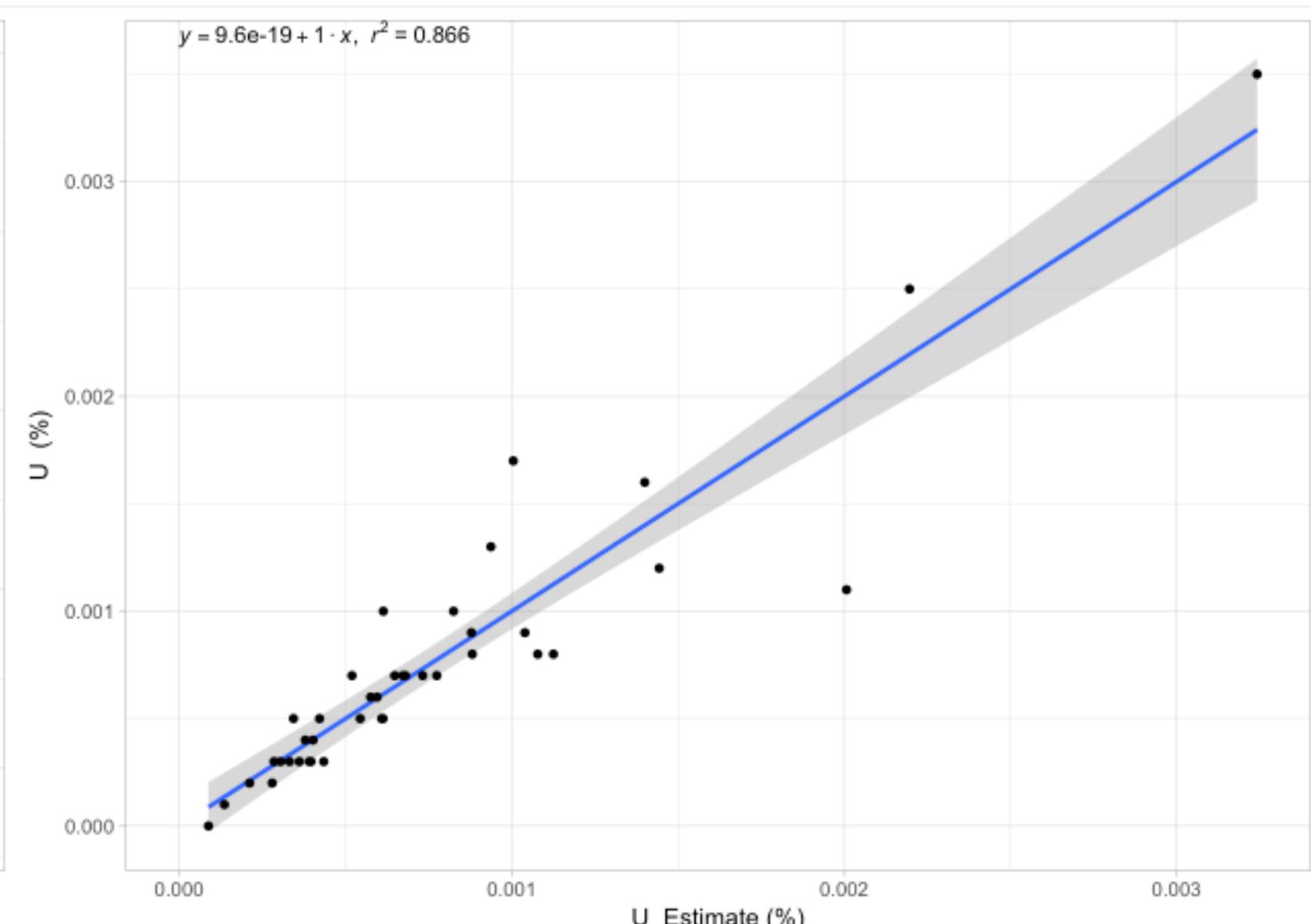
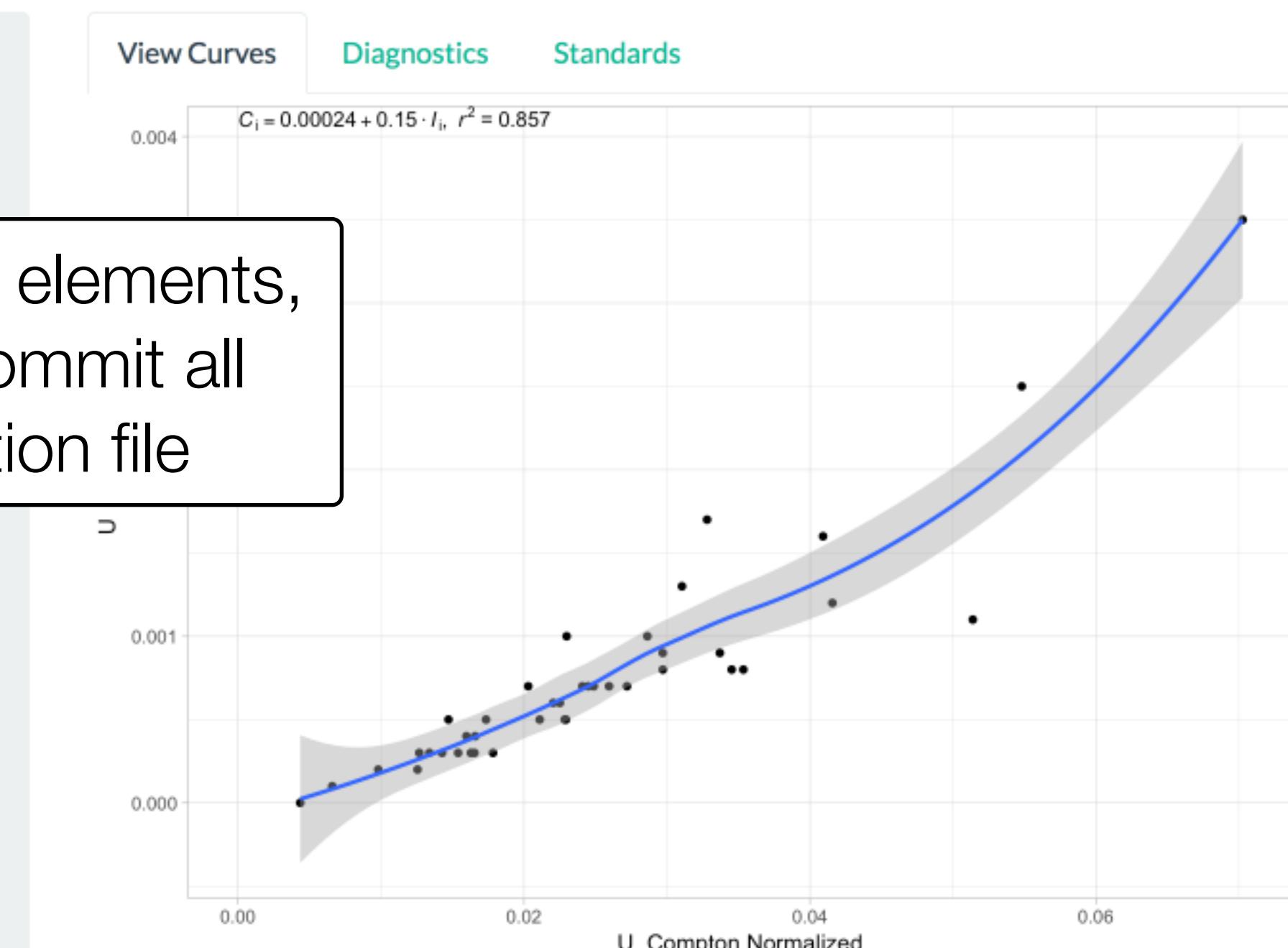
Normalization
 Time
 Total Counts
 Compton

Min
19.5

Max
22

Intercept

K.K.alpha
 Ca.K.alpha
 Ti.K.alpha
 Mn.K.alpha
 Fe.K.alpha
 Co.K.alpha
 Cu.K.alpha
 Zn.K.alpha
 Ga.K.alpha
 Al.K.alpha



When you are done with all elements, you can click 'Save' to commit all elements to the calibration file

X-Ray Fluorescence Calibration

Calibration Name

ObsidianT3S1716

Process Data

Plot Spectrum

Plot

Choose Spectra

Browse...

40 files

Upload complete

Spectra

Net

Element:

(Fe) Iron

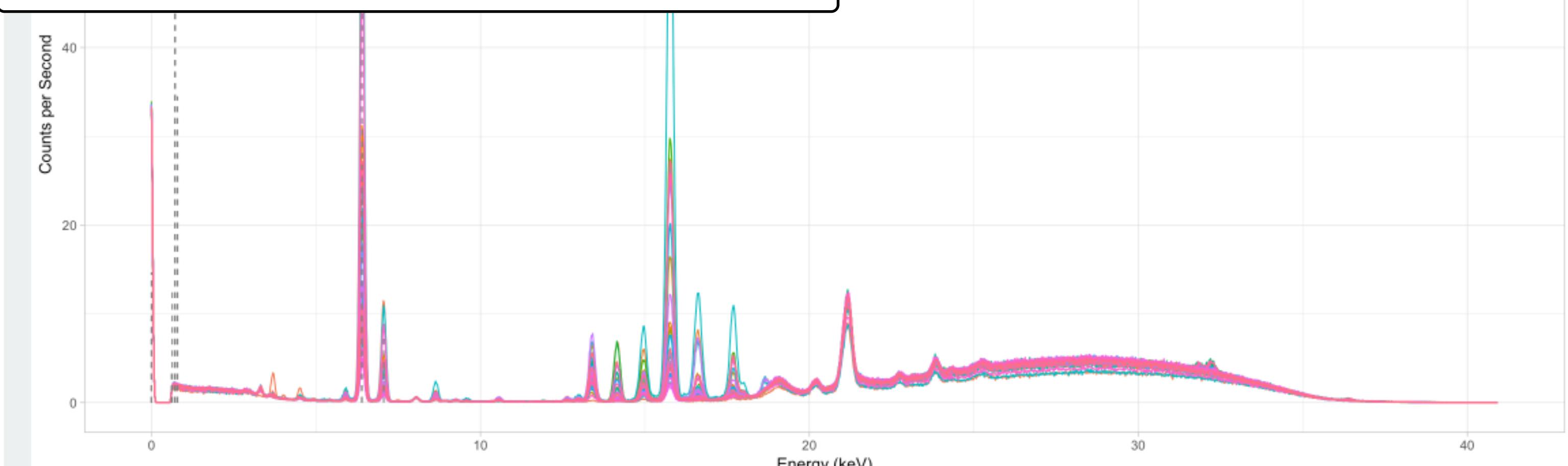
Load Cal File

Browse...

No file selected

Use Cal File

Name your file something diagnostic -
like the name of the material + the
instrument's serial number



Spectrum

- | | | | | |
|------------------------------|------------------------------|------------------------------|---------------------------|------------------------------|
| OB40Archibarca35.CSV | OB40Chickahominy26.CSV | OB40Gregory_Creek38.CSV | OB40Mono_Craters07.CSV | OB40Sarikamis37.CSV |
| OB40Basaltic_Plateau20.CSV | OB40Cougar_Mountain29.CSV | OB40Guadalupe_Victoria02.CSV | OB40Mule_Creek19.CSV | OB40Timber_Butte01.CSV |
| OB40Big_Southern_Butte06.CSV | OB40Davis_Creek27.CSV | OB40Inman_Creek14.CSV | OB40Obsidian_Cliffs39.CSV | OB40Tucker_Hill11.CSV |
| OB40Blue_Mountain04.CSV | OB40East_Medicine_lake12.CSV | OB40KES_276_18.CSV | OB40Pachuca30.CSV | OB40VNN-2_25.CSV |
| OB40Burns_Green15.CSV | OB40El_Paraiso24.CSV | OB40KES_362_17.CSV | OB40Paredon34.CSV | OB40West_New_Britain1_05.CSV |
| OB40Cannonball1_22.CSV | OB40El_Peceno40.CSV | OB40La_Joya16.CSV | OB40Polvadera31.CSV | OB40Whitewater_Ridge09.CSV |
| OB40Casa_Diablo10.CSV | OB40Glass_Buttes03.CSV | OB40McDaniel_Tank21.CSV | OB40RS_Hill08.CSV | OB40Witham_Creek23.CSV |
| OB40Cerro_del_Medio28.CSV | OB40Grasshopper_Flat13.CSV | OB40Meydan_Tepe36.CSV | OB40San_Leonel32.CSV | OB40Zacualtipan33.CSV |

Plot Update Save 

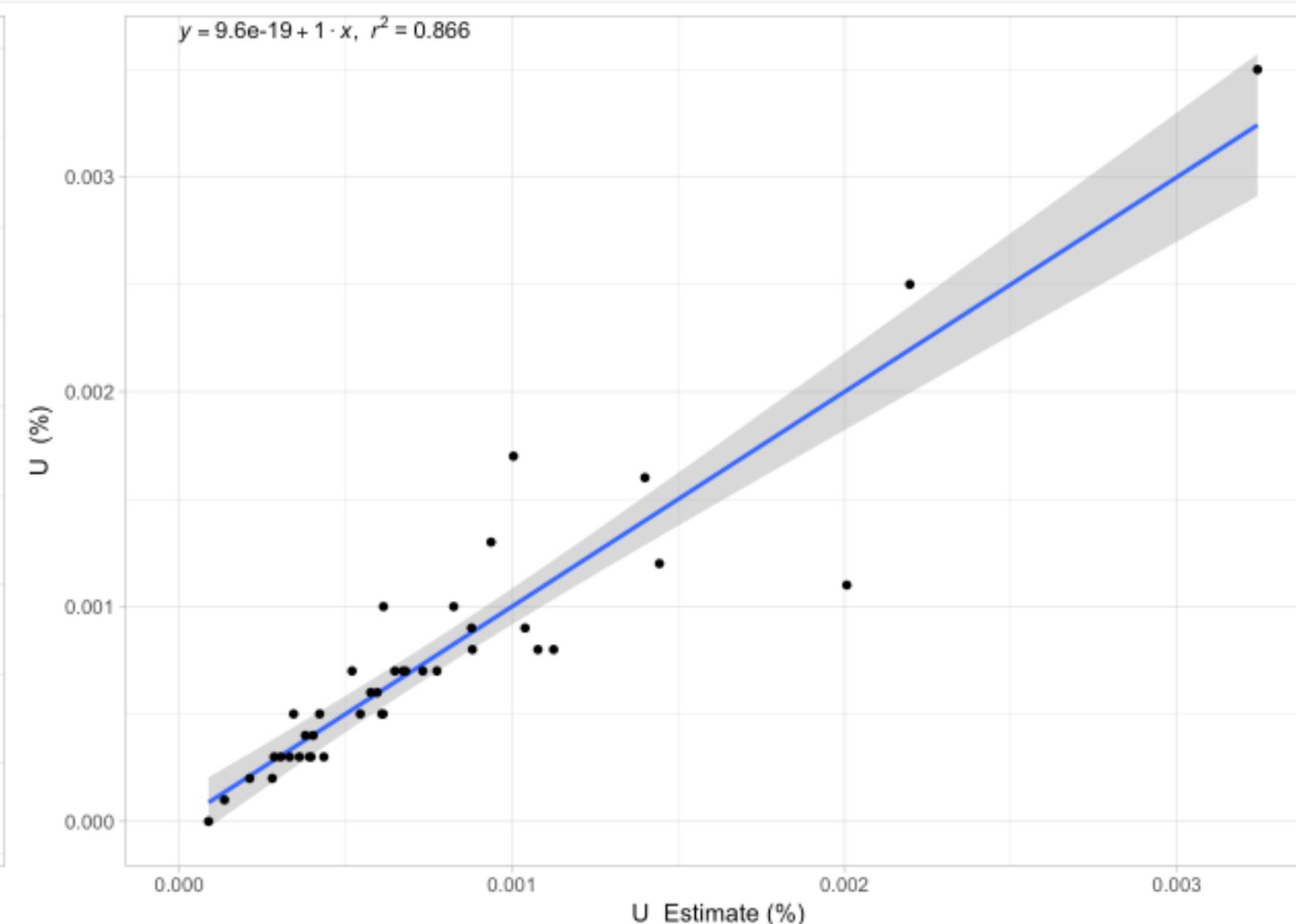
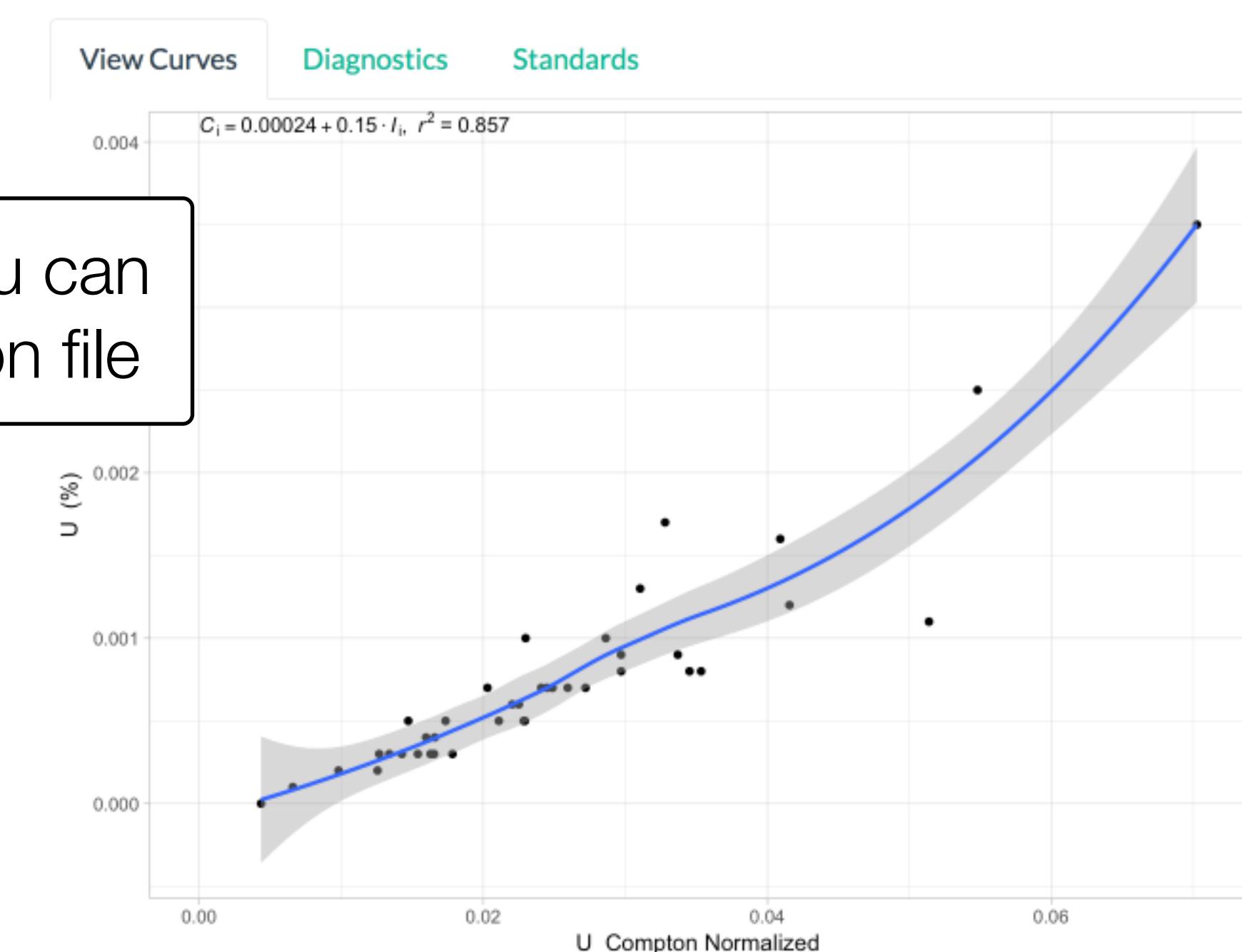
Element
U.L.alpha

Calibration Curve
 Linear
 Non-Linear
 Lucas-Tooth

Normalization
 Time
 Total Counts
 Compton

Min
19.5

Max
22



Once you are ready, you can download the calibration file

The calibration will be saved with the name you gave it as a .quant file

X-Ray Fluorescence Calibration

Calibration Name

myCalibration

Process Data

Plot Spectrum

Plot

Choose Spectra

Browse...

No file selected

Spectra

Net

Elem

(Fe) In the future, you can edit this .quant
file by loading it on the 'Spectrum'
page, and clicking 'Use Cal File'

Load Cal File

Browse...

No file selected

Use Cal File

If you edit a cal file, you will need to
follow all steps to save changes



CloudCal Spectrum Counts Add Concentrations Cal Curves **Apply Calibration**

Quantify

You can test the cal, or apply it to
large data sets, on 'Apply Calibration'

Choose Spectra

Browse...

No file selected

Results

Load Cal File

Browse...

No file selected

CloudCal

127.0.0.1:3315

CloudCal Spectrum Counts Add Concentrations Cal Curves Apply Cal

Quantify

Choose Spectra

Browse... No file selected

Load Cal File

Browse... No file selected

OB40Tucker_Hill11.CSV
OB40Tucker_Hill11.pdz
OB40VNN-2_25.CSV
OB40VNN-2_25.pdz
OB40West_...itain1_05.CSV
OB40West_...ritain1_05.pdz
OB40Whitew...Ridge09.CSV
OB40Whitew...Ridge09.pdz
OB40Witham_Creek23.CSV
OB40Witham_Creek23.pdz
OB40Zacualtipan33.CSV
OB40Zacualtipan33.pdz
Validation
OB40KES_362_17-00.CSV
OB40KES_362_17-00.pdz
OB40KES_362_17-01.CSV
OB40KES_362_17-01.pdz
OB40KES_362_17-02.CSV
OB40KES_362_17-02.pdz
OB40KES_362_17-03.CSV
OB40KES_362_17-03.pdz
OB40KES_362_17-04.CSV
OB40KES_362_17-04.pdz
OB40KES_362_17-05.CSV
OB40KES_362_17-05.pdz
OB40KES_362_17-06.CSV

Format: Custom Files

Options Cancel Open

ObsidianT3S1716.quant

Show All

Lee

The screenshot shows the CloudCal software interface. At the top, there's a navigation bar with tabs: CloudCal, Spectrum, Counts, Add Concentrations, Cal Curves, and Apply Cal. Below this is a main area with several sections: 'Quantify' (button), 'Choose Spectra' (with 'Browse...' button and 'No file selected' text), 'Load Cal File' (with 'Browse...' button and 'No file selected' text), and a large central area for file selection. A modal window titled 'Validation' is open, showing a list of files under 'Favorites' and 'Validation'. The 'Validation' section contains 14 entries, all ending in .CSV or .pdz. The modal also has a 'Format:' dropdown set to 'Custom Files', 'Options' and 'Cancel' buttons, and an 'Open' button. A callout bubble points from the 'Choose Spectra' section towards the 'Validation' modal. The bottom of the screen shows a file list at the bottom left and a toolbar with various icons at the bottom right.

Quantify

Choose Spectra

Browse... 30 files

Upload complete

Validation Counts

The status of the upload will be indicated here

Load Cal File

Browse... No file selected

CloudCal

127.0.0.1:3315

CloudCal Spectrum Counts Add Concentrations Cal Curves Apply C

Quantify

Choose Spectra

Browse... 30 files

Upload complete

Results

Load Cal File

Browse... No file selected

Browse for your .quant file here - you will need to have an existing file

ObsidianT3S1716.quant

Show All

Lee

A screenshot of a software application window titled "CloudCal". The window has a dark header bar with the title and a toolbar with various icons. The main interface is divided into several sections: "Choose Spectra" with a "Browse..." button showing "30 files" uploaded, "Results" (with a download icon), and "Load Cal File" with a "Browse..." button showing "No file selected". A callout box with a black border and white background points from the "Load Cal File" section to the "ObsidianT3S1716.quant" file in a file browser window. The file browser shows a list of files under the folder "Obsidian1716", with "ObsidianT3S1716.quant" highlighted. The file list includes various file types such as pdz, CSV, and SQL files, along with some image and document files.

CloudCal 127.0.0.1:3315 Lee

CloudCal Spectrum Counts Add Concentrations Cal Curves Apply Calibration

Quantify

When everything is ready, click 'Quantify'

Validation Counts

Choose Spectra

Browse... 30 files Upload complete

Results

Load Cal File

Browse... ObsidianT3S1716.quant Upload complete

Show All

ObsidianT3S1716.quant

Quantify

Choose Spectra

Browse... 30 files Upload complete

Results

Load Cal File

Browse... ObsidianT3S1716.quant Upload complete

Validation Counts

Show 10 entries Search:

	Spectrum	As.K.alpha	Ba.K.alpha	Ca.K.alpha	Co.K.alpha	Cu.K.alpha	Fe.K.alpha
1	OB40KES_362_17-00.CSV	0.000463815295623188	-0.064370668587769	0.306205801868475	0.00237023275057566	0.00131966273556775	5.7924834480857
2	OB40KES_362_17-01.CSV	0.000322096800981703	-0.0610441906197434	0.317104466145157	0.00300956958636639	0.000139274358499416	6.23039281488607
3	OB40KES_362_17-02.CSV	0.000315432693311329	-0.0610002951449382	-0.135324989048458	0.00187021927541898	0.000178072071125506	5.70154025119591
4	OB40KES_362_17-03.CSV	-0.000113820147964187	0.00685861778769903	0.371006235386069	0.00333625432297064	0.000322612915564326	5.64466483574534
5	OB40KES_362_17-04.CSV	0.000686491088709026	-0.0793083119274642	0.858100270567482	0.00346760366923405	0.000871850939981579	6.06797227540175
6	OB40KES_362_17-05.CSV	0.00036797545848524	-0.0649268528143068	0.313844891759042	0.00199153754250897	0.000407589231228397	6.14956098518842
7	OB40KES_362_17-06.CSV	0.000527937188744937	-0.0844006731489197	0.366748434920399	0.00278301669649239	0.000362639048515526	5.71349639773715
8	OB40KES_362_17-07.CSV	-0.000186530532410853	-0.0386374268963758	-0.108502135352386	-0.00051692485675101	0.00184082559843627	5.05393602605671
9	OB40KES_362_17-08.CSV	0.0000362648935217126	-0.0495280682863057	0.11805452462534	0.00193892752507277	0.0002435240599248	5.65518595019268
10	OB40KES_362_17-09.CSV	0.000643491416236847	-0.0506811695464275	0.328223673872163	0.00141975060266103	0.0019063462070858	5.6506440639408

Showing 1 to 10 of 30 entries

Previous 1 2 3 Next

A table of values will appear.
 Note that some values may be negative - this indicates that the data fall outside the range of the calibration, or are below the detection limit

Quantify

Choose Spectra

Browse... 30 files

Upload complete

Results

Load Cal File

Browse... ObsidianT3S1716.quant

Upload complete

Validation Counts

Show 10 entries Search:

	Spectrum	As.K.alpha	Ba.K.alpha	Ca.K.alpha	Co.K.alpha	Cu.K.alpha	Fe.K.alpha	Ga.K.alpha
1	OB40KES_362_17-00.CSV	0.232747050302078	6.92422474648682	0.16209169574609	0.511212271199207	0.195341274360673	21.9655028722586	0.112217327824
2	OB40KES_362_17-01.CSV	0.274309023570306	6.82447601064307	0.153779301092444	0.519524665852853	0.257684234263015	22.3063110530581	0.149623103765
3	OB40KES_362_17-02.CSV	0.224434655648432	6.95331812777458	0.108061130497393	0.498743679218739	0.228590852975255	22.1483755546388	0.108061130497
4	OB40KES_362_17-03.CSV	0.22027845832161	7.47699899095426	0.174560287726559	0.569399033774727	0.224434655648432	22.2231871065216	0.120529722477
5	OB40KES_362_17-04.CSV	0.224434655648432	6.67069670955063	0.228590852975255	0.536149455160144	0.228590852975255	21.8325045578003	0.120529722477
6	OB40KES_362_17-05.CSV	0.241059444955724	6.90344375985271	0.157935498419267	0.490431284565093	0.245215642282546	22.5972448659357	0.124685919804
7	OB40KES_362_17-06.CSV	0.253528036936192	6.69147769618474	0.178716485053381	0.561086639121081	0.224434655648432	23.0876761505008	0.128842117131
8	OB40KES_362_17-07.CSV	0.195341274360673	7.248408137979	0.124685919804685	0.403151140701814	0.174560287726559	21.7992549791857	0.157935498419
9	OB40KES_362_17-08.CSV	0.187028879707027	7.11540982352067	0.13299831445833	0.507056073872384	0.207809866341141	22.2689052771167	0.141310709111
10	OB40KES_362_17-09.CSV	0.228590852975255	7.00734869302328	0.166247893072913	0.490431284565093	0.199497471687495	22.5681514846479	0.13299831445

Showing 1 to 10 of 30 entries

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Click on 'Counts' to see the counts per second values for each element

Quantify

Choose Spectra

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Results

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Validation Counts

Show 10 entries Search:

	Spectrum	As.K.alpha	Ba.K.alpha	Ca.K.alpha	Co.K.alpha	Cu.K.alpha	Fe.K.alpha	Ga.K.alpha
1	OB40KES_362_17-00.CSV	0.232747050302078	6.92422474648682	0.16209169574609	0.511212271199207	0.195341274360673	21.9655028722586	0.112217327824
2	OB40KES_362_17-01.CSV	0.274309023570306	6.82447601064307	0.153779301092444	0.519524665852853	0.257684234263015	22.3063110530581	0.149623103765
3	OB40KES_362_17-02.CSV	0.224434655648432	6.95331812777458	0.108061130497393	0.498743679218739	0.228590852975255	22.1483755546388	0.108061130497
4	OB40KES_362_17-03.CSV	0.22027845832161	7.47699899095426	0.174560287726559	0.569399033774727	0.224434655648432	22.2231871065216	0.120529722477
5	OB40KES_362_17-04.CSV	0.224434655648432	6.67069670955063	0.228590852975255	0.536149455160144	0.228590852975255	21.8325045578003	0.120529722477
6	OB40KES_362_17-05.CSV	0.241059444955724	6.90344375985271	0.157935498419267	0.490431284565093	0.245215642282546	22.5972448659357	0.124685919804
7	OB40KES_362_17-06.CSV	0.253528036936192	6.69147769618474	0.178716485053381	0.561086639121081	0.224434655648432	23.0876761505008	0.128842117131
8	OB40KES_362_17-07.CSV	0.195341274360673	7.248408137979	0.124685919804685	0.403151140701814	0.174560287726559	21.7992549791857	0.157935498419
9	OB40KES_362_17-08.CSV	0.187028879707027	7.11540982352067	0.13299831445833	0.507056073872384	0.207809866341141	22.2689052771167	0.141310709111
10	OB40KES_362_17-09.CSV	0.228590852975255	7.00734869302328	0.166247893072913	0.490431284565093	0.199497471687495	22.5681514846479	0.13299831445

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