Appendix B
December 2001 Off-Post Sampling
Camp Stanley Storage Activity

Sample ID	FO-J1				HS-2				HS-3				I10-2				I10-4			
Sample Date	12/20/01				12/19/01				12/19/01				12/18/01				12/19/01			
Sample Type	N				N				N				N				N			
Lab Sample ID	AP26876				AP26772				AP26773				AP26730				AP26778			
Method **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																				
Total Dissolved Solids																				
E160.2 (mg/L)																				
Suspended Solids																				
SW8260 (UG/L)																				
Acetone																				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-																				
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	**0.11**	F	1	0.6	**0.11**	F	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	**0.30**	F	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.38**	F	1	1.0	**0.40**	F	1	1.0	**0.38**	F	1	1.0	**0.21**	F	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	0.11	U	1	1.4	**0.16**	F	1	1.4	0.11	U	1	1.4	**0.16**	F	1	1.4	**0.12**	F	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.22**	F	1	1.0	0.14	U	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																				
pH																				

Sample ID	JW-9				JW-14				JW-14				LS-1			
Sample Date	12/19/01				12/20/01				12/20/01				12/19/01			
Sample Type	N				N				FD				N			
Lab Sample ID	AP26779				AP26877				AP26878				AP26774			
Method **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																
Total Dissolved Solids																
E160.2 (mg/L)																
Suspended Solids																
SW8260 (UG/L)																
Acetone																
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	M	1	1.1	0.36	M	1	1.1	0.36	U	1	1.1
Butanone, 2-																
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	**1.83**		1	0.3	**1.83**		1	0.3	0.27	F	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	**0.33**	F	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	M	1	3.5	0.53	M	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	**0.93**	F	1	1.0	**0.40**	F	1	1.0	**0.22**	F	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	**0.65**	F	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.26**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																
pH																

Sample ID	LS-2				LS-2				LS-2				LS-2			
Sample Date	09/27/01				10/04/01				11/14/01				12/19/01			
Sample Type	N				N				N				N			
Lab Sample ID	AP22601				AP22869				AP25267				AP26777			
Method **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																
Total Dissolved Solids	450		1	15.0												
E160.2 (mg/L)																
Suspended Solids	0		1	10.0												
SW8260 (UG/L)																
Acetone	1.45	U	1	5.0	1.45	U	1	5.0								
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	**0.31**	F	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-	1.65	U	1	5.0	1.65	U	1	5.0								
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	**0.2**	F	1	0.3	**0.65**		1	0.3	**0.15**	F	1	0.3	**0.10**	F	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	**0.22**	F	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.1	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.1	U	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.2	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	**4.6**		1	1.4	**3.2**		1	1.4	**2.7**		1	1.4	**3.45**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**0.17**	F	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.20**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																
pH	7.		1													

Tables present all laboratory results for analytes detected above the method detection limit.
Results from all laboratory analysis are presented in Appendix B.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 36485, 36538, 36527, 36600, 36730, 36936, 37183, 37184, 37198, 37199, 37209, 38228.

Abbreviations/Notes:
Bolded and shaded samples indicate results greater than MCL standards
-- No risk reduction standard or background level available
FD1 Field Duplicate
MDL Method Detection Limit
N1 Environmental Sample
RL Reporting Limit
SQL Sample Quantitation Limit
MCL Maximum Contamination Level

Data Qualifiers:
F- The analyte was positively identified but the associated numerical value is below the RL.
J - The analyte was positively identified, the quantitation is an estimation.
U - The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.
R- The data are unusable due to deficiencies in the ability to analyze the sample and meet QC criteria.
M- Matrix Effect Present

Bold Value > or = MCL
Bold MCL > Value > or = RL
Bold RL > Value > MDL

Appendix B
December 2001 Off-Post Sampling
Camp Stanley Storage Activity

Sample ID	LS-2-LS-3/EP				LS-2/LS-3 EP				LS-3				LS-3				LS-3				LS-3			
Sample Date	10/04/01				11/14/01				09/27/01				10/04/01				11/14/01				12/19/01			
Sample Type	N				N				N				N				N				N			
Lab Sample ID	AP22871				AP25269				AP22602				AP22870				AP25268				AP26776			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																								
Total Dissolved Solids									414		1	15.0												
E160.2 (mg/L)																								
Suspended Solids									0		1	10.0												
SW8260 (UG/L)																								
Acetone	**2.2**	F	1	5.0					1.45	U	1	5.0	1.45	U	1	5.0	Result	Flag	Dilution	SQL				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	**6.2**		1	0.8	**2.8**		1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	**4.5**		1	1.2	**2.9**		1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-	1.65	U	1	5.0					1.65	U	1	5.0	1.65	U	1	5.0								
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	**3.1**		1	0.3	**0.92**		1	0.3	**0.13**	F	1	0.3	**0.13**	F	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	**9.4**		1	0.5	**4.9**		1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.1	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.1	U	1	0.6	**0.19**	F	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	**0.29**	F	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.2	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	**0.54**	F	1	1.4	**0.83**	F	1	1.4	**4.8**		1	1.4	**3.7**		1	1.4	**1.3**	F	1	1.4	**4.29**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**0.18**	F	1	1.0	**0.16**	F	1	1.0	**0.43**	F	1	1.0	**0.45**	F	1	1.0	**0.8**	F	1	1.0	**0.27**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																								
pH									7		1													

Sample ID	LS-4				LS-5				LS-5				LS-6				LS-6 PRE				LS-6 POST				LS-7			
Sample Date	12/19/01				10/25/01				12/19/01				09/27/01				12/18/2001				12/18/2001				09/27/01			
Sample Type	N				N				N				N				N				N				N			
Lab Sample ID	AP26775				AP24132				AP26780				AP22603				AP26723				AP26724				AP22604			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																												
Total Dissolved Solids													450		1	15.0									420		1	15.0
E160.2 (mg/L)																												
Suspended Solids													0		1	10.0									0		1	10.0
SW8260 (UG/L)																												
Acetone													1.45	U	1	5.0									1.45	U	1	5.0
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-													1.65	U	1	5.0									1.65	U	1	5.0
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.09	F	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.20**	F	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	**0.33**	F	1	1.0	**0.29**	F	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4
Tetrachloroethene	**0.23**	F	1	1.4	**0.15**	F	1	1.4	0.11	U	1	1.4	**10.**		1	1.4	**1.97**		1	1.4	0.11	U	1	1.4	**3.3**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.0	**0.27**	F	1	1.0	**0.32**	F	1	1.0	**0.19**	F	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.34**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																												
pH													7		1										7			

Tables present all laboratory results for analytes detected above the method detection limit.
Results from all laboratory analysis are presented in Appendix B.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 37197, 37198, 37183, 37184, 37209, and 37228.

Abbreviations/Notes:
Bolded and shaded samples indicate results greater than MCL standards
-- No risk reduction standard or background level available
FD1 Field Duplicate
MDL Method Detection Limit
N1 Environmental Sample
RL Reporting Limit
SQL Sample Quantitation Limit
MCL Maximum Contamination Level

Data Qualifiers:
F- The analyte was positively identified but the associated numerical value is below the RL.
J - The analyte was positively identified, the quantitation is an estimation.
U - The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.
R- The data are unusable due to deficiencies in the ability to analyze the sample and meet QC criteria.
M- Matrix Effect Present

Bold Value > or = MCL
Bold MCL > Value > or = RL
Bold RL > Value > MDL

Appendix B
December 2001 Off-Post Sampling
Camp Stanley Storage Activity

Sample ID	LS-7 PRE				LS-7 POST				OFR-1				OFR-3			
Sample Date	12/18/2001				12/18/2001				12/20/01				10/25/01			
Sample Type	N				N				N				N			
Lab Sample ID	AP26725				AP26726				AP26879				AP24133			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																
Total Dissolved Solids																
E160.2 (mg/L)																
Suspended Solids																
SW8260 (UG/L)																
Acetone																
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-																
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	**0.75**	F	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.39**	F	1	1.0	**0.30**	F	1	1.0	**0.63**	F	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4
Tetrachloroethene	**1.40**		1	1.4	0.11	U	1	1.4	**0.42**	F	1	1.4	**4.2**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**2.2**		1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																
pH																

Sample ID	OFR-3				RFR-9				RFR-10			
Sample Date	12/18/01				12/20/01				10/03/01			
Sample Type	N				N				N			
Lab Sample ID	AP26728				AP26880				AP22809			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)												
Total Dissolved Solids												
E160.2 (mg/L)												
Suspended Solids												
SW8260 (UG/L)												
Acetone									1.45	U	1	5.0
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-									1.65	U	1	5.0
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	**0.39**	F	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	**0.49**	F	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.26**	F	1	1.0	**0.52**	F	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4
Tetrachloroethene	**1.90**		1	1.4	0.11	U	1	1.4	**19.**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**2.20**		1	1.0	0.14	U	1	1.0	**8.7**		1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)												
pH												

Sample ID	RFR-10 TAP-H				RFR-10 TAP-T				RFR-10 PRE			
Sample Date	10/03/01				10/03/01				10/12/01			
Sample Type	N				N				N			
Lab Sample ID	AP22811				AP22810				AP23210			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)												
Total Dissolved Solids												
E160.2 (mg/L)												
Suspended Solids												
SW8260 (UG/L)												
Acetone	**1.7**	F	1	5.0	**1.6**	F	1	5.0				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-	1.65	U	1	5.0	1.65	U	1	5.0				
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	**0.07**	F	1	0.3	**0.07**	F	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.1	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.1	U	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	**0.48**	F	1	1.2	**0.46**	F	1	1.2	**0.49**	F	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.1	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.2	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4
Tetrachloroethene	**17.**		1	1.4	**19.**		1	1.4	**20.89**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**7.5**		1	1.0	**9.4**		1	1.0	**8.21**		1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)												
pH												

Sample ID	RFR-10 PRE				RFR-10 POST-H				RFR-10 POST-T			
Sample Date	12/17/01				10/12/01				10/12/01			
Sample Type	N				N				N			
Lab Sample ID	AP26645				AP23211				AP23214			
Method / **Analyte**	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)												
Total Dissolved Solids												
E160.2 (mg/L)												
Suspended Solids												
SW8260 (UG/L)												
Acetone												
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-												
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	**0.26**	F	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.30**	F	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4
Tetrachloroethene	**9.02**		1	1.4	0.11	U	1	1.4	0.11	U	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**5.85**		1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)												
pH												

Tables present all laboratory results for analytes detected above the method detection limit.
Results from all laboratory analysis are presented in Appendix B.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 37197, 37198, 37183, 37184, 37209, and 37228.

Abbreviations/Notes:
Bolded and shaded samples indicate results greater than MCL standards
-- No risk reduction standard or background level available
FD1 Field Duplicate
MDL Method Detection Limit
N1 Environmental Sample
RL Reporting Limit
SQL Sample Quantitation Limit
MCL Maximum Contamination Level

Data Qualifiers:
F- The analyte was positively identified but the associated numerical value is below the RL.
J - The analyte was positively identified, the quantitation is an estimation.
U - The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.
R- The data are unusable due to deficiencies in the ability to analyze the sample and meet QC criteria.
M- Matrix Effect Present

Bold Value > or = MCL
Bold MCL > Value > or = RL
Bold RL > Value > MDL

Appendix B
December 2001 Off-Post Sampling
Camp Stanley Storage Activity

Sample ID	RFR-10 POST				RFR-11				RFR-11				RFR-11 PRE				RFR-11 POST				RFR-11 POST				RFR-12			
Sample Date	12/17/01				10/04/01				10/12/01				12/17/01				10/25/01				12/17/01				12/18/01			
Sample Type	N				N				N				N				N				N				N			
Lab Sample ID	AP26646				AP22872				AP23212				AP26647				AP24131				AP26648				AP26729			
Method																												
Analyte	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL	Result	Flag	Dilution	SQL
E160.1 (mg/L)																												
Total Dissolved Solids																												
E160.2 (mg/L)																												
Suspended Solids																												
SW8260 (UG/L)																												
Acetone					1.45	U	1	5.0																				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butanone, 2-					1.65	U	1	5.0																				
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	**0.14**	F	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.1	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.1	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	**0.12**	F	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	**0.96**	F	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	**0.24**	F	1	1.0	0.19	U	1	1.0	**0.22**	F	1	1.0	**0.38**	F	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.1	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.2	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	0.11	U	1	1.4	**16.**		1	1.4	**16.73**		1	1.4	**12.44**		1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.0	**0.35**	F	1	1.0	**0.58**	F	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.15**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1
SW9045 (pH units)																												
pH																												

Bold	Value > or = MCL
Bold	MCL > Value > or = RL
Bold	RL > Value > MDL

Tables present all laboratory results for analytes detected above the method detection limit.
Results from all laboratory analysis are presented in Appendix B.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 37197, 37198, 37183, 37184, 37209, and 37228.

Abbreviations/Notes:
Bolded and shaded samples indicate results greater than MCL standards
-- No risk reduction standard or background level available
FD1 Field Duplicate
MDL Method Detection Limit
N1 Environmental Sample
RL Reporting Limit
SQL Sample Quantitation Limit
MCL Maximum Contamination Level

Data Qualifiers:
F- The analyte was positively identified but the associated numerical value is below the RL.
J - The analyte was positively identified, the quantitation is an estimation.
U - The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.
R- The data are unusable due to deficiencies in the ability to analyze the sample and meet QC criteria.
M- Matrix Effect Present