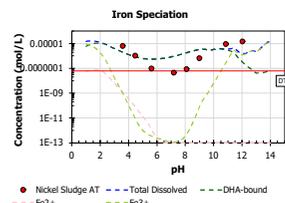
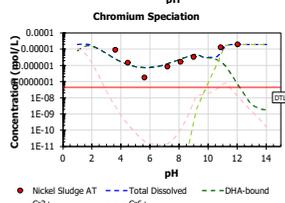
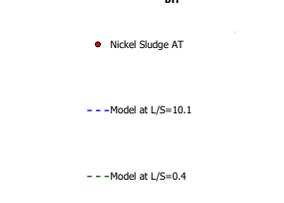
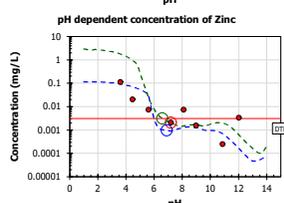
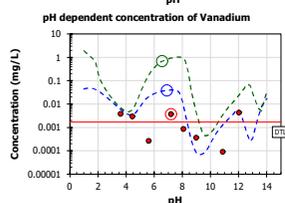
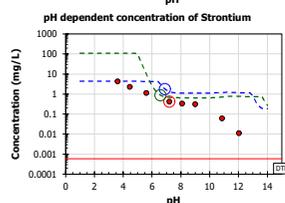
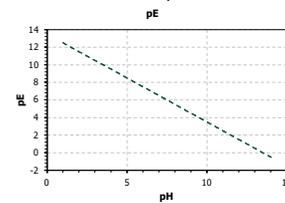
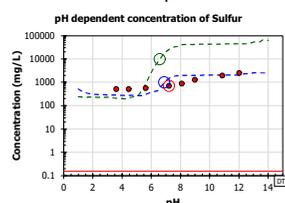
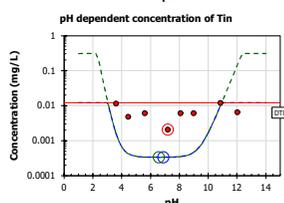
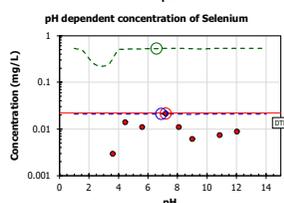
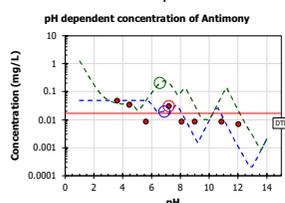
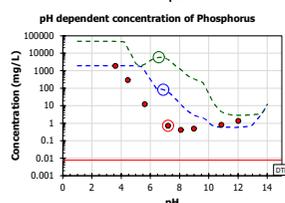
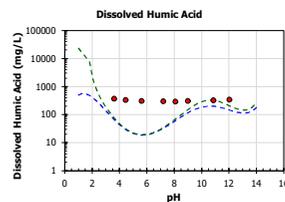
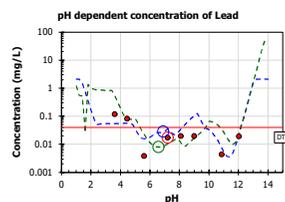
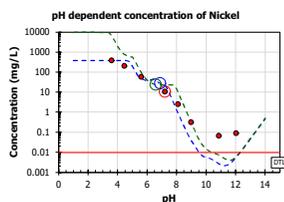
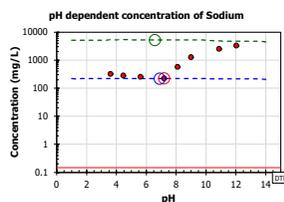
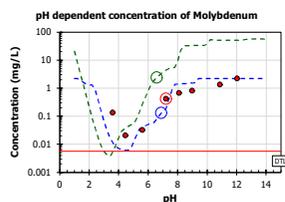
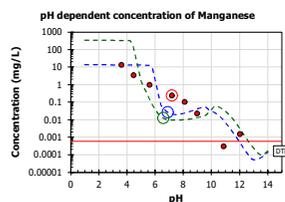
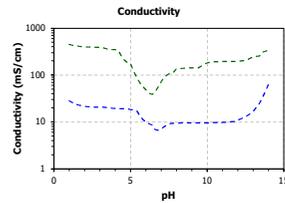
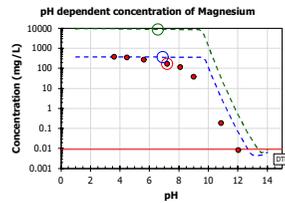
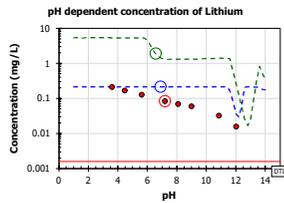
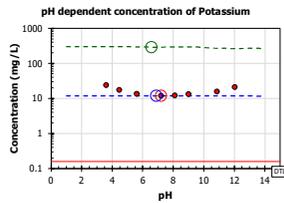
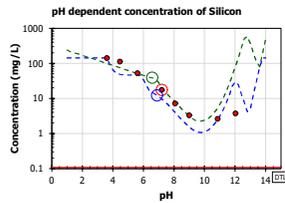
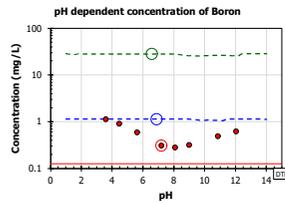
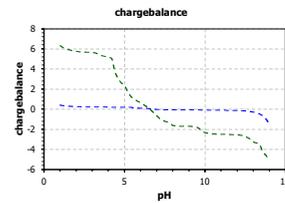
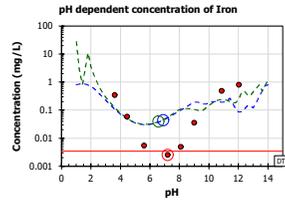
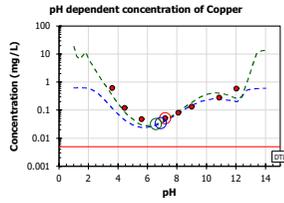
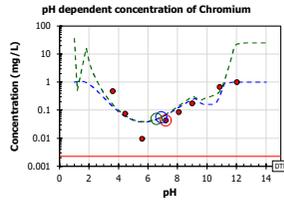
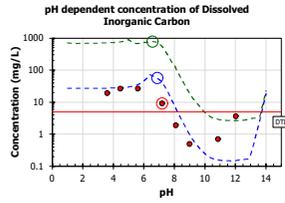
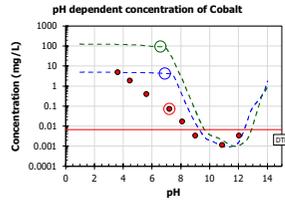
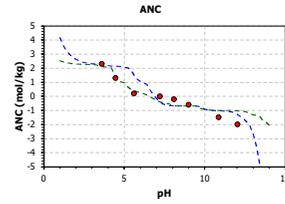
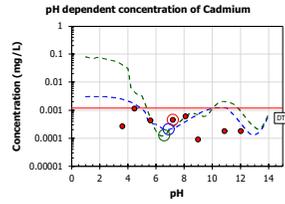
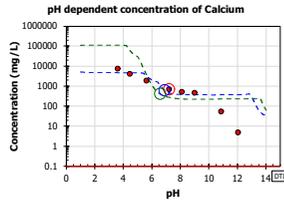
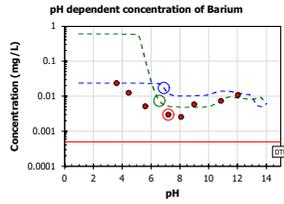
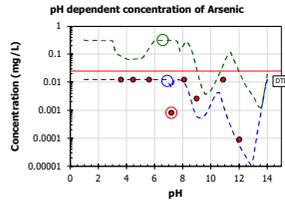
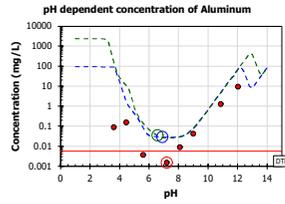


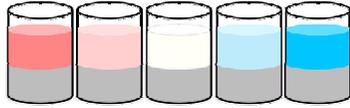
NICKEL SLUDGE AT

COMPARISON pH DEPENDENCE WITH MODEL



**Object** pH Dependent Leaching Test Model  
**Name** Nickel Sludge AT

**pH Dependent Leaching Test Scenario**



Lab Test

Extra L/S Simulation

**Lab Test**

**Model Parameters**

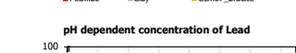
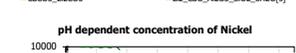
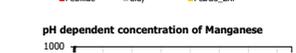
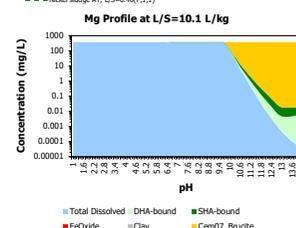
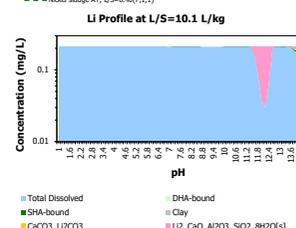
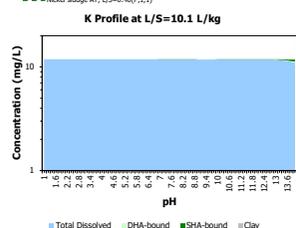
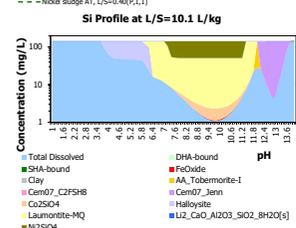
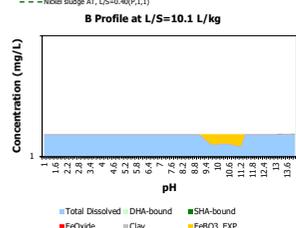
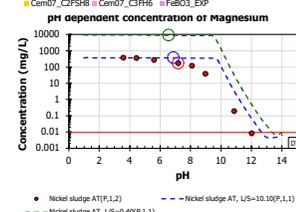
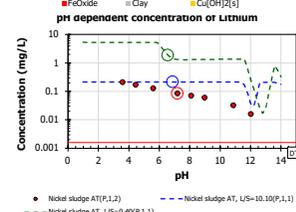
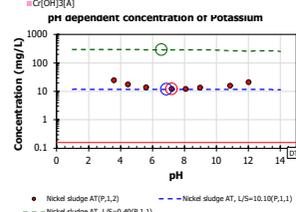
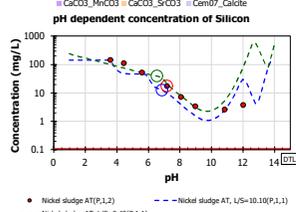
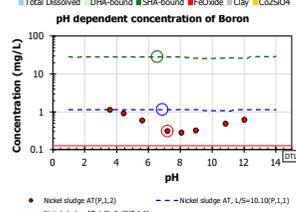
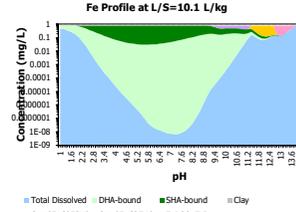
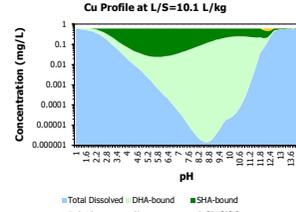
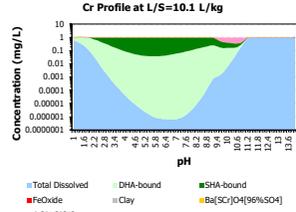
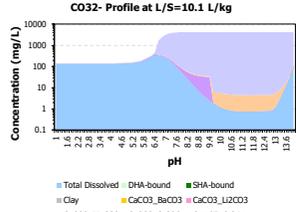
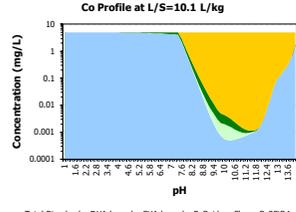
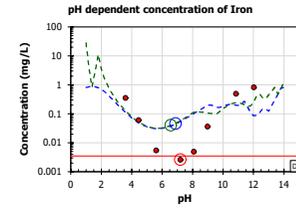
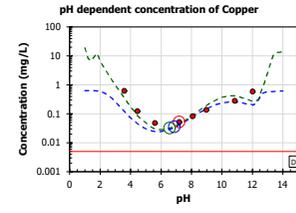
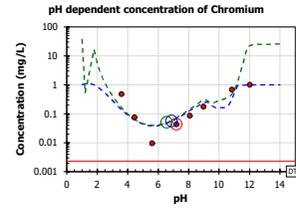
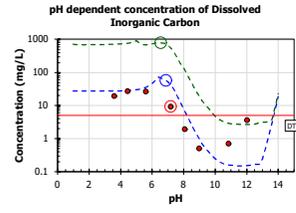
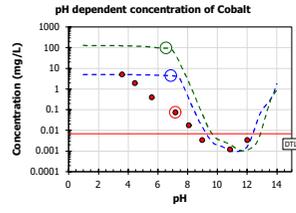
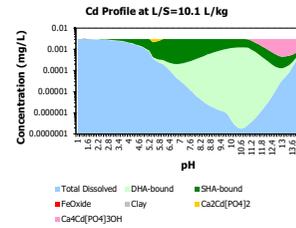
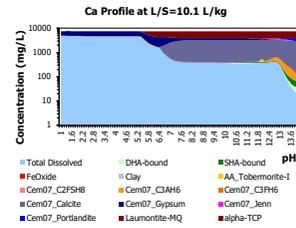
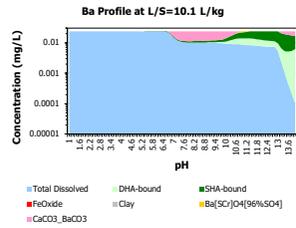
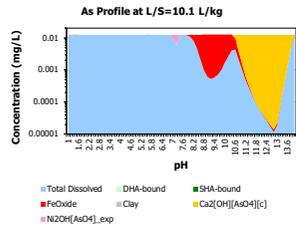
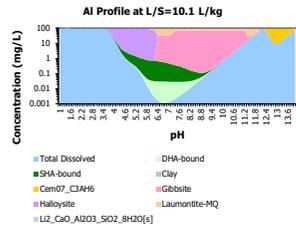
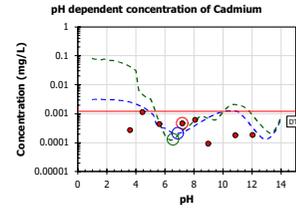
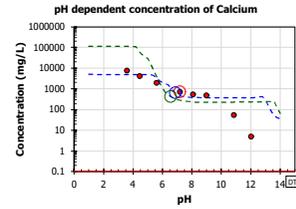
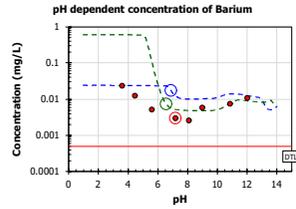
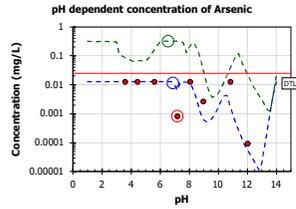
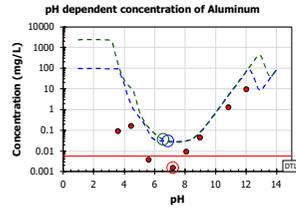
Entity	Unit	Default	Entity	mg/kg	Entity	mg/kg	Entity	mg/kg
c0		-4.514	Al	949.9	B	11.48	Sb	0.4925
c1		1.683	As	0.1274	Si	1444	Se	0.2131
c2		-0.9169	Ba	0.2413	Hg	2.006E-07	Sn	0.1225
c3		0.1653	Ca	7.683E+04	K	119.9	SO4	7.651E+04
c4		-0.01216	Cd	0.03145	Li	2.151	Sr	43.74
c5		0.0003163	Cl	3.545E-08	Mg	3789	V	0.4435
Clay	mg/kg	3000	Co	50.99	Mn	135.9	Zn	1.150
Hydrous Ferric Oxid	mg/kg	50.00	CO32-	4.336E+04	Mo	22.46		
L/S	L/kg	10.10	Cr	10.23	Na	2199		
pE		6.300	Cu	6.312	Ni	3886		
pH		7.200	F	5.000	Pb	21.20		
Solid Humic Acid	mg/kg	5000	Fe	8.204	PO4	5.886E+04		
Simulated Low L/S	L/kg	0.4000						

**Minerals**

Name	Log(K)	Reaction	Name	Log(K)	Reaction
AA_Tobermorite-l	23.86	AA_Tobermorite-l -> 2 Ca+2 + 0.8 H+ + 1.2 H2O + 2.4 H2SiO4-2	Cr(OH)3[A]	68.13	Cr(OH)3[A] + 1 H2O -> 1 CrO4-2 + 5 H+ + 3 e-
alpha-TCP	25.50	alpha-TCP -> 3 Ca+2 + 2 PO4-3	Cu(OH)2[s]	-8.640	Cu(OH)2[s] + 2 H+ -> 1 Cu+2 + 2 H2O
Arsenocrandallite-t	95.56	Arsenocrandallite-therm + 6 H2O -> 3 Al(OH)4- + 2 AsO4-3 + 1 Ca+2 + 7 H+	Fe2[SeO3]3:2l	183.8	Fe2[SeO3]3:2H2O + 7 H2O -> 2 Fe(OH)4- + 14 H+ + 3 SeO4-2 + 6 e-
Ba[Scr]O4[96%SO4]	9.790	Ba[Scr]O4[96%SO4] -> 1 Ba+2 + 0.04 CrO4-2 + 0.96 SO4-2	FeBO3_EXP	32.48	FeBO3_EXP + 2 H2O -> 1 Fe(OH)4- + 2 H+ + 1 H2BO3-
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-	Fluorite	10.96	Fluorite -> 1 Ca+2 + 2 F-
Ca2(OH)[AsO4][c]	4.000	Ca2(OH)[AsO4][c] + 1 H+ -> 1 AsO4-3 + 2 Ca+2 + 1 H2O	Gibbsite	14.71	Gibbsite + 1 H2O -> 1 Al(OH)4- + 1 H+
Ca2Cd[PO4]2	32.95	Ca2Cd[PO4]2 -> 2 Ca+2 + 1 Cd+2 + 2 PO4-3	Halloysite	79.12	Halloysite + 7 H2O -> 2 Al(OH)4- + 6 H+ + 2 H2SiO4-2
Ca4Cd[PO4]3OH	39.23	Ca4Cd[PO4]3OH + 1 H+ -> 4 Ca+2 + 1 Cd+2 + 1 H2O + 3 PO4-3	Laumontite-M	118.0	Laumontite-MQ + 8 H2O -> 2 Al(OH)4- + 1 Ca+2 + 8 H+ + 4 H2SiO4-2
CaCO3_BaCO3	22.00	CaCO3_BaCO3 -> 1 Ba+2 + 2 CO3-2 + 1 Ca+2	LDH_Zn_zc	20.91	LDH_Zn_zc + 3 H+ -> 1 Al(OH)4- + 1 CO3-2 + 3 H2O + 3 Zn+2
CaCO3_Li2CO3	21.30	CaCO3_Li2CO3 -> 2 CO3-2 + 1 Ca+2 + 2 Li+	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+
CaCO3_MnCO3	21.48	CaCO3_MnCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Mn+2	Magadiite	165.6	Magadiite + 9 H2O -> 13 H+ + 7 H2SiO4-2 + 1 Na+
CaCO3_SrCO3	19.85	CaCO3_SrCO3 -> 2 CO3-2 + 1 Ca+2 + 1 Sr+2	Manganite	-25.27	Manganite + 3 H+ + 1 e- -> 2 H2O + 1 Mn+2
CaMoO4[c]	7.940	CaMoO4[c] -> 1 Ca+2 + 1 MoO4-2	Ni(OH)2.Zn[OI]	-12.00	Ni(OH)2.Zn(OH)2-Exp + 4 H+ -> 4 H2O + 1 Ni+2 + 1 Zn+2
CaSb(OH)6[s]2_exp	19.41	CaSb(OH)6[s]2_exp -> 1 Ca+2 + 2 Sb(OH)6-	Ni(OH)2[s]	-10.80	Ni(OH)2[s] + 2 H+ -> 2 H2O + 1 Ni+2
Cem07_Brucite	-16.83	Cem07_Brucite + 2 H+ -> 2 H2O + 1 Mg+2	Ni2OH[AsO4]	14.00	Ni2OH[AsO4]_exp + 1 H+ -> 1 AsO4-3 + 1 H2O + 2 Ni+2
Cem07_C2FSH8	21.41	Cem07_C2FSH8 -> 2 Ca+2 + 2 Fe(OH)4- + 3 H2O + 1 H2SiO4-2	Ni2SiO4	5.498	Ni2SiO4 + 2 H+ -> 1 H2SiO4-2 + 2 Ni+2
Cem07_C3AH6	-35.14	Cem07_C3AH6 + 4 H+ -> 2 Al(OH)4- + 3 Ca+2 + 4 H2O	Ni3[PO4]2	31.30	Ni3[PO4]2 -> 3 Ni+2 + 2 PO4-3
Cem07_C3FH6	-30.82	Cem07_C3FH6 + 4 H+ -> 3 Ca+2 + 2 Fe(OH)4- + 4 H2O	NiMoO4	11.15	NiMoO4 -> 1 MoO4-2 + 1 Ni+2
Cem07_Calcite	8.485	Cem07_Calcite -> 1 CO3-2 + 1 Ca+2	Pb(OH)2[C]	-8.150	Pb(OH)2[C] + 2 H+ -> 2 H2O + 1 Pb+2
Cem07_Gypsum	4.583	Cem07_Gypsum -> 1 Ca+2 + 2 H2O + 1 SO4-2	PbMoO4[c]	15.80	PbMoO4[c] -> 1 MoO4-2 + 1 Pb+2
Cem07_Jenn	-7.799	Cem07_Jenn + 1.3333333 H+ -> 1.6666667 Ca+2 + 1.7666666 H2O + 1 H2SiO4-2	PbOH[Sb(OH)6]	12.00	PbOH[Sb(OH)6]_exp1 + 1 H+ -> 1 H2O + 1 Pb+2 + 1 Sb(OH)6-
Cem07_Portlandite	-22.79	Cem07_Portlandite + 2 H+ -> 1 Ca+2 + 2 H2O	Sn(OH)2[s]	1.447	Sn(OH)2[s] + 2 H+ -> 2 H2O + 1 Sn+2
Co2SiO4	5.289	Co2SiO4 + 2 H+ -> 2 Co+2 + 1 H2SiO4-2	Strengite	47.97	Strengite + 2 H2O -> 1 Fe(OH)4- + 4 H+ + 1 PO4-3

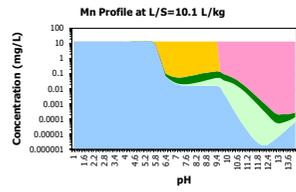
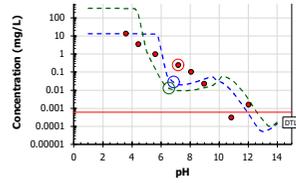
NICKEL SLUDGE AT

COMPARISON AND PARTITIONING

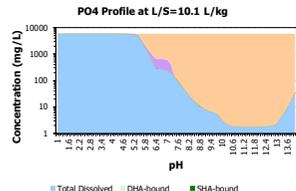
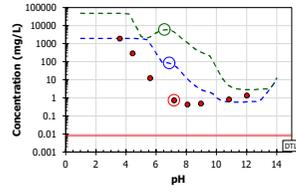


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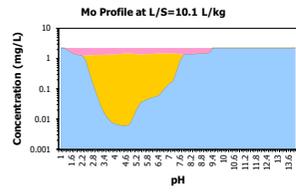
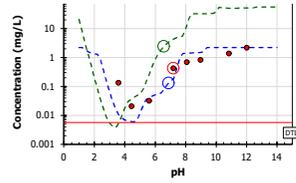
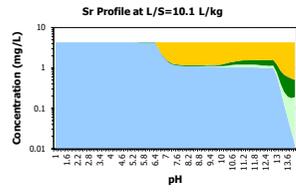
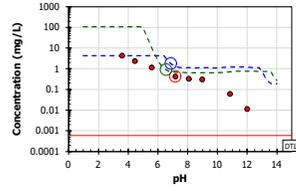
COMPARISON AND PARTITIONING



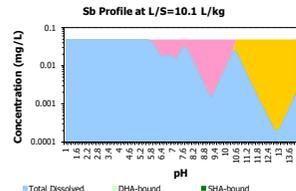
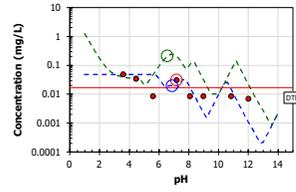
pH dependent concentration of Phosphorus



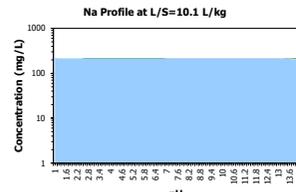
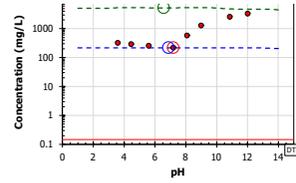
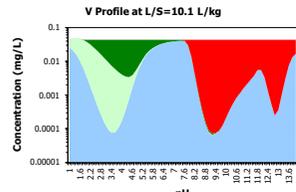
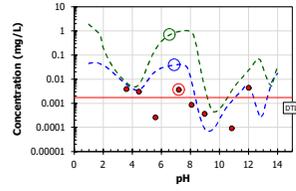
pH dependent concentration of Strontium



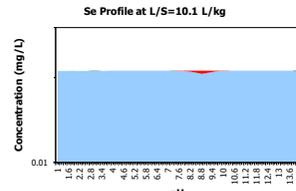
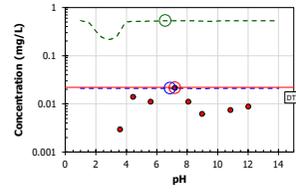
pH dependent concentration of Antimony



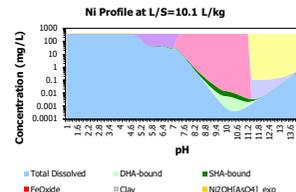
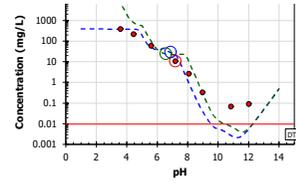
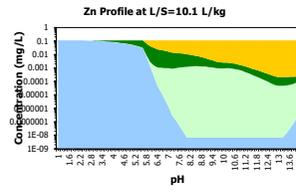
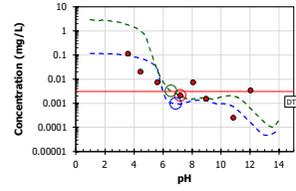
pH dependent concentration of Vanadium



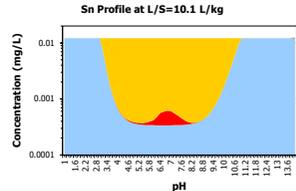
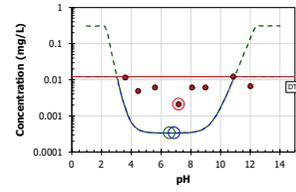
pH dependent concentration of Selenium



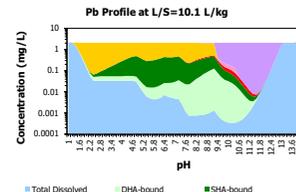
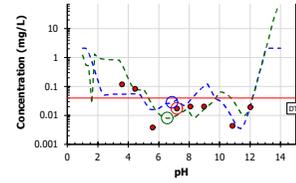
pH dependent concentration of Zinc



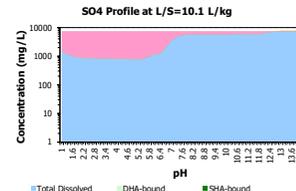
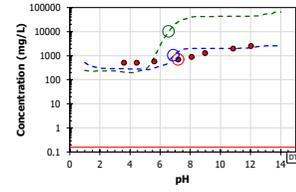
pH dependent concentration of Tin



pH dependent concentration of Sulfur



pH dependent concentration of Strontium



## Model Comparison: residuals - Concentration

Name **Ni Sludge AT**

### 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	8	7	6	5	4	3	2	1	Total Avg	
pH	3.60	4.46	5.61	7.20	8.09	8.99	10.9	12.0		Deviation
Al	3.00	1.03	1.33	1.28	0.53	0.26	0.59	0.88		0.48
As	0.01	0.01	0.01	0.89	-0.11	-0.63	-0.89	-0.23		0.18
Ba	0.00	0.28	0.66	0.61	0.60	0.25	0.28	0.07		0.15
Ca	-0.20	0.07	0.22	-0.18	-0.14	-0.09	0.85	1.89		0.26
Cd	0.98	0.17	-0.08	-0.30	-0.18	0.91	0.84	0.26		0.21
Cl	-	-	-	-	-	-	-	-		-
Co	0.00	0.41	1.08	1.76	1.28	0.64	-0.05	-0.30		0.32
CO32-	-	-	-	-	-	-	-	-		-
Cr	-0.54	-0.09	0.61	0.19	0.17	0.14	-0.31	0.00		0.12
Cu	-0.71	-0.43	-0.29	-0.10	-0.01	0.04	-0.04	-0.45		0.12
Fe	-0.50	-0.09	0.75	1.33	1.32	0.73	-0.39	-0.97		0.31
B	0.01	0.10	0.29	0.57	0.61	0.55	0.34	0.27		0.14
Si	0.01	-0.36	-0.05	-0.25	-0.26	-0.33	-0.01	0.89		0.13
Hg	-	-	-	-	-	-	-	-		-
K	-0.31	-0.17	-0.06	-0.01	-0.01	-0.05	-0.12	-0.25		0.06
Li	0.00	0.10	0.23	0.41	0.49	0.55	0.82	0.36		0.16
Mg	0.00	0.04	0.14	0.33	0.49	0.97	1.27	0.89		0.24
Mn	0.00	0.58	1.11	-1.05	-0.63	0.26	1.48	-0.52		0.30
Mo	-1.08	-0.52	0.11	-0.36	0.32	0.27	0.21	0.00		0.17
Na	-0.17	-0.12	-0.07	-0.01	-0.42	-0.77	-1.06	-1.18		0.23
Ni	0.00	0.26	-0.08	0.37	-0.45	-0.92	-1.40	-1.20		0.27
Pb	-0.33	-0.16	0.62	0.32	0.24	0.72	0.23	0.07		0.14
PO4	-	-	-	-	-	-	-	-		-
Sb	0.01	0.16	0.75	-0.29	0.30	-0.57	0.23	-0.82		0.17
Se	0.86	0.18	0.28	0.00	0.28	0.53	0.45	0.38		0.16
Sn	-0.85	-1.06	-1.25	-0.79	-1.24	-1.11	-0.09	0.27		0.33
SO4	-	-	-	-	-	-	-	-		-
Sr	0.00	0.27	0.58	0.50	0.53	0.57	1.30	2.03		0.33
V	0.20	0.08	1.91	1.05	0.93	-0.64	1.16	0.07		0.34
Zn	-0.07	0.55	0.61	-0.36	-0.77	-0.09	0.44	-1.32		0.23
<b>Avg Deviation</b>	0.15	0.08	0.14	0.14	0.12	0.12	0.15	0.17		0.22

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