

Object Name pH Dependent Leaching Test Model
MSWI Bottom ash DE

pH Dependent Leaching Test Scenario



Lab Test

Extra L/S Simulation

Lab Test

Model Parameters

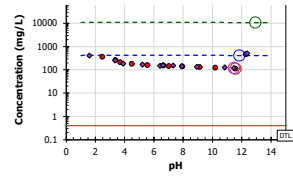
Entity	Unit	Default	Entity	mg/kg	Entity	mg/kg	Entity	mg/kg
c0		-3.401	Acetic acid	2.220E-08	B	1.081E-09	Sb	1.218E-08
c1		-1.683	Al	5437	Si	7712	Se	0.01345
c2		0.3256	As	1.635	Hg	0.009140	Sn	1.187E-08
c3		-0.02570	Ba	11.56	K	2179	SO4	9554
c4		0.0007423	Ca	4.348E+04	Li	6.941E-10	Sr	99.16
c5		-2.906E-07	Cd	6.997	Mg	2528	V	72.44
Clay	mg/kg	3000	Cl	1973	Mn	299.0	Zn	1371
Hydrous Ferric Oxide	mg/kg	150.0	Co	5.893E-09	Mo	9.594E-09		
L/S	L/kg	10.60	CO32-	5.000E+04	Na	4417		
pE		2.400	Cr	13.89	Ni	25.39		
pH		11.56	Cu	729.6	Pb	324.6		
Solid Humic Acid	mg/kg	500.0	F	1.900E-09	PO4	2300		
Simulated Low L/S	L/kg	0.4000	Fe	6841				

Solid Solutions

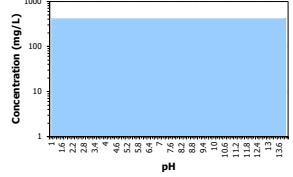
Name	End Member	Log(K)	Reaction
CSHii_ss	Cem07_Jenn_ss	-7.799	Cem07_Jenn_ss + 1.33333 H+ -> 1 CSHii_ss + 1.66667 Ca+2 + 1.76667 H2O + 1 H2SiO4-2
	Cem07_Tob_II_ss	10.36	Cem07_Tob_II_ss -> 1 CSHii_ss + 0.83333 Ca+2 + 0.33333 H+ + 0.16667 H2O + 1 H2SiO4-2
ettr_ss	AsO4_Ettringite_ss	26.79	AsO4_Ettringite_ss + 1 H+ + 8 H2O -> 2 Al[OH]4- + 3 AsO4-3 + 6 Ca+2 + 1 ettr_ss
	Ba_Ettringite_ss	4.008	Ba_Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ba+2 + 3 SO4-2 + 1 ettr_ss
	BO3_Ettringite_ss	-46.87	BO3_Ettringite_ss + 7 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 H2BO3- + 1 ettr_ss
	CrO4_Ettringite_ss	-8.592	CrO4_Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 CrO4-2 + 1 ettr_ss
	Ettringite_ss	-10.99	Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 SO4-2 + 1 ettr_ss
	MoO4_Ettringite_ss	-9.592	MoO4_Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 MoO4-2 + 1 ettr_ss
	PO4_Ettringite_ss	39.10	PO4_Ettringite_ss + 1 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 PO4-3 + 1 ettr_ss
	SeO4-2_Ettringite_ss	-8.592	SeO4-2_Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 SeO4-2 + 1 ettr_ss
	Sr_Ettringite_ss	4.008	Sr_Ettringite_ss + 4 H+ + 8 H2O -> 2 Al[OH]4- + 3 SO4-2 + 6 Sr+2 + 1 ettr_ss
	VO3_Ettringite_ss	-53.79	VO3_Ettringite_ss + 13 H+ + 2 H2O -> 2 Al[OH]4- + 6 Ca+2 + 3 VO2+ + 1 ettr_ss

Minerals

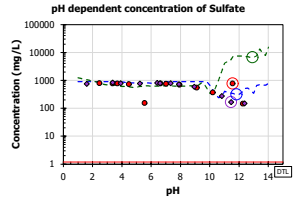
Name	Log(K)	Reaction	Name	Log(K)	Reaction
AA_Fe[OH]3[am]	16.60	AA_Fe[OH]3[am] + 1 H2O -> 1 Fe[OH]4- + 1 H+	Cem07_CAH10	7.505	Cem07_CAH10 -> 2 Al[OH]4- + 1 Ca+2 + 6 H2O
Adamite-therm	12.64	Adamite-therm + 1 H+ -> 1 AsO4-3 + 1 H2O + 2 Zn+2	Cem07_Calcite	8.485	Cem07_Calcite -> 1 CO3-2 + 1 Ca+2
alpha-TCP	25.50	alpha-TCP -> 3 Ca+2 + 2 PO4-3	Cem07_Gypsum	4.583	Cem07_Gypsum -> 1 Ca+2 + 2 H2O + 1 SO4-2
Austinite-therm	11.47	Austinite-therm + 1 H+ -> 1 AsO4-3 + 1 Ca+2 + 1 H2O + 1 Zn+2	Cr[OH]3[A]	68.13	Cr[OH]3[A] + 1 H2O -> 1 CrO4-2 + 5 H+ + 3 e-
Ba[SCr]O4[96%SO4]	9.790	Ba[SCr]O4[96%SO4] -> 1 Ba+2 + 0.04 CrO4-2 + 0.96 SO4-2	Cu[OH]2[s]	-8.640	Cu[OH]2[s] + 2 H+ -> 1 Cu+2 + 2 H2O
BaSrSO4[50%Ba]	8.221	BaSrSO4[50%Ba] -> 0.5 Ba+2 + 1 SO4-2 + 0.5 Sr+2	Exp_CaCO3_MgCO3	19.00	Exp_CaCO3_MgCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Mg+2
Ca2Cd[PO4]2	32.95	Ca2Cd[PO4]2 -> 2 Ca+2 + 1 Cd+2 + 2 PO4-3	Exp2_LDH_Pb_zc	61.00	Exp2_LDH_Pb_zc + 1 H2O -> 1 Al[OH]4- + 3 CO3-2 + 1 H+ + 3 Pb+2
Ca3[OH]2[CrO4]2[ccc]	1.000	Ca3[OH]2[CrO4]2[ccc] + 2 H+ -> 3 Ca+2 + 2 CrO4-2 + 2 H2O	Laumontite	118.0	Laumontite + 8 H2O -> 2 Al[OH]4- + 1 Ca+2 + 8 H+ + 4 H2SiO4-2
Ca4Cd[PO4]3OH	39.23	Ca4Cd[PO4]3OH + 1 H+ -> 4 Ca+2 + 1 Cd+2 + 1 H2O + 3 PO4-3	Manganite	-25.27	Manganite + 3 H+ + 1 e- -> 2 H2O + 1 Mn+2
Ca5[OH][VO4]3[cx]	-54.00	Ca5[OH][VO4]3[cx] + 13 H+ -> 5 Ca+2 + 7 H2O + 3 VO2+	Ni[OH]2[s]	-10.80	Ni[OH]2[s] + 2 H+ -> 2 H2O + 1 Ni+2
CaCO3_SrCO3	19.85	CaCO3_SrCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2	NiHPO4	25.00	NiHPO4 -> 1 H+ + 1 Ni+2 + 1 PO4-3
Cd[OH]2[C]	-13.65	Cd[OH]2[C] + 2 H+ -> 1 Cd+2 + 2 H2O	Pb[OH]2[C]	-8.150	Pb[OH]2[C] + 2 H+ -> 2 H2O + 1 Pb+2
CdHPO4	26.48	CdHPO4 -> 1 Cd+2 + 1 H+ + 1 PO4-3	Pb2V2O7	0.9500	Pb2V2O7 + 3 H+ -> 1.5 H2O + 1 Pb+2 + 1 VO2+
Cem07_Al[OH]3[am]	13.76	Cem07_Al[OH]3[am] + 1 H2O -> 1 Al[OH]4- + 1 H+	Pb3[VO4]2	-3.070	Pb3[VO4]2 + 4 H+ -> 2 H2O + 1.5 Pb+2 + 1 VO2+
Cem07_Brucite	-16.83	Cem07_Brucite + 2 H+ -> 2 H2O + 1 Mg+2	Rhodochrosite	10.41	Rhodochrosite -> 1 CO3-2 + 1 Mn+2
Cem07_C3AH6	-35.14	Cem07_C3AH6 + 4 H+ -> 2 Al[OH]4- + 3 Ca+2 + 4 H2O	Strengite	48.00	Strengite + 2 H2O -> 1 Fe[OH]4- + 4 H+ + 1 PO4-3
Cem07_C3FH6	-30.82	Cem07_C3FH6 + 4 H+ -> 3 Ca+2 + 2 Fe[OH]4- + 4 H2O	Willemite	6.289	Willemite + 2 H+ -> 1 H2SiO4-2 + 2 Zn+2



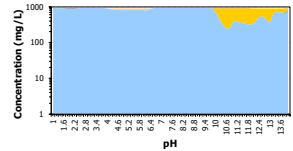
Na Profile at L/S=10.6 L/kg



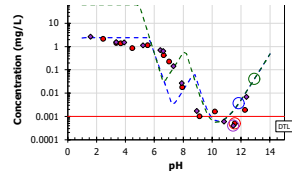
Na Profile at L/S=10.6 L/kg



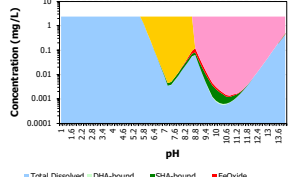
SO4 Profile at L/S=10.6 L/kg



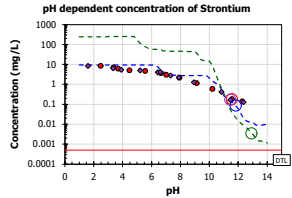
SO4 Profile at L/S=10.6 L/kg



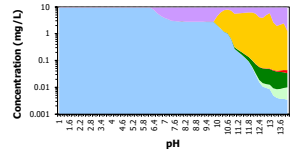
Ni Profile at L/S=10.6 L/kg



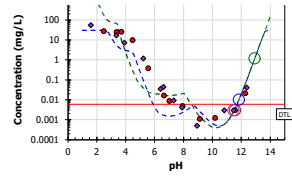
Ni Profile at L/S=10.6 L/kg



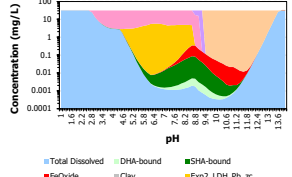
Sr Profile at L/S=10.6 L/kg



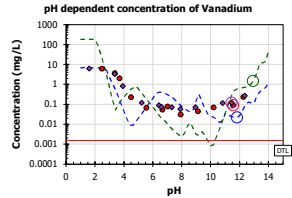
Sr Profile at L/S=10.6 L/kg



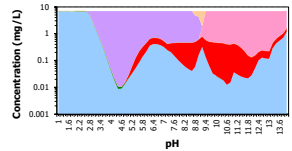
Pb Profile at L/S=10.6 L/kg



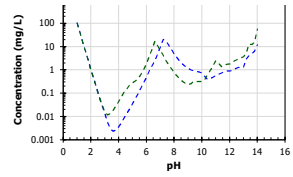
Pb Profile at L/S=10.6 L/kg



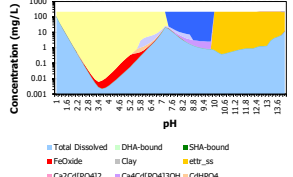
V Profile at L/S=10.6 L/kg



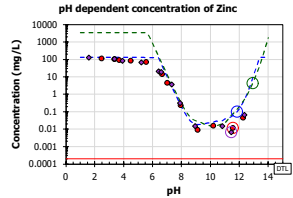
V Profile at L/S=10.6 L/kg



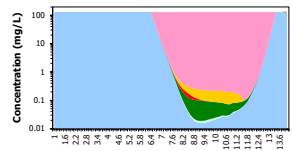
PO4 Profile at L/S=10.6 L/kg



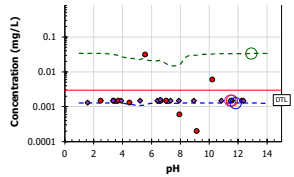
PO4 Profile at L/S=10.6 L/kg



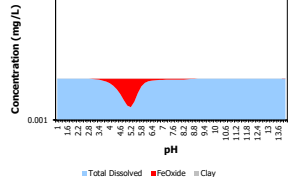
Zn Profile at L/S=10.6 L/kg



Zn Profile at L/S=10.6 L/kg



Se Profile at L/S=10.6 L/kg



Se Profile at L/S=10.6 L/kg

Model Comparison: residuals - Concentration

Name MSWI Bottom ash DE

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)													
Fraction	9	8	7	6	1	11	10	4	3	2	5	12	Total Avg
pH	2.47	3.36	3.69	4.48	5.55	6.65	7.02	7.93	9.12	10.2	11.6	12.3	Deviation
Al	0.05	0.19	0.20	-0.07	-1.49	-0.72	-0.86	-0.10	-0.68	0.11	0.54	0.88	0.19
As	-0.19	0.07	0.30	0.80	0.80	0.43	0.55	0.13	-0.28	0.14	0.70	1.22	0.17
Ba	0.03	-0.03	0.14	-0.14	-0.32	-0.47	-0.79	-0.79	-0.75	-0.40	-0.57	-0.59	0.14
Br	-	-	-	-	-	-	-	-	-	-	-	-	-
Ca	0.05	0.14	0.15	0.17	0.14	-0.11	-0.21	-0.15	0.06	0.06	-0.26	-0.54	0.06
Cd	0.22	0.38	0.02	0.54	0.12	-1.37	-1.65	-1.56	-0.85	-1.00	-0.08	0.41	0.26
Cl	0.05	0.08	0.09	0.11	0.21	0.12	0.13	0.15	0.15	0.13	0.10	0.17	0.04
Co	-	-	-	-	-	-	-	-	-	-	-	-	-
CO32-	-	-	-	-	-	-	-	-	-	-	-	-	-
Cr	-0.07	0.23	0.46	0.86	1.34	0.93	0.59	-0.33	-0.71	-0.28	0.15	0.68	0.19
Cu	0.05	0.32	0.32	0.66	1.04	1.18	0.76	-0.36	-0.77	-0.97	-0.67	0.03	0.20
F	-	-	-	-	-	-	-	-	-	-	-	-	-
Fe	-0.05	0.31	-0.05	-1.15	-1.93	0.20	-0.49	-0.51	-0.65	-0.23	1.02	1.71	0.27
B	-	-	-	-	-	-	-	-	-	-	-	-	-
Si	0.05	0.31	0.15	-0.48	-0.77	-0.75	-0.74	-0.63	-0.29	-0.62	0.32	1.02	0.17
Hg	-0.35	-0.87	-1.21	-1.69	-1.71	-1.61	-1.57	-1.47	-1.35	-1.27	-1.16	-1.14	0.39
K	0.05	0.18	0.27	0.33	0.39	0.43	0.47	0.49	0.54	0.58	0.67	0.58	0.13
Li	-	-	-	-	-	-	-	-	-	-	-	-	-
Mg	0.05	0.18	0.25	0.35	0.37	-0.87	-0.93	-0.83	-0.31	0.52	-0.04	-0.37	0.15
Mn	0.05	0.14	0.22	0.29	0.26	0.45	0.55	0.61	0.30	0.75	-0.24	-0.48	0.12
Mo	-	-	-	-	-	-	-	-	-	-	-	-	-
Na	0.05	0.20	0.29	0.36	0.42	0.43	0.46	0.47	0.50	0.52	0.56	-0.05	0.11
Ni	0.05	0.22	0.24	0.44	0.32	-1.07	-1.48	-0.18	1.18	-0.37	0.63	0.69	0.21
NO3	-	-	-	-	-	-	-	-	-	-	-	-	-
Pb	0.05	-0.66	-0.86	-0.65	-1.22	-0.99	-0.76	-0.31	0.36	-0.35	0.08	0.47	0.19
PO4	-	-	-	-	-	-	-	-	-	-	-	-	-
Sb	-	-	-	-	-	-	-	-	-	-	-	-	-
Se	-0.07	-0.08	-0.08	-0.04	-1.43	-0.08	-0.08	0.32	0.80	-2.03	-1.49	-1.36	0.28
Sn	-	-	-	-	-	-	-	-	-	-	-	-	-
SO4	0.05	0.06	0.07	0.05	0.72	0.06	0.08	0.11	0.21	0.09	-0.36	0.46	0.08
Sr	0.05	0.13	0.19	0.26	0.30	0.17	0.11	0.11	0.38	0.41	-0.14	-0.71	0.09
Th	-	-	-	-	-	-	-	-	-	-	-	-	-
U	-	-	-	-	-	-	-	-	-	-	-	-	-
V	0.04	-0.76	-1.09	-1.41	0.05	0.87	0.60	0.46	0.74	-0.50	-0.58	-0.56	0.21
Zn	0.05	0.10	0.14	0.19	0.26	0.49	0.33	0.04	0.33	0.22	0.72	0.91	0.12
Avg Deviat	0.02	0.07	0.09	0.14	0.19	0.16	0.17	0.13	0.14	0.15	0.13	0.17	0.17

Yellow = own pH All residuals within + 1 or - 1 are considered to represent a good fit.