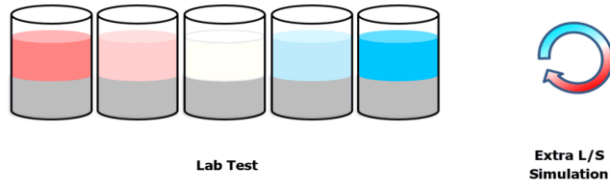


**Object
Name**

**pH Dependent Leaching Test Model
Drinkingwater Sludge NL**

pH Dependent Leaching Test Scenario



Lab Test

Model Parameters

Entity	Unit	Default	Available Content					
			Entity	mg/kg				
c0		-3.647	Acetic acid	2.220E-08	F	4.790	Pb	0.5900
c1		-0.5858	Ag	1.079E-08	Fe	1.208E+04	PO4	57.17
c2		-0.1176	Al	111.9	B	57.06	Sb	0.4467
c3		0.05073	As	0.3178	Si	1261	Se	0.4905
c4		-0.004665	Ba	245.8	Hg	2.006E-08	Sn	0.1335
c5		0.0001337	Br	28.21	K	308.6	S	809.4
Clay	mg/kg	3000	Ca	5.448E+04	Li	0.1392	Sr	246.6
Hydrous Ferric Oxide	mg/kg	750.0	Cd	0.1760	Mg	626.4	Th	2.320E-08
L/S	L/kg	10.50	Cl	268.4	Mn	796.6	U	2.380E-08
pE		5.200	Co	1.581	Mo	0.1869	V	0.8850
pH		7.720	CO32-	8.200E+04	Na	166.0	Zn	10.67
Solid Humic Acid	mg/kg	1.200E+04	Cr	1.747	Ni	6.808		
Simulated Low L/S	L/kg	0.4000	Cu	6.689	NO3	6.200E-09		

Solid Solutions

Name	End Member	Log(K)	Reaction
None			

Minerals

Name	Log(K)	Reaction	Name	Log(K)	Reaction
BaSrSO4[50%Ba]	8.221	BaSrSO4[50%Ba] -> 0.5 Ba+2 + 1 SO4-2 + 0.5 Sr+2	Cem07_SiO2[;]	24.21	Cem07_SiO2[am] + 2 H2O -> 2 H+ + 1 H2SiO4-2
beta-TCP	28.93	beta-TCP -> 3 Ca+2 + 2 PO4-3	FeBO3_EXP	32.48	FeBO3_EXP + 2 H2O -> 1 Fe[OH]4- + 2 H+ + 1 H2BO3-
Ca2[OH][AsO4][c]	4.000	Ca2[OH][AsO4][c] + 1 H+ -> 1 AsO4-3 + 2 Ca+2 + 1 H2O	Fe2[MoO4]3[;]	82.02	Fe2[MoO4]3[1] + 8 H2O -> 2 Fe[OH]4- + 8 H+ + 3 MoO4-2
Ca2[OH]2.2Sb[OH]6[c]_exp1	5.000	Ca2[OH]2.2Sb[OH]6[c]_exp1 + 2 H+ -> 2 Ca+2 + 2 H2O + 2 Sb[OH]6-	Ferrihydrite	18.41	Ferrihydrite + 1 H2O -> 1 Fe[OH]4- + 1 H+
CaCO3_BaCO3_exp	21.30	CaCO3_BaCO3_exp -> 1 Ba+2 + 2 CO3-2 + 1 Ca+2	Fluorite	10.96	Fluorite -> 1 Ca+2 + 2 F-
CaCO3_MgCO3	19.00	CaCO3_MgCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Mg+2	Kaolinite	81.80	Kaolinite + 7 H2O -> 2 Al[OH]4- + 6 H+ + 2 H2SiO4-2
CaCO3_MnCO3	21.48	CaCO3_MnCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Mn+2	Laumontite-exp	116.0	Laumontite-exp + 8 H2O -> 2 Al[OH]4- + 1 Ca+2 + 8 H+ + 4 H2SiO4-2
CaCO3_SrCO3_exp	19.85	CaCO3_SrCO3_exp -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2	Li-Albite[low2]	86.27	Li-Albite[low2] + 8 H2O -> 1 Al[OH]4- + 6 H+ + 3 H2SiO4-2 + 1 Li+
CaMoO4[c]	7.940	CaMoO4[c] -> 1 Ca+2 + 1 MoO4-2	NiHPO4	25.00	NiHPO4 -> 1 H+ + 1 Ni+2 + 1 PO4-3
Cem07_Brucite	-16.83	Cem07_Brucite + 2 H+ -> 2 H2O + 1 Mg+2	Pb[OH]2[C]	-8.150	Pb[OH]2[C] + 2 H+ -> 2 H2O + 1 Pb+2
Cem07_Calcite	8.485	Cem07_Calcite -> 1 CO3-2 + 1 Ca+2	Pb[Sb[OH]6]2	29.00	Pb[Sb[OH]6]2_exp -> 1 Pb+2 + 2 Sb[OH]6-
Cem07_Gypsum	4.583	Cem07_Gypsum -> 1 Ca+2 + 2 H2O + 1 SO4-2	Strengite	48.00	Strengite + 2 H2O -> 1 Fe[OH]4- + 4 H+ + 1 PO4-3
			ZnSiO3	18.69	ZnSiO3 + 1 H2O -> 1 H2SiO4-2 + 1 Zn+2

Model Comparison: residuals - Concentration

Sample

Name Drinkingwater Sludge NL

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

Fraction pH	Residuals as log(model/sample)									Total Avg Deviation
	8	7	6	5	4	3	2	1		
Al	-0.01	0.30	1.22	-0.58	-0.84	-1.06	-1.27	1.07	0.32	
As	0.28	0.07	-0.44	-0.27	0.34	0.92	0.48	0.01	0.15	
B	0.30	0.09	0.02	0.01	0.03	0.05	-0.05	0.00	0.06	
Ba	-0.33	0.16	0.00	-0.83	-0.93	-0.81	-1.00	-0.54	0.24	
Ca	0.00	0.09	0.20	0.01	-0.11	0.05	0.11	0.37	0.06	
Cd	-0.01	0.38	1.67	0.82	0.62	1.07	1.10	1.26	0.35	
Cl	0.14	0.27	0.08	0.04	-0.02	0.10	0.06	0.26	0.05	
Co	0.00	0.30	0.97	1.69	0.98	0.92	0.91	0.90	0.34	
CO32-	-	-	-	-	-	-	-	-	-	
Cr	-1.76	-0.75	-0.21	0.32	0.08	0.42	0.30	0.49	0.26	
Cu	-0.35	-0.21	0.12	-0.01	0.12	0.37	0.19	-0.16	0.08	
F	0.03	0.78	0.68	0.79	0.80	0.49	0.32	0.02	0.20	
Fe	-0.01	-0.96	-0.34	-0.09	-1.40	-1.34	-1.81	-2.08	0.44	
Hg	-	-	-	-	-	-	-	-	-	
K	0.01	0.21	0.12	0.19	0.21	0.28	0.23	0.17	0.07	
Li	0.01	0.09	0.28	0.69	0.84	1.11	0.56	0.30	0.21	
Mg	0.00	0.06	0.27	-0.54	-0.68	0.32	0.53	1.02	0.19	
Mn	0.00	0.21	-0.96	-0.16	-0.27	0.29	0.32	0.66	0.16	
Mo	1.81	1.72	1.79	2.00	1.22	0.61	0.80	0.02	0.50	
Na	-0.07	-0.06	-0.02	0.08	0.12	-1.20	-1.48	-2.06	0.35	
Ni	0.00	0.39	0.95	0.63	0.37	0.26	0.27	0.27	0.17	
Pb	-0.65	-0.87	-0.72	0.29	0.66	0.37	0.05	0.19	0.19	
PO4	-	-	-	-	-	-	-	-	-	
S	1.37	0.44	0.20	0.26	0.19	0.17	0.09	0.02	0.19	
Sb	-0.52	0.99	1.97	1.36	1.49	1.22	1.20	0.92	0.45	
Se	0.00	0.11	0.07	0.74	0.49	0.57	0.37	0.05	0.14	
Si	-0.22	0.30	0.91	1.13	1.13	1.12	0.70	0.40	0.29	
Sn	-0.04	0.46	-0.17	0.01	-0.43	1.53	0.49	0.64	0.23	
Sr	-0.18	-0.16	0.06	-0.43	-0.62	-0.21	-0.01	0.45	0.12	
U	-	-	-	-	-	-	-	-	-	
V	-0.36	0.17	1.03	0.62	-0.11	0.20	0.16	-0.71	0.18	
Zn	0.00	0.54	1.50	-0.05	0.00	0.33	0.11	0.66	0.22	
Avg Deviation	0.12	0.11	0.17	0.15	0.14	0.15	0.14	0.15	0.24	

Yellow = own pH All residuals within +1 or -1 are considered to represent a good fit. 0 of course the best.

B Phases missing over entire pH range
 Fe Pretty good. SHA - DHA adjustment??
 Mo Phase missing or not strong enough. Fe-Mo phase. Possible effect of very high Fe load?
 Sb Phases missing pH 4- 12
 Si No phases for Si at pH around neutral