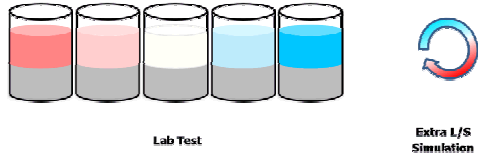


Object pH Dependent Leaching Test Model
Name Biomass wood ash NL

pH Dependent Leaching Test Scenario



Lab Test

Model Parameters

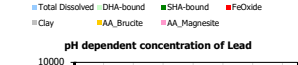
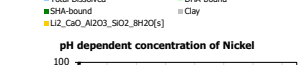
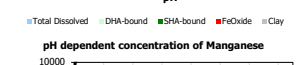
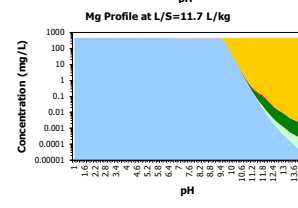
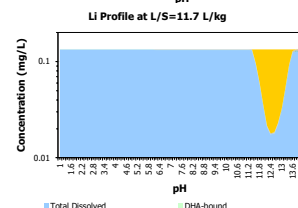
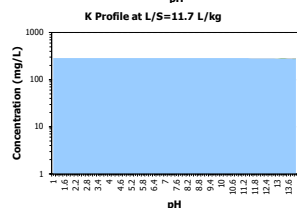
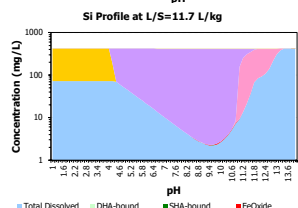
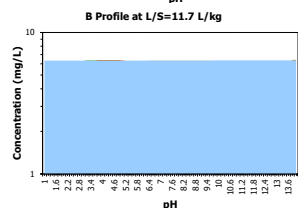
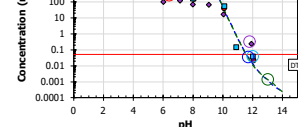
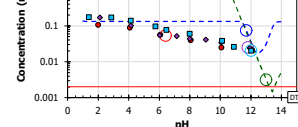
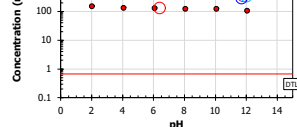
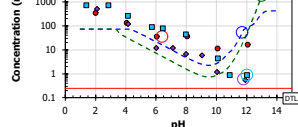
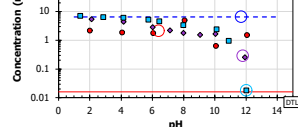
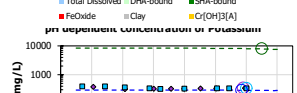
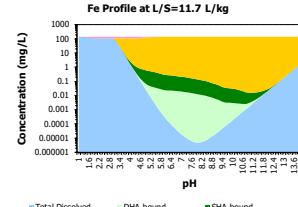
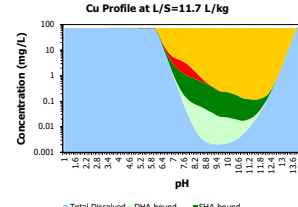
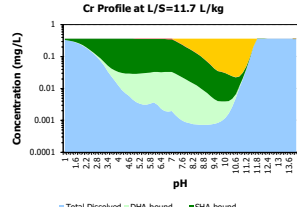
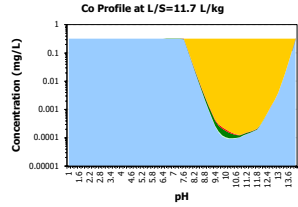
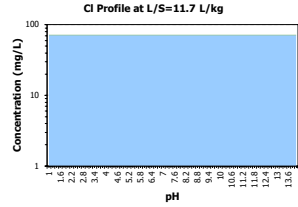
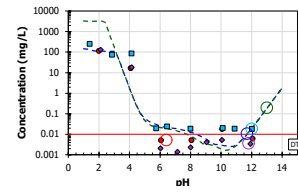
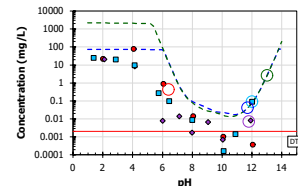
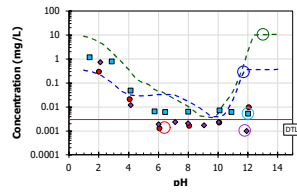
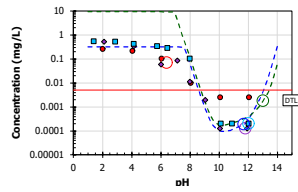
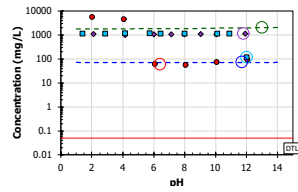
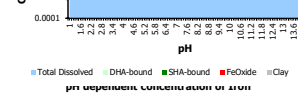
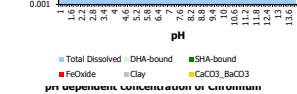
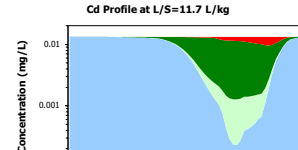
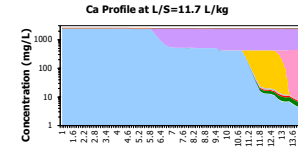
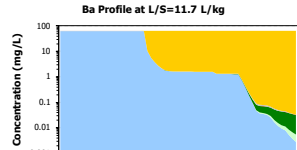
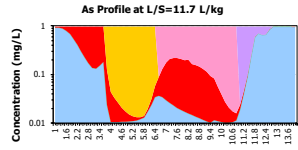
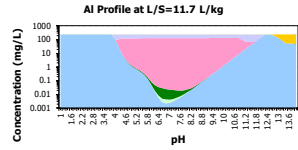
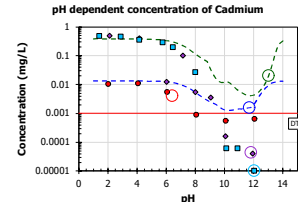
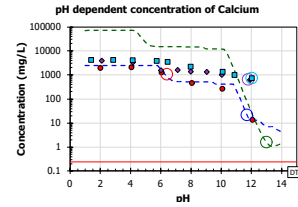
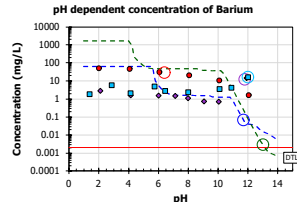
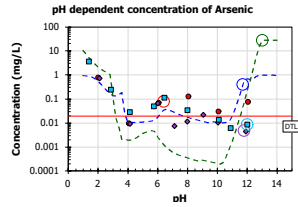
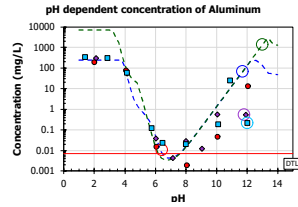
Entity	Unit	Default	Entity	mg/kg	Entity	mg/kg	Entity	mg/kg
c0		-5.538	Al	2828	Si	4932	Sn	0.08540
c1		-0.5263	As	11.25	Hg	2.006E-07	S	41.00
c2		0.1980	Ba	733.1	K	3380	Sr	129.2
c3		-0.02918	Ca	2.863E+04	Li	1.559	V	7.845
c4		0.001916	Cd	0.1552	Mg	5405	Zn	478.1
c5		-4.675E-05	Cl	832.8	Mn	494.1		
Clay	mg/kg	2.000E+04	Co	3.768	Mo	1.568		
Hydrous Ferric Oxide	mg/kg	327.0	CO32-	3.520E+04	Na	2072		
L/S	L/kg	11.66	Cr	4.259	Ni	15.12		
pE		1.000	Cu	869.9	Pb	803.9		
pH		12.00	F	10.00	PO4	630.2		
Solid Humic Acid	mg/kg	500.0	Fe	1658	Sb	6.414		
Simulated Low L/S	L/kg	0.4000	B	74.15	Se	0.09894		

Solid Solutions

Name	End Member	Log(K) Reaction
CSHi_ss	Cem07_SiO2[am]_s	24.21 Cem07_SiO2[am]_ss + 2 H2O -> 1 CSHi_ss + 2 H+ + 1 H2SiO4-2
	Cem07_Tob_I_ss	23.87 Cem07_Tob_I_ss -> 1 CSHi_ss + 2 Ca+2 + 0.8 H+ + 1.2 H2O + 2.4 H2SiO4-2
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

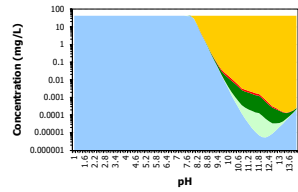
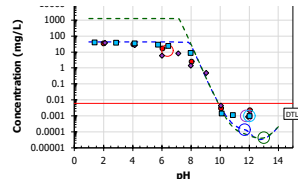
Minerals

Name	Log(K)	Reaction	Name	Log(K)	Reaction
AA_3CaO_Al2O3_6H2O]	-33.53	AA_3CaO_Al2O3_6H2O[s] + 4 H+ -> 2 Al[OH]4- + 3 Ca+2 + 4 H2O	CaSb[OH]6[s]2	18.41	CaSb[OH]6[s]2 + 6 H2O -> 1 Ca+2 + 2 Sb[OH]6-
AA_Brucite	-16.84	AA_Brucite + 2 H+ -> 2 H2O + 1 Mg+2	CoSiO4	6.289	CoSiO4 + 2 H+ -> 2 Co+2 + 1 H2SiO4-2
AA_Calcite	9.481	AA_Calcite -> 1 CO3-2 + 1 Ca+2	Cr[OH]3[A]	68.13	Cr[OH]3[A] + 1 H2O -> 1 CrO4-2 + 5 H+ + 3 e-
AA_Fe[OH]3[microcr]	18.60	AA_Fe[OH]3[microcr] + 1 H2O -> 1 Fe[OH]4- + 1 H+	Cu[OH]2[s]	-8.640	Cu[OH]2[s] + 2 H+ -> 1 Cu+2 + 2 H2O
AA_Magnesite	9.359	AA_Magnesite -> 1 CO3-2 + 1 Mg+2	Fe_Vanadate	19.18	Fe_Vanadate + 1 H2O -> 0.5 Fe[OH]4- + 1 VO2+ + 0.5 e-
Albite[low]	85.27	Albite[low] + 8 H2O -> 1 Al[OH]4- + 6 H+ + 3 H2SiO4-2 + 1 Na+	Fluorite	10.96	Fluorite -> 1 Ca+2 + 2 F-
Antimocrandallite-exp	63.00	Antimocrandallite-exp + 8 H2O -> 3 Al[OH]4- + 1 Ca+2 + 3 H+ + 2 Sb[OH]6-	Li2_CaO_Al2O3_SiO2_8H2O[s]	22.69	Li2_CaO_Al2O3_SiO2_8H2O[s] -> 2 Al[OH]4- + 1 Ca+2 + 3 H2O + 1 H2SiO4-2 + 2 Li+
Arsenocrandallite-therm	95.56	Arsenocrandallite-therm + 6 H2O -> 3 Al[OH]4- + 2 AsO4-3 + 1 Ca+2 + 7 H+	Manganite	-25.27	Manganite + 3 H+ + 1 e- -> 2 H2O + 1 Mn+2
Austinite-therm	11.47	Austinite-therm + 1 H+ -> 1 AsO4-3 + 1 Ca+2 + 1 H2O + 1 Zn+2	Ni[OH]2[s]	-10.80	Ni[OH]2[s] + 2 H+ -> 2 H2O + 1 Ni+2
BaSrSO4[50%Ba]	8.221	BaSrSO4[50%Ba] -> 0.5 Ba+2 + 1 SO4-2 + 0.5 Sr+2	Pb[OH]2[C]	-8.150	Pb[OH]2[C] + 2 H+ -> 2 H2O + 1 Pb+2
beta-TCP	28.93	beta-TCP -> 3 Ca+2 + 2 PO4-3	Pb2VO7	0.9500	Pb2VO7 + 3 H+ -> 1.5 H2O + 1 Pb+2 + 1 VO2+
Boehmite	14.42	Boehmite + 2 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+
Ca[OH]Sb[OH]6[s]	2.000	Ca[OH]Sb[OH]6[s] + 1 H+ -> 1 Ca+2 + 1 H2O + 1 Sb[OH]6-	PbMoO4[c]	15.80	PbMoO4[c] -> 1 MoO4-2 + 1 Pb+2
Ca5[OH][AsO4]3[c]	26.13	Ca5[OH][AsO4]3[c] + 1 H+ -> 3 AsO4-3 + 5 Ca+2 + 1 H2O	Sn[OH]2[s]	1.447	Sn[OH]2[s] + 2 H+ -> 2 H2O + 1 Sn+2
CaCO3_BaCO3	22.00	CaCO3_BaCO3 -> 1 Ba+2 + 2 CO3-2 + 1 Ca+2	Strengite	48.00	Strengite + 2 H2O -> 1 Fe[OH]4- + 4 H+ + 1 PO4-3
CaCO3_SrCO3	19.85	CaCO3_SrCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2	Willemite	6.289	Willemite + 2 H+ -> 1 H2SiO4-2 + 2 Zn+2
CaMoO4[c]	7.940	CaMoO4[c] -> 1 Ca+2 + 1 MoO4-2	Zn3[SbO4]2[s]	3.000	Zn3[SbO4]2[s] + 4 H+ + 4 H2O -> 2 Sb[OH]6- + 3 Zn+2
CaPb2[PO4]2	40.76	CaPb2[PO4]2 -> 1 Ca+2 + 2 PO4-3 + 2 Pb+2			

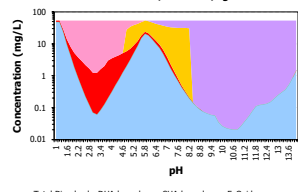
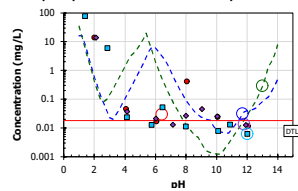


BIOMASS ASH WOOD PELLETS NL

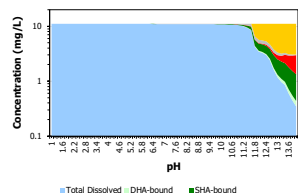
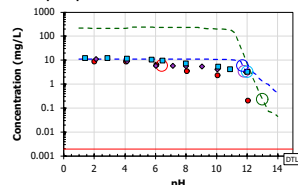
COMPARISON AND PARTITIONING



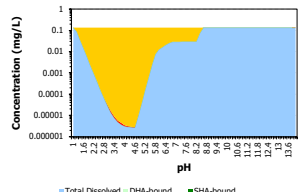
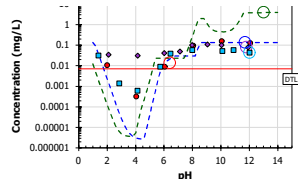
Mn Profile at L/S=11.7 L/kg



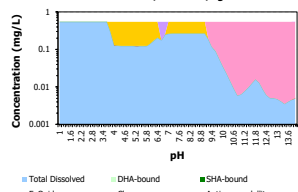
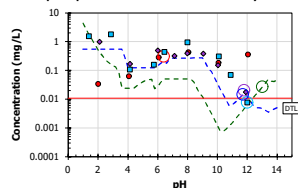
PO4 Profile at L/S=11.7 L/kg



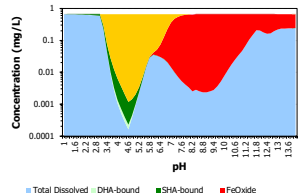
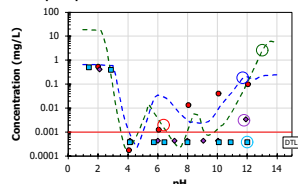
Sr Profile at L/S=11.7 L/kg



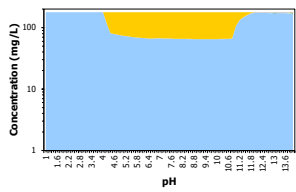
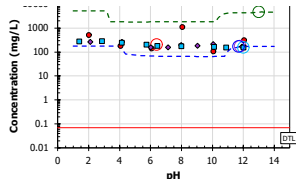
Mo Profile at L/S=11.7 L/kg



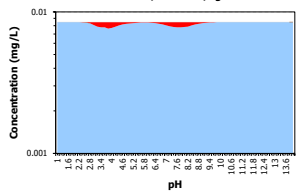
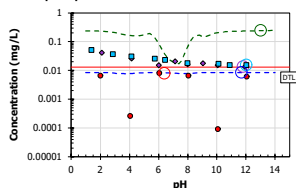
Sb Profile at L/S=11.7 L/kg



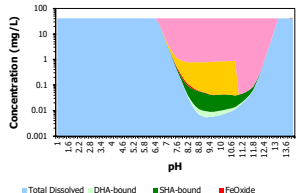
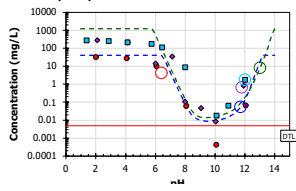
V Profile at L/S=11.7 L/kg



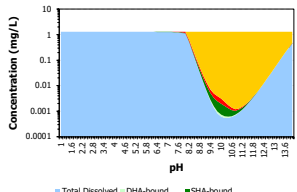
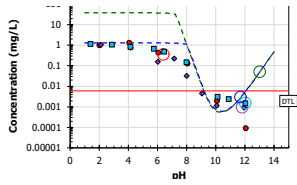
Na Profile at L/S=11.7 L/kg



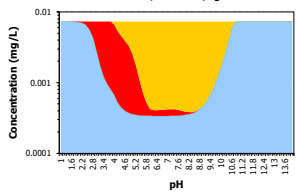
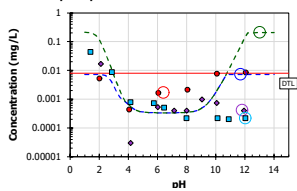
Se Profile at L/S=11.7 L/kg



Zn Profile at L/S=11.7 L/kg



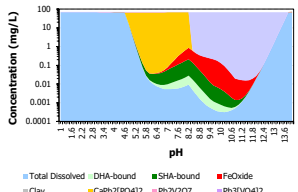
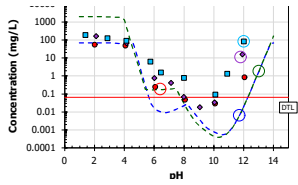
Ni Profile at L/S=11.7 L/kg



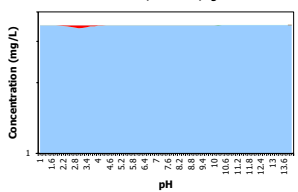
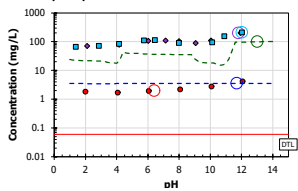
Sn Profile at L/S=11.7 L/kg



V Profile at L/S=11.7 L/kg



Pb Profile at L/S=11.7 L/kg



S Profile at L/S=11.7 L/kg



Sr Profile at L/S=11.7 L/kg

Model Comparison: residuals - Concentration

Name **Biomass wood ash 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

Residuals as log(model/sample)							
Fraction	6	5	4	3	2	1	Total Avg
pH	2.01	4.07	6.07	8.07	10.1	12.1	Deviation
Al	0.10	-0.05	0.38	0.97	1.54	1.07	0.36
As	-0.15	0.05	-0.55	-0.87	-0.45	0.96	0.25
Ba	0.10	0.14	-0.81	-1.10	-0.90	-1.60	0.38
Ca	0.10	0.08	0.08	0.05	0.19	0.01	0.04
Cd	0.10	0.08	0.35	0.76	0.38	0.54	0.18
Cl	-1.89	-1.79	0.07	0.12	0.00	-0.11	0.43
Co	0.10	0.18	0.50	0.66	-1.39	-0.87	0.31
Cr	-0.10	0.20	1.42	1.00	0.26	1.58	0.40
Cu	0.54	-0.01	1.82	0.79	1.35	2.40	0.57
Fe	0.00	-1.09	0.68	0.38	-0.78	0.57	0.28
B	0.48	0.54	0.57	0.12	1.01	0.64	0.25
Si	-0.66	-0.26	-0.22	-0.89	-0.59	0.74	0.25
K	0.28	0.35	0.36	0.38	0.38	0.44	0.15
Li	0.10	0.19	0.38	0.53	0.73	0.13	0.17
Mg	0.10	0.10	0.23	0.44	-0.07	-0.28	0.10
Mn	0.10	0.15	0.42	0.93	0.01	-1.46	0.30
Mo	-0.68	-2.01	0.20	-0.49	-0.06	0.03	0.37
Na	-0.45	-0.07	-0.31	-1.21	-0.20	-0.25	0.23
Ni	0.13	0.00	0.47	0.83	-0.44	1.82	0.35
Pb	0.10	0.13	-1.21	-0.29	-1.59	-1.50	0.42
PQ4	-	-	-	-	-	-	-
Sb	1.22	0.32	-0.26	-0.19	-0.78	-1.50	0.36
Se	0.11	1.47	0.02	0.08	1.97	0.14	0.41
Sn	0.14	0.15	-0.68	-0.78	-0.58	-0.06	0.20
S	0.28	0.31	0.26	0.20	0.11	-0.07	0.09
Sr	0.10	0.10	0.23	0.50	0.67	1.23	0.25
V	0.08	0.78	1.43	-0.65	-0.63	0.30	0.32
Zn	0.10	0.19	0.60	0.01	1.38	0.45	0.26
Avg Deviat	0.10	0.13	0.13	0.13	0.17	0.19	0.28

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