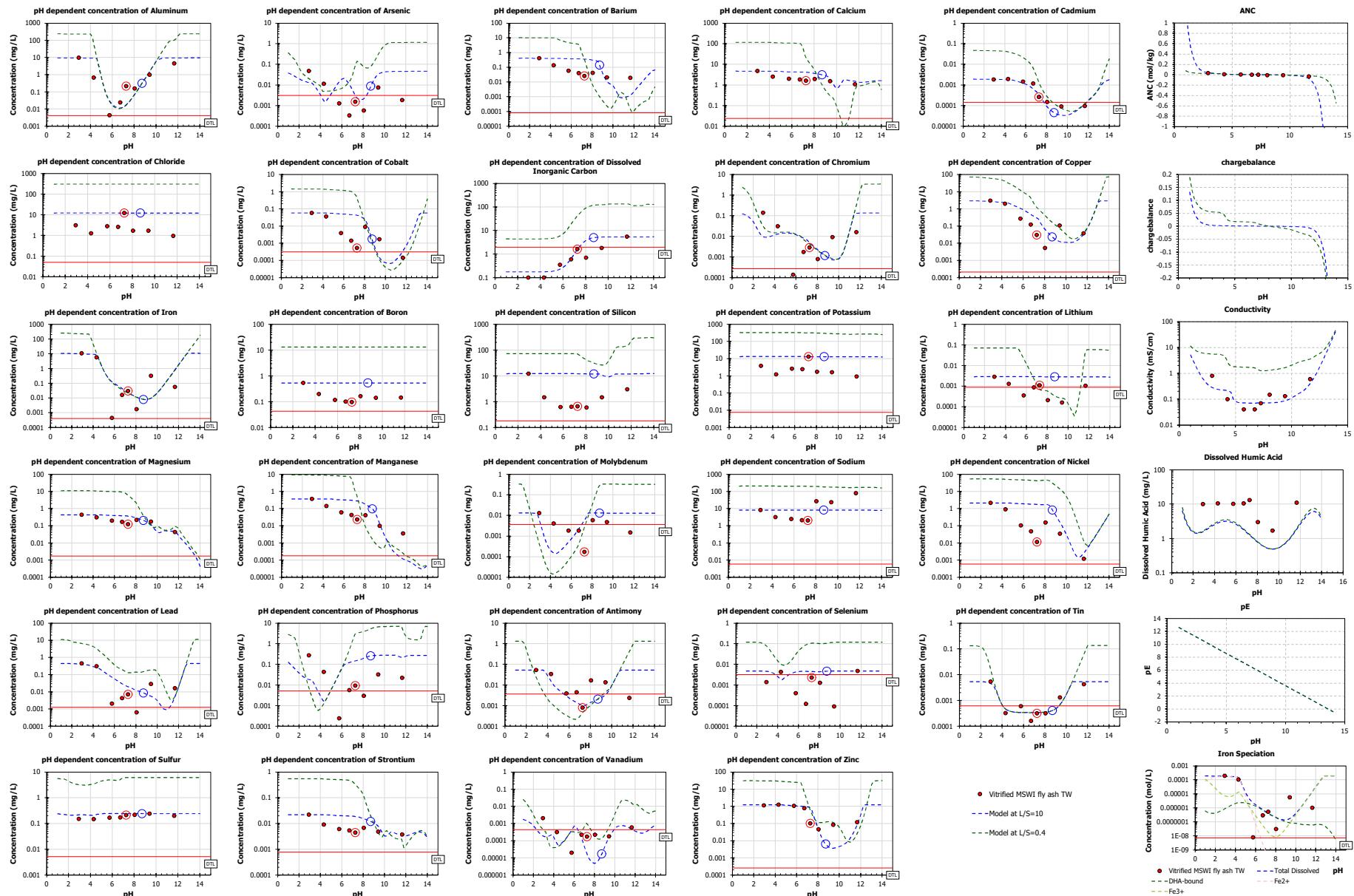


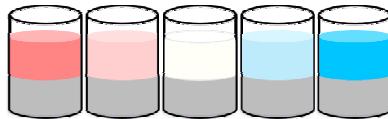
VITRIFIED MSWI FLY ASH

COMPARISON pH DEPENDENCE WITH MODEL



Object pH Dependent Leaching Test Model
Name Vitrified MSWI Fly ash TW

pH Dependent Leaching Test Scenario



Lab Test

Extra L/S Simulation

Lab Test

Model Parameters

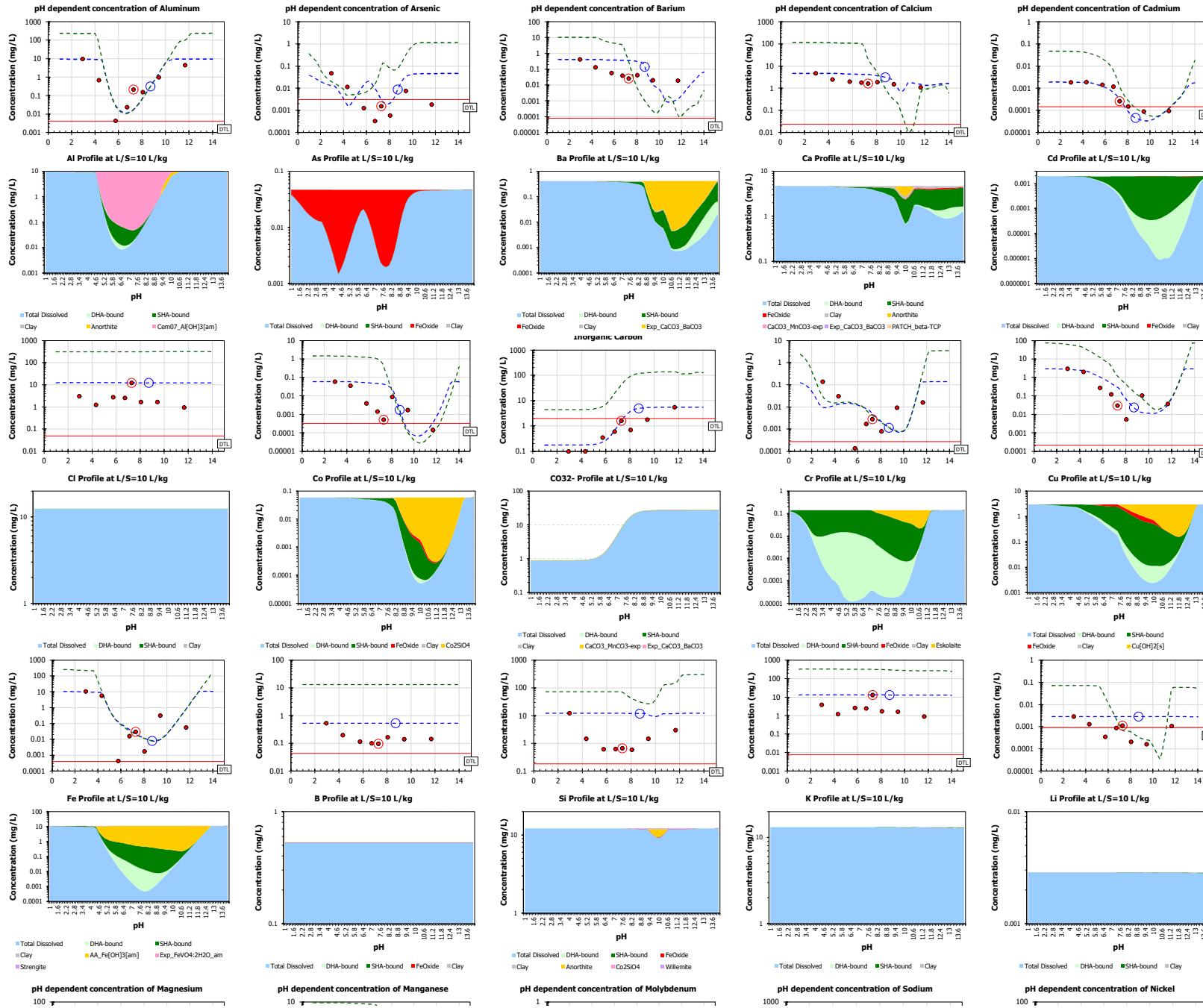
Entity	Unit	Available Content							
		Default	Entity	mg/kg	Entity	mg/kg	Entity	mg/kg	
c0		-3.484	Al	96.03	Fe	107.7	P	2.743	
c1		-3.047	As	0.4709	B	5.293	Sb	0.5287	
c2		1.221	Ba	4.159	Si	122.1	Se	0.04775	
c3		-0.2073	Br	7.990E-09	Hg	2.006E-08	Sn	0.05350	
c4		0.01534	Ca	47.32	K	132.1	S	2.430	
c5		-0.0004075	Cd	0.01920	Li	0.02880	Sr	0.2204	
Clay	mg/kg	1000	Cl	124.3	Mg	4.381	Th	2.320E-08	
Hydrous Ferric Oxide	mg/kg	60.00	Co	0.5856	Mn	3.662	U	2.380E-08	
L/S	L/kg	10.02	CO32-	275.0	Mo	0.1313	V	0.02030	
pE		6.330	Cr	1.394	Na	81.78	Zn	12.46	
pH		7.270	Cu	29.63	Ni	21.41			
Solid Humic Acid	mg/kg	300.0	F	1.900E-09	Pb	4.453			
Simulated Low L/S	L/kg	0.4000							

Minerals

Name	Log(K)	Reaction	Name	Log(K)	Reaction
AA_Fe[OH]3[am]	16.60	AA_Fe[OH]3[am] + 1 H2O -> 1 Fe[OH]4- + 1 H+	Eskolaite	139.5	Eskolaite + 5 H2O -> 2 CrO4-2 + 10 H+ + 6 e-
Anorthite	63.81	Anorthite + 8 H2O -> 2 Al[OH]4- + 1 Ca+2 + 4 H+ + 2 H2SiO4-2	Exp_CaCO3_BaCO3	21.30	Exp_CaCO3_BaCO3 -> 1 Ba+2 + 2 CO3-2 + 1 Ca+2
Arsenocrandallite-therm	95.56	Arsenocrandallite-therm + 6 H2O -> 3 Al[OH]4- + 2 AsO4-3 + 1 Ca+2 + 7 H+	Exp_FeVO4:2H2O_am	23.48	Exp_FeVO4:2H2O_am + 2 H2O -> 1 Fe[OH]4- + 1 VO2+
Ba[Scr]O4[96%SO4]	9.790	Ba[Scr]O4[96%SO4] -> 1 Ba+2 + 0.04 CrO4-2 + 0.96 SO4-2	Li-Albite[low2]	86.27	Li-Albite[low2] + 8 H2O -> 1 Al[OH]4- + 6 H+ + 3 H2SiO4-2 + 1 Li+
CaCO3_Li2CO3	21.30	CaCO3_Li2CO3 -> 2 CO3-2 + 1 Ca+2 + 2 Li+	Manganite	-25.27	Manganite + 3 H+ + 1 e- -> 2 H2O + 1 Mn+2
CaCO3_MgCO3-exp	18.02	CaCO3_MgCO3-exp -> 2 CO3-2 + 1 Ca+2 + 1 Mg+2	Ni[OH]2[s]	-10.80	Ni[OH]2[s] + 2 H+ -> 2 H2O + 1 Ni+2
CaCO3_MnCO3-exp	20.78	CaCO3_MnCO3-exp -> 2 CO3-2 + 1 Ca+2 + 1 Mn+2	PATCH_beta-TCP	28.93	PATCH_beta-TCP -> 3 Ca+2 + 2 PO4-3
CaCO3_SrCO3	19.85	CaCO3_SrCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2	Pb[OH]2[C]	-8.150	Pb[OH]2[C] + 2 H+ -> 2 H2O + 1 Pb+2
Cem07_Al[OH]3[am]	13.76	Cem07_Al[OH]3[am] + 1 H2O -> 1 Al[OH]4- + 1 H+	PbMoO4[c]	15.80	PbMoO4[c] -> 1 MoO4-2 + 1 Pb+2
Cem07_Brucite	-16.83	Cem07_Brucite + 2 H+ -> 2 H2O + 1 Mg+2	PbOH[Sb[OH]6]_exp1	12.00	PbOH[Sb[OH]6]_exp1 + 1 H+ -> 1 H2O + 1 Pb+2 + 1 Sb[OH]6-
Cem07_C4FcH12	-20.47	Cem07_C4FcH12 + 4 H+ -> 1 CO3-2 + 4 Ca+2 + 2 Fe[OH]4- + 10 H2O	Sn[OH]2[s]	1.447	Sn[OH]2[s] + 2 H+ -> 2 H2O + 1 Sn+2
Cem07_Calcite	8.485	Cem07_Calcite -> 1 CO3-2 + 1 Ca+2	Strengite	48.00	Strengite + 2 H2O -> 1 Fe[OH]4- + 4 H+ + 1 PO4-3
Cem07_SiO2[am]	24.21	Cem07_SiO2[am] + 2 H2O -> 2 H+ + 1 H2SiO4-2	Wairakite	113.6	Wairakite + 10 H2O -> 2 Al[OH]4- + 1 Ca+2 + 8 H+ + 4 H2SiO4-2
Co2SiO4	6.289	Co2SiO4 + 2 H+ -> 2 Co+2 + 1 H2SiO4-2	Willemite	6.289	Willemite + 2 H+ -> 1 H2SiO4-2 + 2 Zn+2
Cu[OH]2[s]	-8.640	Cu[OH]2[s] + 2 H+ -> 1 Cu+2 + 2 H2O			

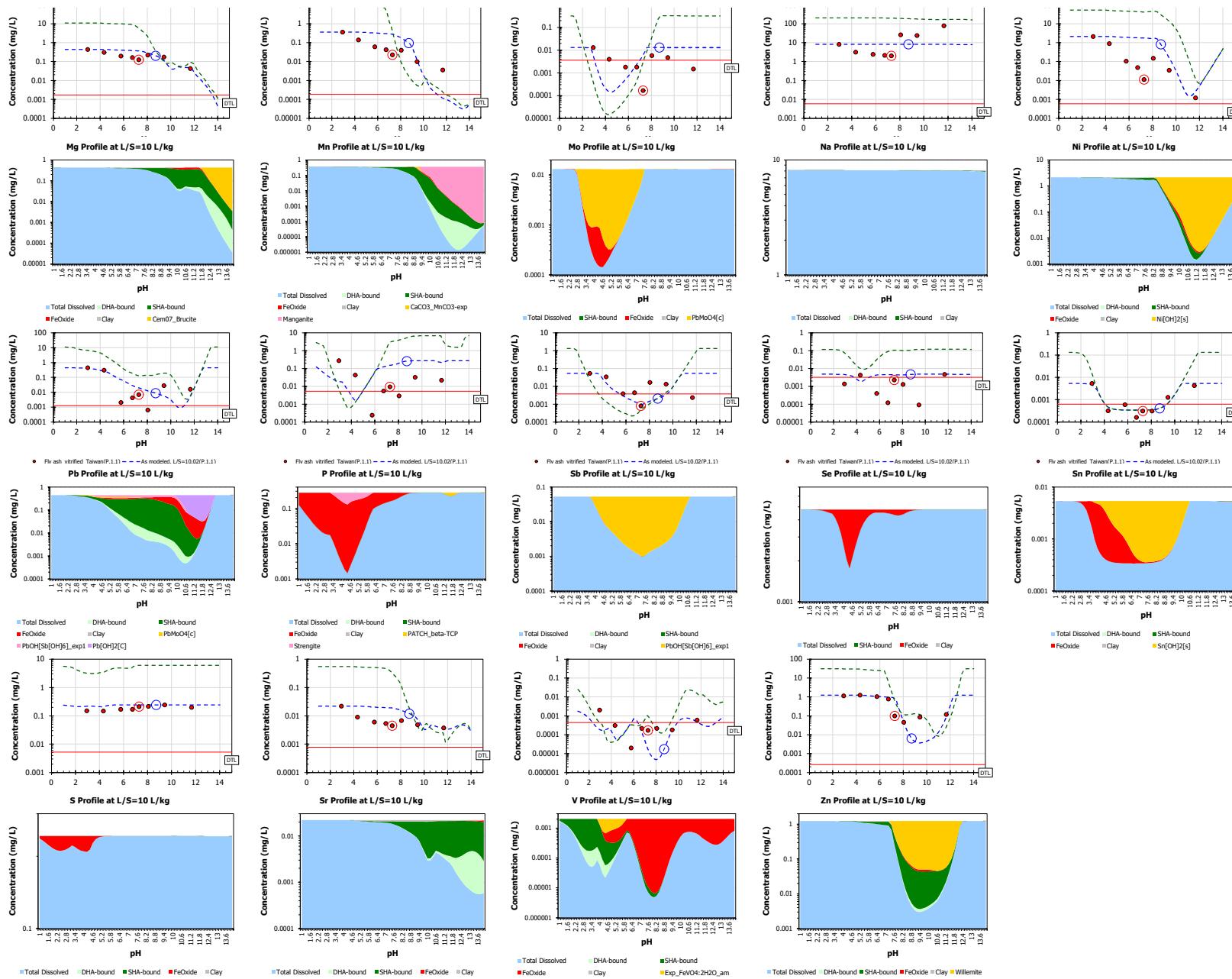
VITRIFIED MSWI FLY ASH

COMPARISON AND PARTITIONING



VITRIFIED MSWI FLY ASH

COMPARISON AND PARTITIONING



Model Comparison: residuals - Concentration

Name Vitrified MSWI Fly ash TW

Legend

Total Average Deviation Square root of the sum of the squared values of residuals divided by the number of values, over the entire X range.

User Average Deviation Square root of the sum of the squared values of residuals divided by the number of values, over the user defined X range.

Fractional Average Deviation: Square root of the sum of the squared values of residuals divided by the number of values, over the fraction.

Note that the Total and User Average Deviation columns are averages as well.

Residual details, concentrations

Residuals as log(model/sample)

	8	7	6	5	4	3	2	1	Total Avg
Fraction									Deviation
pH	2.93	4.31	5.75	6.72	7.27	8.05	9.42	11.7	0.23
Al	-0.01	1.13	0.84	-0.31	-1.09	-0.33	0.21	0.33	0.23
As	-0.55	-0.82	1.05	1.48	0.30	0.58	0.65	1.39	0.33
Ba	0.00	0.48	0.82	0.97	1.15	0.87	-0.28	-1.17	0.29
Ca	0.00	0.27	0.34	0.37	0.41	0.28	0.11	0.14	0.10
Cd	0.01	-0.02	-0.07	-0.26	0.12	-0.23	-0.42	0.10	0.07
Cl	0.60	0.99	0.65	0.69	0.00	0.87	0.86	1.11	0.28
Co	0.00	0.21	1.12	1.52	1.92	0.42	-1.06	0.38	0.37
CO32-	-	-	-	-	-	-	-	-	-
Cr	-1.15	-0.37	1.98	0.68	0.25	0.44	-1.09	0.92	0.36
Cu	-0.03	0.07	0.60	0.59	0.99	1.07	-0.93	-0.03	0.24
F	-	-	-	-	-	-	-	-	-
Fe	-0.04	0.16	2.47	0.42	-0.12	0.83	-1.51	1.17	0.41
B	0.00	0.43	0.66	0.72	0.74	0.51	0.57	0.56	0.20
Si	0.00	0.92	1.30	1.29	1.27	1.30	0.87	0.60	0.37
Hg	-	-	-	-	-	-	-	-	-
K	0.54	1.04	0.70	0.73	0.00	0.88	0.90	1.16	0.29
Li	0.00	0.35	0.91	0.52	0.41	1.13	1.25	0.43	0.26
Mg	0.00	0.15	0.31	0.36	0.47	0.11	-0.18	0.04	0.09
Mn	0.00	0.41	0.73	0.85	1.08	0.64	0.12	-1.48	0.29
Mo	-0.34	-1.42	-0.55	0.10	1.55	0.35	0.43	0.94	0.31
Na	0.00	0.41	0.52	0.58	0.60	-0.51	-0.47	-0.99	0.20
Ni	0.00	0.37	1.26	1.58	2.19	1.05	0.65	0.34	0.41
Pb	-0.05	-0.07	1.55	0.83	0.45	1.28	-0.71	-0.48	0.30
P	-1.17	-1.42	2.16	1.35	1.21	1.83	0.93	1.01	0.51
Sb	0.00	-0.34	-0.12	-0.50	0.11	-1.05	-0.59	1.35	0.24
Se	0.52	-0.36	1.03	1.58	0.29	0.53	1.72	0.00	0.34
Sn	-0.07	0.17	-0.24	0.32	0.04	0.05	-0.27	0.10	0.07
S	0.15	0.16	0.16	0.15	0.06	0.05	0.00	0.08	0.04
Sr	0.00	0.38	0.53	0.56	0.63	0.36	0.20	-0.02	0.14
U	-	-	-	-	-	-	-	-	-
V	-1.03	-0.65	1.32	-0.21	-1.01	-1.65	-0.06	-0.12	0.33
Zn	0.04	-0.01	0.01	0.10	0.91	-0.03	-1.36	-0.13	0.21
Avg Deviation	0.08	0.12	0.20	0.16	0.17	0.16	0.15	0.15	0.26

Yellow = own pH

All residuals within + 1 or - 1 are considered to represent a good fit.