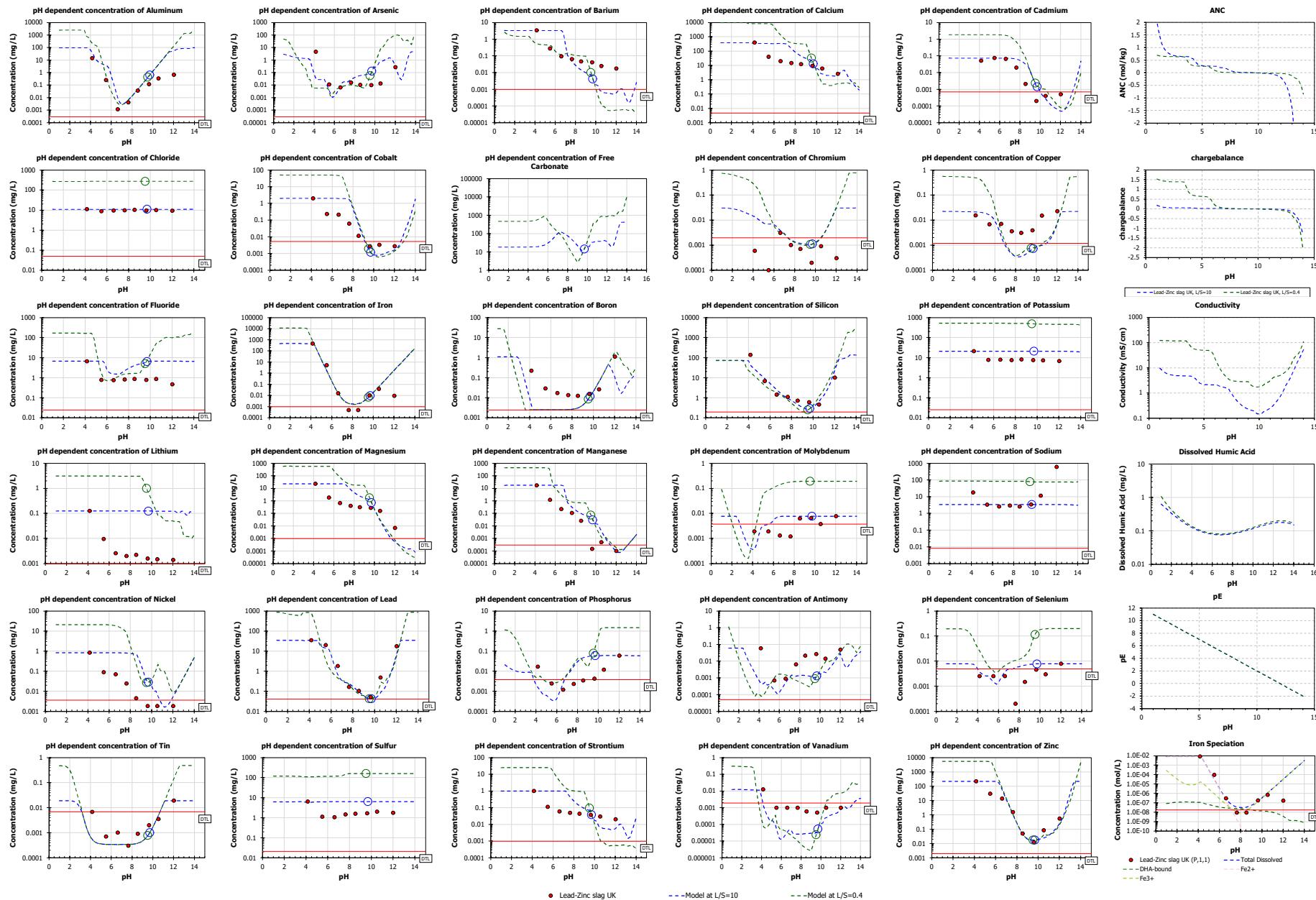
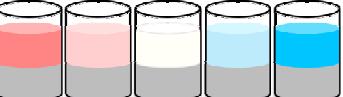


LEAD-ZINC SLAG UK

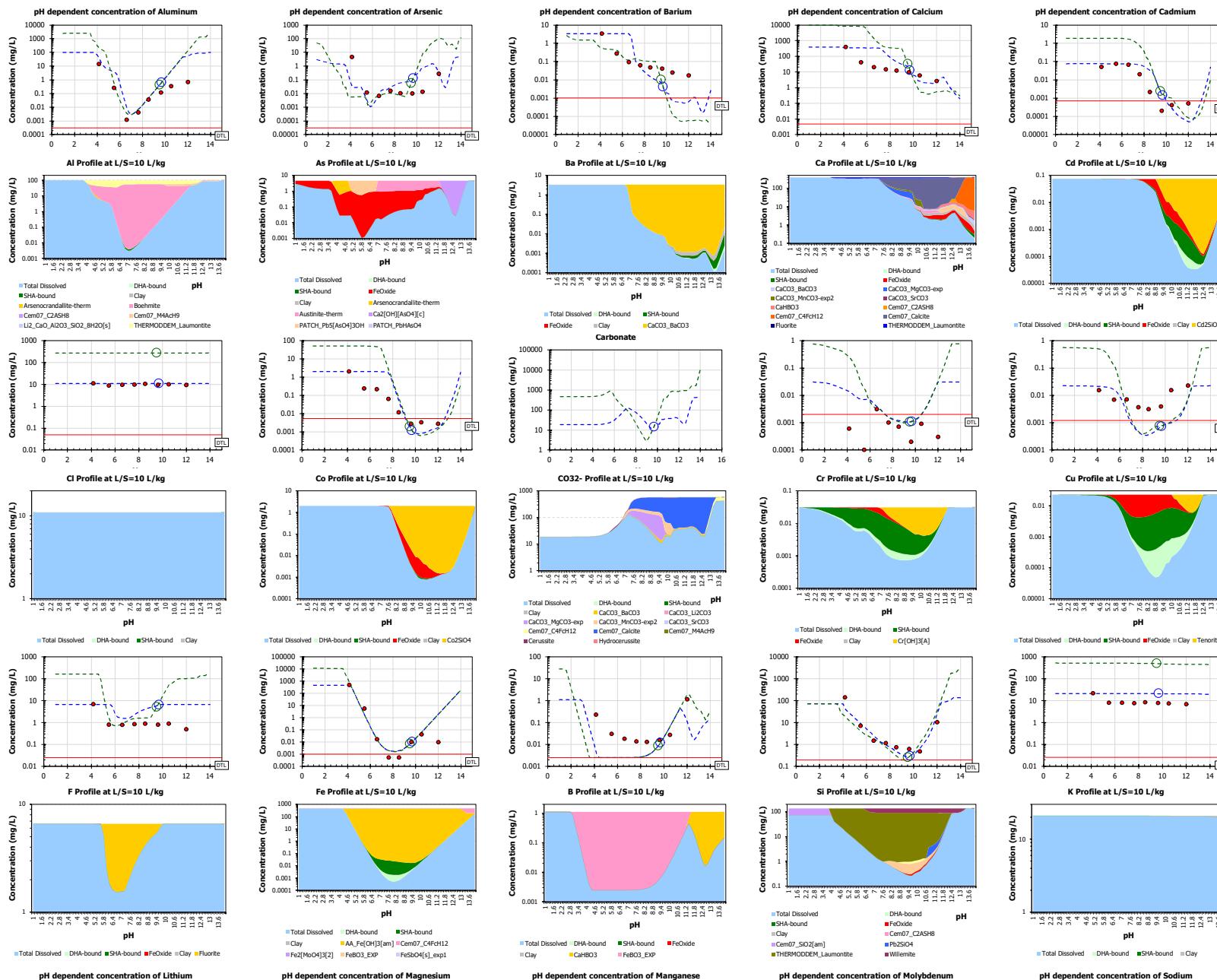
COMPARISON pH DEPENDENCE WITH MODEL



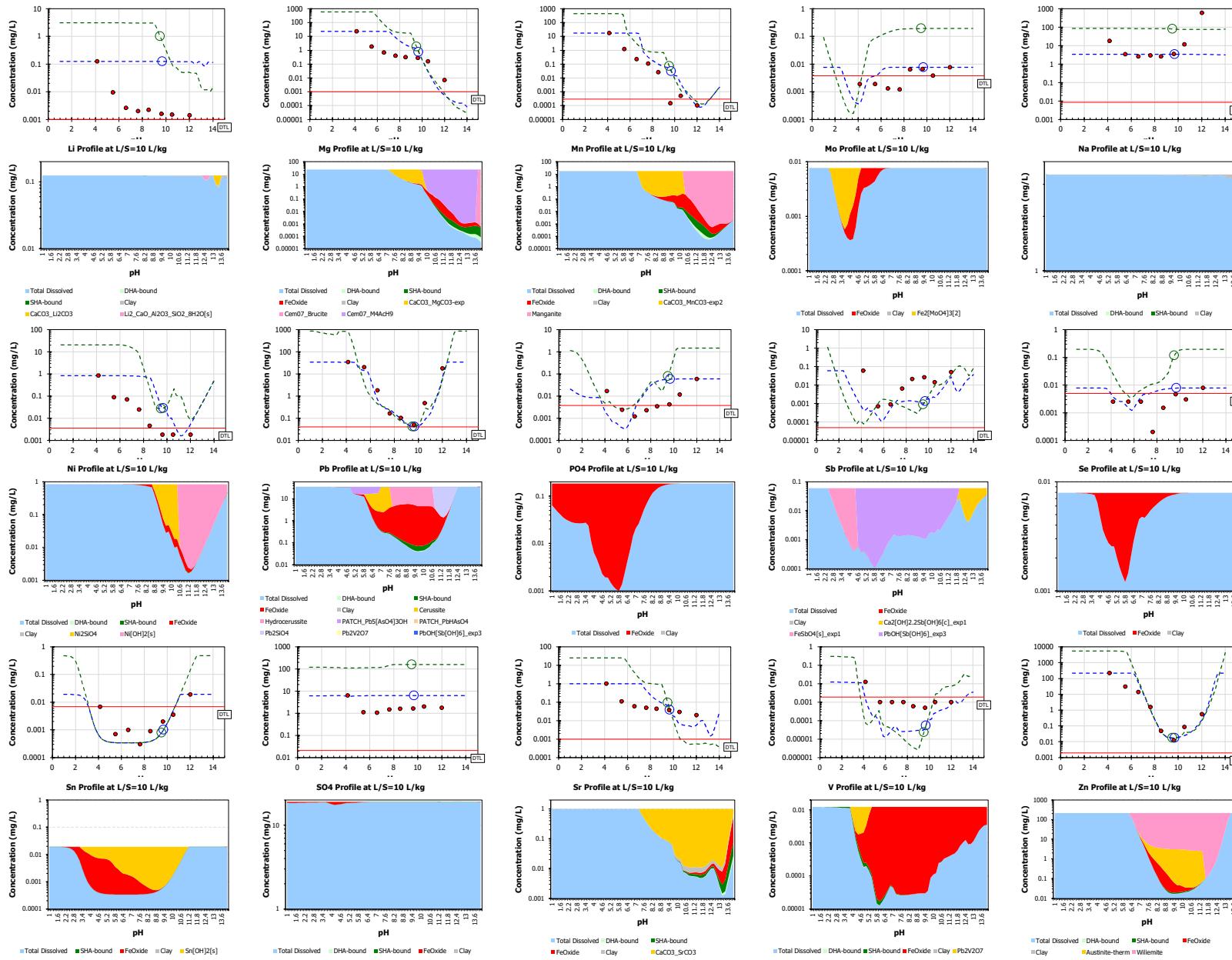
Object Name	pH Dependent Leaching Test Model														
	Lead-Zinc slag UK														
pH Dependent Leaching Test Scenario															
															
	Lab Test				Extra L/S Simulation										
Lab Test															
Model Parameters															
Available Content															
Entity	Unit	Default	Entity	mg/kg	Entity	mg/kg	Entity	mg/kg							
c0		-6.197	Al	981.1	B	11.16	Sb	0.5960							
c1		-0.3078	As	46.00	Si	1358	Se	0.07900							
c2		0.003221	Ba	33.36	K	209.2	Sn	0.1910							
c3		0.003526	Ca	3879	Li	1.246	SO4	189.7							
c4		-0.0001801	Cd	0.7480	Mg	232.2	Sr	9.993							
c5		0	Cl	110.1	Mn	176.0	V	0.1230							
Clay	mg/kg	3000	Co	20.19	Mo	0.07700	Zn	2155							
Hydrous Ferric Oxide	mg/kg	300.0	CO32-	5800	Na	34.00									
L/S	L/kg	10.00	Cr	0.3100	Ni	8.364									
pE		0.2000	Cu	0.2250	NO3	6.200E-09									
pH		11.80	F	66.00	Pb	345.2									
Solid Humic Acid	mg/kg	15.00	Fe	4597	PO4	1.828									
Simulated Low L/S	L/kg	0.4000													
Minerals															
Name	Log(K)	Reaction													
AA_Fe[OH]3[am]	16.60	AA_Fe[OH]3[am] + 1 H2O -> 1 Fe[OH]4- + 1 H+													
Adamite-therm	12.64	Adamite-therm + 1 H+ -> 1 AsO4-3 + 1 H2O + 2 Zn+2													
Arsenocrandallite-therm	95.56	Arsenocrandallite-therm + 6 H2O -> 3 Al[OH]4- + 2 AsO4-3 + 1 Ca+2 + 7 H+													
Austinite-therm	11.47	Austinite-therm + 1 H+ -> 1 AsO4-3 + 1 Ca+2 + 1 H2O + 1 Zn+2													
Boehmite	14.42	Boehmite + 2 H2O -> 1 Al[OH]4- + 1 H+													
Ca2[OH][AsO4][c]	4.000	Ca2[OH][AsO4][c] + 1 H+ -> 1 AsO4-3 + 2 Ca+2 + 1 H2O													
Ca2[OH]2.2Sb[OH]6[c]_e:	5.000	Ca2[OH]2.2Sb[OH]6[c]_exp1 + 2 H+ -> 2 Ca+2 + 2 H2O + 2 Sb[OH]6-													
Ca2Cd[PO4]2	32.95	Ca2Cd[PO4]2 -> 2 Ca+2 + 1 Cd+2 + 2 PO4-3													
CaCO3_BaCO3	21.00	CaCO3_BaCO3 -> 1 Ba+2 + 2 CO3-2 + 1 Ca+2													
CaCO3_Li2CO3	21.30	CaCO3_Li2CO3 -> 2 CO3-2 + 1 Ca+2 + 2 Li+													
CaCO3_MgCO3-exp	18.02	CaCO3_MgCO3-exp -> 2 CO3-2 + 1 Ca+2 + 1 Mg+2													
CaCO3_MnCO3-exp2	19.78	CaCO3_MnCO3-exp2 -> 2 CO3-2 + 1 Ca+2 + 1 Mn+2													
CaCO3_SrCO3	19.85	CaCO3_SrCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2													
CaHBO3	-2.097	CaHBO3 + 1 H+ -> 1 Ca+2 + 1 H2BO3-													
Cd2SiO4	6.059	Cd2SiO4 + 2 H+ -> 2 Cd+2 + 1 H2SiO4-2													
Cem07_Brucite	-16.83	Cem07_Brucite + 2 H+ -> 2 H2O + 1 Mg+2													
Cem07_C2ASH8	17.40	Cem07_C2ASH8 -> 2 Al[OH]4- + 2 Ca+2 + 3 H2O + 1 H2SiO4-2													
Cem07_C4FcH12	-20.47	Cem07_C4FcH12 + 4 H+ -> 1 CO3-2 + 4 Ca+2 + 2 Fe[OH]4- + 10 H2O													
Cem07_Calcite	8.485	Cem07_Calcite -> 1 CO3-2 + 1 Ca+2													
Cem07_M4AcH9	-4.823	Cem07_M4AcH9 + 4 H+ -> 2 Al[OH]4- + 1 CO3-2 + 7 H2O + 4 Mg+2													
Cem07_SiO2[am]	24.21	Cem07_SiO2[am] + 2 H2O -> 2 H+ + 1 H2SiO4-2													
Cerussite	13.20	Cerussite -> 1 CO3-2 + 1 Pb+2													
Co2SiO4	5.289	Co2SiO4 + 2 H+ -> 2 Co+2 + 1 H2SiO4-2													
			Name	Log(K)	Reaction										
			Cr[OH]3[A]	68.13	Cr[OH]3[A] + 1 H2O -> 1 CrO4-2 + 5 H+ + 3 e-										
			Fe2[MoO4]3[86.35	Fe2[MoO4]3[2] + 8 H2O -> 2 Fe[OH]4- + 8 H+ + 3 MoO4-2										
			FeBO3_EXP	32.48	FeBO3_EXP + 2 H2O -> 1 Fe[OH]4- + 2 H+ + 1 H2BO3-										
			FeSbO4[s]_ex	32.48	FeSbO4[s]_exp1 + 6 H2O -> 1 Fe[OH]4- + 2 H+ + 1 Sb[OH]6-										
			Fluorite	10.96	Fluorite -> 1 Ca+2 + 2 F-										
			Hydrocerussit	18.77	Hydrocerussite + 2 H+ -> 2 CO3-2 + 2 H2O + 3 Pb+2										
			Li2_CaO_Al2C	22.69	Li2_CaO_Al2O3_SiO2_8H2O[s] -> 2 Al[OH]4- + 1 Ca+2 + 3 H2O + 1 H2SiO4-2 + 2 Li+										
			Manganite	-25.27	Manganite + 3 H+ + 1 e- -> 2 H2O + 1 Mn+2										
			Ni[OH]2[s]	-10.80	Ni[OH]2[s] + 2 H+ -> 2 H2O + 1 Ni+2										
			Ni2SiO4	5.498	Ni2SiO4 + 2 H+ -> 1 H2SiO4-2 + 2 Ni+2										
			PATCH_BaCa!	7.412	PATCH_BaCaSO4[50%Ba] -> 0.5 Ba+2 + 0.5 Ca+2 + 1 SO4-2										
			PATCH_beta:-	28.93	PATCH_beta-TCP -> 3 Ca+2 + 2 PO4-3										
			PATCH_Pb5[A	62.13	PATCH_Pb5[AsO4]3OH + 1 H+ -> 3 AsO4-3 + 1 H2O + 5 Pb+2										
			PATCH_PbHA	23.97	PATCH_PbHAsO4 -> 1 AsO4-3 + 1 H+ + 1 Pb+2										
			Pb2SiO4	6.289	Pb2SiO4 + 2 H+ -> 1 H2SiO4-2 + 2 Pb+2										
			Pb2V2O7	0.9500	Pb2V2O7 + 3 H+ -> 1.5 H2O + 1 Pb+2 + 1 VO2+										
			PbOH[Sb]O]	11.00	PbOH[Sb]O]_exp3 + 1 H+ -> 1 H2O + 1 Pb+2 + 1 Sb[OH]6-										
			Sn[OH]2[s]	1.447	Sn[OH]2[s] + 2 H+ -> 2 H2O + 1 Sn+2										
			Strengite	48.00	Strengite + 2 H2O -> 1 Fe[OH]4- + 4 H+ + 1 PO4-3										
			Tenorite	-7.644	Tenorite + 2 H+ -> 1 Cu+2 + 1 H2O										
			THERMODDE	120.8	THERMODDEM_Laumontite + 8 H2O -> 2 Al[OH]4- + 1 Ca+2 + 8 H+ + 4 H2SiO4-2										
			Willemite	6.289	Willemite + 2 H+ -> 1 H2SiO4-2 + 2 Zn+2										

LEAD-ZINC SLAG UK

COMPARISON AND PARTITIONING



COMPARISON AND PARTITIONING



Model Comparison: residuals - Concentration

Name Lead-Zinc slag UK

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									
pH	4.17	5.50	6.61	7.64	8.55	9.63	10.5	12.0	Deviation
Al	0.45	1.17	0.90	0.22	0.13	0.68	1.14	1.93	0.35
As	-2.15	-0.50	0.12	0.25	0.76	0.99	1.72	0.07	0.38
Ba	0.00	1.09	1.55	-0.12	-0.51	-0.87	-1.49	-1.44	0.37
Ca	-0.04	0.93	1.24	1.01	0.57	0.26	-0.36	-0.06	0.25
Cd	0.17	0.00	0.05	0.51	1.19	0.88	-0.18	-1.00	0.23
Cl	0.00	0.10	0.07	0.05	0.03	0.07	0.04	0.07	0.02
Co	0.00	0.93	0.97	1.43	0.64	-0.30	-0.61	-0.20	0.27
CO32-	-	-	-	-	-	-	-	-	-
Cr	1.36	1.84	0.10	0.21	0.23	0.74	0.34	2.01	0.40
Cu	0.14	0.36	-0.50	-0.90	-0.93	-0.73	-1.19	-0.49	0.26
F	0.00	0.93	0.31	0.45	0.67	0.90	0.89	1.15	0.27
Fe	-0.01	-0.47	-0.01	0.60	0.59	-0.05	0.21	2.31	0.31
B	-1.72	-1.07	-0.84	-0.71	-0.59	-0.17	0.40	-1.02	0.33
Si	-0.49	0.12	0.25	-0.06	-0.18	-0.34	0.17	0.42	0.10
K	0.00	0.44	0.43	0.45	0.41	0.44	0.46	0.49	0.15
Li	0.00	1.12	1.68	1.79	1.75	1.89	1.91	1.94	0.58
Mg	0.00	1.10	1.54	1.31	0.91	0.54	-0.75	-1.12	0.36
Mn	0.00	1.17	1.89	0.47	0.59	2.42	0.89	-0.02	0.44
Mo	-0.70	0.30	0.76	0.81	0.09	0.07	0.31	0.00	0.17
Na	-0.72	0.00	0.11	0.07	0.12	-0.01	-0.54	-2.25	0.30
Ni	0.00	0.97	1.07	1.51	2.18	1.19	0.57	0.48	0.42
Pb	-0.01	-0.15	-0.43	0.18	-0.04	-0.09	-0.76	-0.40	0.12
PO4	-	-	-	-	-	-	-	-	-
Sb	-2.11	-0.59	-0.24	-0.70	-1.18	-1.36	-0.81	-0.28	0.38
Se	0.18	-0.22	0.17	1.44	0.67	0.23	0.42	0.00	0.21
Sn	-1.10	-0.31	-0.47	0.06	-0.37	-0.33	0.16	0.00	0.17
SO4	-	-	-	-	-	-	-	-	-
Sr	0.00	0.95	1.23	0.94	0.49	0.14	-0.61	-0.54	0.26
V	-0.90	-1.24	-1.46	-1.59	-1.33	-1.09	-0.58	-0.08	0.40
Zn	0.00	0.85	0.60	-0.19	-0.03	0.16	-0.53	-0.33	0.16
Avg Deviation	0.15	0.16	0.17	0.16	0.16	0.16	0.16	0.20	0.28

Yellow = own pH

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