APPENDIX A—SUMMARY OF ANALYTICAL RESULTS

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW6010	NONE	ARSENIC	=	4.6	2	1	MG/KG			
				BARIUM	=	111	0.4	1	MG/KG			
				CADMIUM	=	0.27	0.2	1	MG/KG			
				CHROMIUM	=	10.6	1	1	MG/KG			
				COPPER	=	4.4	2	1	MG/KG			
				NICKEL	TR	0.8	2	1	MG/KG	F	B	
				LEAD	=	4.9	2	1	MG/KG			
				ZINC	=	24.5	2	1	MG/KG			
		SW7471		MERCURY	ND	0	0.1	1	MG/KG	U	U	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW8260	NONE	DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	TR	0.002	0.02	1	MG/KG	F	J	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag

Flag 2 = Laboratory data qualifier

Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW8260	NONE	n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW8270	NONE	4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW8270	NONE	2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
BORROW_08282002	8/28/2002	SW8270	NONE	2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CS-B3-ET-1	9/24/2002	SW6010	SW1311	SILVER	TR	0.003	0.01	1	MG/L	F	B	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.3217	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0382	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.019	0.02	1	MG/L	F	B	
				LEAD	TR	0.031	0.05	1	MG/L	F		
				ANTIMONY	TR	0.0326	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	M	U	
		SW8260		BENZENE	ND	0	0.5	20	MG/L	U	U	
				BENZENE	ND	0	10	400	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.5	20	MG/L	U	U	
				CHLOROBENZENE	ND	0	10	400	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.5	20	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	10	400	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.5	20	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	10	400	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.5	20	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	10	400	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	1	20	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	20	400	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	=	34.2	0.5	20	MG/L		E	
				TETRACHLOROETHYLENE(PCE)	=	45.1	10	400	MG/L		D	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-1	9/24/2002	SW8260	SW1311	TRICHLOROETHYLENE (TCE)	=	2.2	0.5	20	MG/L			
				TRICHLOROETHYLENE (TCE)	TR	4.5	10	400	MG/L	F	DJ	
				CHLOROFORM	ND	0	0.5	20	MG/L	U	U	
				CHLOROFORM	ND	0	10	400	MG/L	U	U	
				VINYL CHLORIDE	ND	0	1	20	MG/L	U	U	
				VINYL CHLORIDE	ND	0	20	400	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	=	5870	50	1	MG/KG			
CS-B3-ET-2		SW6010	SW1311	SILVER	TR	0.002	0.01	1	MG/L	F	B	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.5922	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0696	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.018	0.02	1	MG/L	F	B	
				LEAD	=	0.057	0.05	1	MG/L			
				ANTIMONY	TR	0.0186	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.0049	0.025	1	MG/L	M	J	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.2	0.025	1	MG/L	M		
				TRICHLOROETHYLENE (TCE)	=	0.029	0.025	1	MG/L			
				CHLOROFORM	TR	0.0085	0.025	1	MG/L	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-2	9/24/2002	SW8260	SW1311	VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	=	391	50	1	MG/KG			
CS-B3-ET-20	10/2/2002	SW8260		TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
CS-B3-ET-21				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
CS-B3-ET-22		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-22	10/2/2002	SW8270	NONE	4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-22	10/2/2002	SW8270	NONE	NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CS-B3-ET-23		SW8260		TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
CS-B3-ET-24	10/8/2002			TETRACHLOROETHYLENE(PCE)	=	0.56	0.5	1	MG/KG		D	
				TETRACHLOROETHYLENE(PCE)	=	0.103	0.005	1	MG/KG		E	J
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.5	1	MG/KG	U	U	
CS-B3-ET-25	10/17/2002			TETRACHLOROETHYLENE(PCE)	TR	0.0015	0.005	1	MG/KG	F	J	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
CS-B3-ET-DRUM	9/24/2002	SW6010		ARSENIC	TR	0.004	0.01	1	MG/L	F	B	
				BARIUM	=	0.0999	0.002	1	MG/L			
				CADMIUM	=	0.0079	0.001	1	MG/L			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW6010	NONE	CHROMIUM	TR	0.002	0.01	1	MG/L	F	B	
				COPPER	=	0.055	0.01	1	MG/L			
				NICKEL	=	0.028	0.01	1	MG/L			
				LEAD	=	0.041	0.01	1	MG/L			
				ZINC	=	0.152	0.02	1	MG/L			
		SW8260		ACETONE	ND	0	2000	200	UG/L	R	UR	
				ACETONE	ND	0	20000	2000	UG/L	R	UR	
				BROMODICHLOROMETHANE	ND	0	100	200	UG/L	U	U	
				BROMODICHLOROMETHANE	ND	0	1000	2000	UG/L	U	U	
				BROMOBENZENE	ND	0	200	200	UG/L	U	U	
				BROMOBENZENE	ND	0	2000	2000	UG/L	U	U	
				BROMOCHLOROMETHANE	ND	0	200	200	UG/L	U	U	
				BROMOCHLOROMETHANE	ND	0	2000	2000	UG/L	U	U	
				BROMOMETHANE	ND	0	600	200	UG/L	U	U	
				BROMOMETHANE	ND	0	6000	2000	UG/L	U	U	
				n-BUTYLBENZENE	ND	0	200	200	UG/L	U	U	
				n-BUTYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				SEC-BUTYLBENZENE	ND	0	200	200	UG/L	U	U	
				SEC-BUTYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				t-BUTYLBENZENE	ND	0	200	200	UG/L	U	U	
				t-BUTYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				BENZENE	ND	0	80	200	UG/L	U	U	
				BENZENE	ND	0	800	2000	UG/L	U	U	
				TOLUENE	ND	0	200	200	UG/L	U	U	
				TOLUENE	ND	0	2000	2000	UG/L	U	U	
				CHLOROBENZENE	ND	0	100	200	UG/L	U	U	
				CHLOROBENZENE	ND	0	1000	2000	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8260	NONE	2-CHLOROTOLUENE	ND	0	200	200	UG/L	U	U	
				2-CHLOROTOLUENE	ND	0	2000	2000	UG/L	U	U	
				4-CHLOROTOLUENE	ND	0	200	200	UG/L	U	U	
				4-CHLOROTOLUENE	ND	0	2000	2000	UG/L	U	U	
				CHLOROETHANE	ND	0	200	200	UG/L	U	U	
				CHLOROETHANE	ND	0	2000	2000	UG/L	U	U	
				1-CHLOROHEXANE	ND	0	200	200	UG/L	U	U	
				1-CHLOROHEXANE	ND	0	2000	2000	UG/L	U	U	
				CHLOROMETHANE	=	316.8	200	200	UG/L	B	B	
				CHLOROMETHANE	=	3144	2000	2000	UG/L		D	
				CARBON TETRACHLORIDE	ND	0	200	200	UG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	2000	2000	UG/L	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	200	200	UG/L	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	2000	2000	UG/L	U	U	
				DIBROMOCHLOROMETHANE	ND	0	100	200	UG/L	U	U	
				DIBROMOCHLOROMETHANE	ND	0	1000	2000	UG/L	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	400	200	UG/L	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	4000	2000	UG/L	U	U	
				DIBROMOMETHANE	ND	0	200	200	UG/L	U	U	
				DIBROMOMETHANE	ND	0	2000	2000	UG/L	U	U	
				1,1-DICHLOROETHANE	ND	0	200	200	UG/L	U	U	
				1,1-DICHLOROETHANE	ND	0	2000	2000	UG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	100	200	UG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	1000	2000	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	200	200	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	200	200	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8260	NONE	1,3-DICHLOROBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	100	200	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	1000	2000	UG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	200	200	UG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	2000	2000	UG/L	U	U	
				cis-1,2-DICHLOROETHYLENE	TR	170.6	200	200	UG/L	F	J	
				cis-1,2-DICHLOROETHYLENE	ND	0	2000	2000	UG/L	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	200	200	UG/L	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	2000	2000	UG/L	U	U	
				1,1-DICHLOROPROPENE	ND	0	200	200	UG/L	U	U	
				1,1-DICHLOROPROPENE	ND	0	2000	2000	UG/L	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	100	200	UG/L	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	1000	2000	UG/L	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	200	200	UG/L	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	2000	2000	UG/L	U	U	
				1,2-DICHLOROPROPANE	ND	0	200	200	UG/L	U	U	
				1,2-DICHLOROPROPANE	ND	0	2000	2000	UG/L	U	U	
				1,3-DICHLOROPROPANE	ND	0	80	200	UG/L	U	U	
				1,3-DICHLOROPROPANE	ND	0	800	2000	UG/L	U	U	
				2,2-DICHLOROPROPANE	ND	0	200	200	UG/L	U	U	
				2,2-DICHLOROPROPANE	ND	0	2000	2000	UG/L	U	U	
				ETHYLBENZENE	ND	0	200	200	UG/L	U	U	
				ETHYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	200	200	UG/L	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	2000	2000	UG/L	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	200	200	UG/L	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	2000	2000	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8260	NONE	DICHLORODIFLUOROMETHANE	ND	0	200	200	UG/L	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	2000	2000	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	120	200	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	1200	2000	UG/L	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	200	200	UG/L	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	2000	2000	UG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	2000	200	UG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	20000	2000	UG/L	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	2000	200	UG/L	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	20000	2000	UG/L	U	U	
				METHYLENE CHLORIDE	ND	0	200	200	UG/L	U	U	
				METHYLENE CHLORIDE	ND	0	2000	2000	UG/L	U	U	
				NAPHTHALENE	ND	0	200	200	UG/L	U	U	
				NAPHTHALENE	ND	0	2000	2000	UG/L	U	U	
				n-PROPYLBENZENE	ND	0	200	200	UG/L	U	U	
				n-PROPYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	100	200	UG/L	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	1000	2000	UG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	=	25630	200	200	UG/L		E	
				TETRACHLOROETHYLENE(PCE)	=	20500	2000	2000	UG/L		D	
				STYRENE	ND	0	200	200	UG/L	U	U	
				STYRENE	ND	0	2000	2000	UG/L	U	U	
				BROMOFORM	ND	0	200	200	UG/L	U	U	
				BROMOFORM	ND	0	2000	2000	UG/L	U	U	
				tert-BUTYL METHYL ETHER	ND	0	1000	200	UG/L	U	U	
				tert-BUTYL METHYL ETHER	ND	0	10000	2000	UG/L	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	100	200	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8260	NONE	1,1,1,2-TETRACHLOROETHANE	ND	0	1000	2000	UG/L	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	200	200	UG/L	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	2000	2000	UG/L	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	200	200	UG/L	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	2000	2000	UG/L	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	200	200	UG/L	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	200	200	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	2000	2000	UG/L	U	U	
				TRICHLOROETHYLENE (TCE)	=	1744	200	200	UG/L			
				TRICHLOROETHYLENE (TCE)	=	3244	2000	2000	UG/L		D	
				CHLOROFORM	ND	0	60	200	UG/L	U	U	
				CHLOROFORM	ND	0	600	2000	UG/L	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	200	200	UG/L	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	2000	2000	UG/L	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	200	200	UG/L	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	2000	2000	UG/L	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	200	200	UG/L	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	2000	2000	UG/L	U	U	
				VINYL CHLORIDE	ND	0	200	200	UG/L	U	U	
				VINYL CHLORIDE	ND	0	2000	2000	UG/L	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	400	200	UG/L	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	4000	2000	UG/L	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	200	200	UG/L	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	2000	2000	UG/L	U	U	
		SW8270		ACENAPHTHENE	ND	0	10	1	UG/L	U	U	
				ACENAPHTHENE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8270	NONE	ACENAPHTHYLENE	ND	0	10	1	UG/L	U	U	
				ACENAPHTHYLENE	ND	0	10	1	UG/L	U	U	
				ANTHRACENE	ND	0	10	1	UG/L	U	U	
				ANTHRACENE	ND	0	10	1	UG/L	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	10	1	UG/L	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	TR	2.3	10	1	UG/L	F	J	
				bis(2-ETHYLHEXYL) PHTHALATE	TR	2.66	10	1	UG/L	F	J	
				4-BROMOPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				BENZO(a)ANTHRACENE	ND	0	10	1	UG/L	U	U	
				BENZO(a)ANTHRACENE	ND	0	10	1	UG/L	U	U	
				BENZOIC ACID	ND	0	100	1	UG/L	R	UR	
				BENZOIC ACID	ND	0	100	1	UG/L	R	UR	
				BENZO(a)PYRENE	ND	0	10	1	UG/L	M	U	
				BENZO(a)PYRENE	ND	0	10	1	UG/L	M	U	
				BENZO(b)FLUORANTHENE	ND	0	10	1	UG/L	M	U	
				BENZO(b)FLUORANTHENE	ND	0	10	1	UG/L	M	U	
				BENZO(g,h,i)PERYLENE	ND	0	10	1	UG/L	M	U	
				BENZO(g,h,i)PERYLENE	ND	0	10	1	UG/L	M	U	
				BENZO(k)FLUORANTHENE	ND	0	10	1	UG/L	M	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8270	NONE	BENZO(k)FLUORANTHENE	ND	0	10	1	UG/L	M	U	
				BENZYL ALCOHOL	ND	0	20	1	UG/L	U	U	
				BENZYL ALCOHOL	ND	0	20	1	UG/L	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	20	1	UG/L	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	20	1	UG/L	U	U	
				CHRYSENE	ND	0	10	1	UG/L	U	U	
				CHRYSENE	ND	0	10	1	UG/L	U	U	
				4-CHLOROANILINE	ND	0	20	1	UG/L	U	U	
				4-CHLOROANILINE	ND	0	20	1	UG/L	U	U	
				2-CHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				2-CHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				2-CHLORONAPHTHALENE	ND	0	10	1	UG/L	U	U	
				2-CHLORONAPHTHALENE	ND	0	10	1	UG/L	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	10	1	UG/L	M	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	10	1	UG/L	M	U	
				DIBENZOFURAN	ND	0	10	1	UG/L	U	U	
				DIBENZOFURAN	ND	0	10	1	UG/L	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	20	1	UG/L	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	20	1	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8270	NONE	2,4-DICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				2,4-DICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				DIETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				DIETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				2,4-DIMETHYLPHENOL	ND	0	10	1	UG/L	U	U	
				2,4-DIMETHYLPHENOL	ND	0	10	1	UG/L	U	U	
				DIMETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				DIMETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	50	1	UG/L	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	50	1	UG/L	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	10	1	UG/L	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	10	1	UG/L	U	U	
				2,4-DINITROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4-DINITROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				2,4-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				2,6-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				2,6-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				FLUORENE	ND	0	10	1	UG/L	U	U	
				FLUORENE	ND	0	10	1	UG/L	U	U	
				FLUORANTHENE	ND	0	10	1	UG/L	U	U	
				FLUORANTHENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROBENZENE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8270	NONE	HEXACHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROETHANE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROETHANE	ND	0	10	1	UG/L	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	10	1	UG/L	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	10	1	UG/L	U	U	
				ISOPHORONE	ND	0	10	1	UG/L	U	U	
				ISOPHORONE	ND	0	10	1	UG/L	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	10	1	UG/L	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	10	1	UG/L	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	50	1	UG/L	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	50	1	UG/L	U	U	
				2-METHYLNAPHTHALENE	ND	0	10	1	UG/L	U	U	
				2-METHYLNAPHTHALENE	ND	0	10	1	UG/L	U	U	
				NAPHTHALENE	TR	3.9	10	1	UG/L	F	J	
				NAPHTHALENE	TR	3.97	10	1	UG/L	F	J	
				N-NITROSODIPHENYLAMINE	ND	0	10	1	UG/L	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	10	1	UG/L	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	10	1	UG/L	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	10	1	UG/L	U	U	
				2-NITROANILINE	ND	0	50	1	UG/L	U	U	
				2-NITROANILINE	ND	0	50	1	UG/L	U	U	
				3-NITROANILINE	ND	0	50	1	UG/L	U	U	
				3-NITROANILINE	ND	0	50	1	UG/L	U	U	
				4-NITROANILINE	ND	0	50	1	UG/L	U	U	
				4-NITROANILINE	ND	0	50	1	UG/L	U	U	
				NITROBENZENE	ND	0	10	1	UG/L	U	U	
				NITROBENZENE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-DRUM	9/24/2002	SW8270	NONE	2-NITROPHENOL	ND	0	10	1	UG/L	U	U	
				2-NITROPHENOL	ND	0	10	1	UG/L	U	U	
				4-NITROPHENOL	ND	0	50	1	UG/L	U	U	
				4-NITROPHENOL	ND	0	50	1	UG/L	U	U	
				PENTACHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				PENTACHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				PHENANTHRENE	ND	0	10	1	UG/L	U	U	
				PHENANTHRENE	ND	0	10	1	UG/L	U	U	
				PHENOL	TR	6.04	10	1	UG/L	F	J	
				PHENOL	TR	6.14	10	1	UG/L	F	J	
				PYRENE	ND	0	10	1	UG/L	U	U	
				PYRENE	ND	0	10	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
		TNRCC 1		TOTAL PETROLEUM HYDROCARBONS	ND	0	5	1	MG/L	U	U	
CS-B3-ET-WC03	9/30/2002	SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.185	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-WC03	9/30/2002	SW6010	SW1311	ANTIMONY	ND	0	0.045	1	MG/L	U	U	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	UN	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				BENZENE	ND	0	0.1	4	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.1	4	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.1	4	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.1	4	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.1	4	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.2	4	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.39	0.025	1	MG/L	M	E	
				TETRACHLOROETHYLENE(PCE)	=	0.54	0.1	4	MG/L	M	D	
				TRICHLOROETHYLENE (TCE)	TR	0.01	0.025	1	MG/L	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.011	0.1	4	MG/L	F	DJ	
				CHLOROFORM	TR	0.0099	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.012	0.1	4	MG/L	F	DJ	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
				VINYL CHLORIDE	ND	0	0.2	4	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CS-B3-ET-WC04	10/2/2002	SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.142	0.005	1	MG/L			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CS-B3-ET-WC04	10/2/2002	SW6010	SW1311	BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	ND	0	0.045	1	MG/L	U	U	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	M	UN	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	=	0.13	0.025	1	MG/L	R		
				TRICHLOROETHYLENE (TCE)	TR	0.0027	0.025	1	MG/L	R	J	
				CHLOROFORM	TR	0.011	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 ET-1	9/17/2002	SW6010		ARSENIC	=	3.3	2	1	MG/KG		B	
				BARIUM	=	27.1	0.4	1	MG/KG			
				CADMIUM	TR	0.16	0.2	1	MG/KG	F	B	
				CHROMIUM	=	5.7	1	1	MG/KG			
				COPPER	=	5.1	2	1	MG/KG			
				NICKEL	=	5.7	2	1	MG/KG			
				LEAD	=	7.8	2	1	MG/KG			
				ZINC	=	10.1	2	1	MG/KG			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-1	9/17/2002	SW8260	NONE	ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-1	9/17/2002	SW8260	NONE	cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	TR	0.0009	0.005	1	MG/KG	F	J	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.0708	0.005	1	MG/KG			
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-1	9/17/2002	SW8260	NONE	1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	=	0.0062	0.005	1	MG/KG			
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-10		SW6010		ARSENIC	TR	1.3	2	1	MG/KG	F	B	
				BARIUM	=	8.7	0.4	1	MG/KG			
				CADMIUM	TR	0.07	0.2	1	MG/KG	F	B	
				CHROMIUM	=	6.7	1	1	MG/KG			
				COPPER	=	3.1	2	1	MG/KG			
				NICKEL	=	5.6	2	1	MG/KG			
				LEAD	=	2.3	2	1	MG/KG		B	
				ZINC	=	8.8	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-10	9/17/2002	SW8260	NONE	TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-10	9/17/2002	SW8260	NONE	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0023	0.005	1	MG/KG	F	J	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-10	9/17/2002	SW8260	NONE	O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-11		SW6010		ARSENIC	=	3.1	2	1	MG/KG		B	
				BARIUM	=	21.4	0.4	1	MG/KG			
				CADMIUM	=	0.23	0.2	1	MG/KG		B	
				CHROMIUM	=	6.1	1	1	MG/KG			
				COPPER	=	32.2	2	1	MG/KG			
				NICKEL	=	5.1	2	1	MG/KG			
				LEAD	=	20.5	2	1	MG/KG			
				ZINC	=	46	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-11	9/17/2002	SW8260	NONE	DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-11	9/17/2002	SW8260	NONE	n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.001	0.005	1	MG/KG	F	J	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-12		SW6010		ARSENIC	=	2.3	2	1	MG/KG		B	
				BARIUM	=	8.7	0.4	1	MG/KG			
				CADMIUM	TR	0.07	0.2	1	MG/KG	F	B	
				CHROMIUM	=	3.4	1	1	MG/KG			
				COPPER	=	3.3	2	1	MG/KG			
				NICKEL	=	5.1	2	1	MG/KG			
				LEAD	TR	1.9	2	1	MG/KG	F	B	
				ZINC	=	6.4	2	1	MG/KG			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-12	9/17/2002	SW8260	NONE	ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-12	9/17/2002	SW8260	NONE	cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-12	9/17/2002	SW8260	NONE	1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-13		SW6010		ARSENIC	=	2.1	2	1	MG/KG		B	
				BARIUM	=	10.3	0.4	1	MG/KG			
				CADMIUM	=	0.36	0.2	1	MG/KG		B	
				CHROMIUM	=	4.1	1	1	MG/KG			
				COPPER	=	7.3	2	1	MG/KG			
				NICKEL	=	5.4	2	1	MG/KG			
				LEAD	=	7.2	2	1	MG/KG			
				ZINC	=	20.6	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-13	9/17/2002	SW8260	NONE	TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-13	9/17/2002	SW8260	NONE	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.0111	0.005	1	MG/KG			
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0035	0.005	1	MG/KG	F	J	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-13	9/17/2002	SW8260	NONE	O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-14		SW6010		ARSENIC	TR	1.3	2	1	MG/KG	F	B	
				BARIUM	=	9.7	0.4	1	MG/KG			
				CADMIUM	TR	0.19	0.2	1	MG/KG	F	B	
				CHROMIUM	=	2.8	1	1	MG/KG			
				COPPER	=	3.1	2	1	MG/KG			
				NICKEL	=	3.4	2	1	MG/KG			
				LEAD	=	38	2	1	MG/KG			
				ZINC	=	46.7	2	1	MG/KG			
		SW8260		ACETONE	ND	0	50	10	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	2	10	MG/KG	U	U	
				BROMOBENZENE	ND	0	5	10	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	5	10	MG/KG	U	U	
				BROMOMETHANE	ND	0	10	10	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	5	10	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	5	10	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	5	10	MG/KG	U	U	
				BENZENE	ND	0	2	10	MG/KG	U	U	
				TOLUENE	ND	0	5	10	MG/KG	U	U	
				CHLOROBENZENE	ND	0	2	10	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	5	10	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	5	10	MG/KG	U	U	
				CHLOROETHANE	ND	0	5	10	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	5	10	MG/KG	U	U	
				CHLOROMETHANE	ND	0	5	10	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	5	10	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	6	10	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-14	9/17/2002	SW8260	NONE	DIBROMOCHLOROMETHANE	ND	0	3	10	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	10	10	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	5	10	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	5	10	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	3	10	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	5	10	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	6	10	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	2	10	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	6	10	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	5	10	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	5	10	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	5	10	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	3	10	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	5	10	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	5	10	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	2	10	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	5	10	MG/KG	U	U	
				ETHYLBENZENE	ND	0	5	10	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	5	10	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	5	10	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	5	10	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	3	10	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	5	10	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	20	10	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	20	10	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	5	10	MG/KG	U	U	
				NAPHTHALENE	ND	0	5	10	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-14	9/17/2002	SW8260	NONE	n-PROPYLBENZENE	ND	0	5	10	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	3	10	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	=	1443	100	200	MG/KG		D	
				STYRENE	ND	0	5	10	MG/KG	U	U	
				BROMOFORM	ND	0	6	10	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	20	10	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	3	10	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	5	10	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	5	10	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	5	10	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	5	10	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	=	11.51	5	10	MG/KG			
				CHLOROFORM	ND	0	2	10	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	5	10	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	6	10	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	5	10	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	5	10	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	5	10	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	5	10	MG/KG	U	U	
CSB3 ET-15		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	M	U	M
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	TR	0.19	0.7	1	MG/KG	M	J	M

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-15	9/17/2002	SW8270	NONE	4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	M	U	M
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	M	U	M
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	M	U	M
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	M	U	M

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-15	9/17/2002	SW8270	NONE	2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	M	U	M
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	M	U	M
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-15	9/17/2002	SW8270	NONE	2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3 ET-16		SW6010		ARSENIC	TR	1.9	2	1	MG/KG	F	B	
				BARIUM	=	9.5	0.4	1	MG/KG			
				CADMIUM	=	0.33	0.2	1	MG/KG		B	
				CHROMIUM	=	4.4	1	1	MG/KG			
				COPPER	=	8.6	2	1	MG/KG			
				NICKEL	=	4.8	2	1	MG/KG			
				LEAD	=	6.9	2	1	MG/KG			
				ZINC	=	103	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-16	9/17/2002	SW8260	NONE	P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-16	9/17/2002	SW8260	NONE	NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.0079	0.005	1	MG/KG			
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-16	9/17/2002	SW8270	NONE	bis(2-ETHYLHEXYL) PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-16	9/17/2002	SW8270	NONE	DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-16	9/17/2002	SW8270	NONE	1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3 ET-17	9/25/2002	SW8260		TETRACHLOROETHYLENE(PCE)	TR	0.36	0.6	1	MG/KG	F	DJ	
				TETRACHLOROETHYLENE(PCE)	=	0.1424	0.006	1	MG/KG		E	J
				TRICHLOROETHYLENE (TCE)	ND	0	0.006	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.6	1	MG/KG	U	U	
CSB3 ET-18				TETRACHLOROETHYLENE(PCE)	ND	0	0.006	1	MG/KG	U	U	
CSB3 ET-19		SW6010		COPPER	=	3	2	1	MG/KG			
CSB3 ET-2	9/17/2002	SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-2	9/17/2002	SW8270	NONE	4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-2	9/17/2002	SW8270	NONE	2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3 ET-3		SW6010		ARSENIC	=	2.5	2	1	MG/KG		B	
				BARIUM	=	16.5	0.4	1	MG/KG			
				CADMIUM	=	0.23	0.2	1	MG/KG		B	
				CHROMIUM	=	6.4	1	1	MG/KG			
				COPPER	=	5.9	2	1	MG/KG			
				NICKEL	=	7.2	2	1	MG/KG			
				LEAD	=	4.3	2	1	MG/KG			
				ZINC	=	18.6	2	1	MG/KG			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-3	9/17/2002	SW8260	NONE	ACETONE	ND	0	250	50	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	10	50	MG/KG	U	U	
				BROMOBENZENE	ND	0	25	50	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	25	50	MG/KG	U	U	
				BROMOMETHANE	ND	0	50	50	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	25	50	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	25	50	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	25	50	MG/KG	U	U	
				BENZENE	ND	0	10	50	MG/KG	U	U	
				TOLUENE	ND	0	25	50	MG/KG	U	U	
				CHLOROBENZENE	ND	0	10	50	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	25	50	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	25	50	MG/KG	U	U	
				CHLOROETHANE	ND	0	25	50	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	25	50	MG/KG	U	U	
				CHLOROMETHANE	ND	0	25	50	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	25	50	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	30	50	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	15	50	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	50	50	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	25	50	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	25	50	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	15	50	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	25	50	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	30	50	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	10	50	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	30	50	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-3	9/17/2002	SW8260	NONE	cis-1,2-DICHLOROETHYLENE	ND	0	25	50	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	25	50	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	25	50	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	15	50	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	25	50	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	25	50	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	10	50	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	25	50	MG/KG	U	U	
				ETHYLBENZENE	ND	0	25	50	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	25	50	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	25	50	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	25	50	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	15	50	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	25	50	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	100	50	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	100	50	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	25	50	MG/KG	U	U	
				NAPHTHALENE	TR	18.98	25	50	MG/KG	F	J	
				n-PROPYLBENZENE	ND	0	25	50	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	15	50	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	=	5192	500	1000	MG/KG		D	
				STYRENE	ND	0	25	50	MG/KG	U	U	
				BROMOFORM	ND	0	30	50	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	100	50	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	15	50	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	25	50	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	25	50	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-3	9/17/2002	SW8260	NONE	1,2,3-TRICHLOROBENZENE	ND	0	25	50	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	25	50	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	=	130.8	25	50	MG/KG			
				CHLOROFORM	ND	0	10	50	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	25	50	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	TR	4.57	30	50	MG/KG	F	J	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	25	50	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	25	50	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	25	50	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	25	50	MG/KG	U	U	
CSB3 ET-4		SW6010		ARSENIC	TR	1.9	2	1	MG/KG	F	B	
				BARIUM	=	7.8	0.4	1	MG/KG			
				CADMIUM	TR	0.07	0.2	1	MG/KG	F	B	
				CHROMIUM	=	3.7	1	1	MG/KG			
				COPPER	=	5.7	2	1	MG/KG			
				NICKEL	=	4.5	2	1	MG/KG			
				LEAD	=	4.7	2	1	MG/KG			
				ZINC	=	10.4	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-4	9/17/2002	SW8260	NONE	TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-4	9/17/2002	SW8260	NONE	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0009	0.005	1	MG/KG	F	J	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-4	9/17/2002	SW8260	NONE	O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-5		SW6010		ARSENIC	=	2.6	2	1	MG/KG		B	
				BARIUM	=	8.6	0.4	1	MG/KG			
				CADMIUM	TR	0.12	0.2	1	MG/KG	F	B	
				CHROMIUM	=	4.3	1	1	MG/KG			
				COPPER	=	2.9	2	1	MG/KG			
				NICKEL	=	5.8	2	1	MG/KG			
				LEAD	TR	1.5	2	1	MG/KG	F	B	
				ZINC	=	15.4	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-5	9/17/2002	SW8260	NONE	DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-5	9/17/2002	SW8260	NONE	n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-6		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	M	U	M
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	TR	0.37	0.7	1	MG/KG	M	J	M

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-6	9/17/2002	SW8270	NONE	4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	M	U	M
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	M	U	M
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	M	U	M
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	M	U	M
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	TR	0.12	0.7	1	MG/KG	F	J	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	M	U	M

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-6	9/17/2002	SW8270	NONE	2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	M	U	M
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	TR	0.56	0.7	1	MG/KG	F		
				NAPHTHALENE	=	1.26	3.5	5	MG/KG			F
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	TR	0.11	0.7	1	MG/KG	F	J	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	TR	0.11	0.7	1	MG/KG	M	J	M
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-6	9/17/2002	SW8270	NONE	2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3 ET-7		SW6010		ARSENIC	TR	1.8	2	1	MG/KG	F	B	
				BARIUM	=	10.4	0.4	1	MG/KG			
				CADMIUM	TR	0.08	0.2	1	MG/KG	F	B	
				CHROMIUM	=	5.2	1	1	MG/KG			
				COPPER	=	9.6	2	1	MG/KG			
				NICKEL	=	4.9	2	1	MG/KG			
				LEAD	=	3.7	2	1	MG/KG			
				ZINC	=	25.2	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-7	9/17/2002	SW8260	NONE	P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	TR	0.0012	0.005	1	MG/KG	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-7	9/17/2002	SW8260	NONE	NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0013	0.005	1	MG/KG	F	J	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 ET-8		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-8	9/17/2002	SW8270	NONE	bis(2-ETHYLHEXYL) PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	U
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-8	9/17/2002	SW8270	NONE	DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-8	9/17/2002	SW8270	NONE	1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3 ET-9		SW6010		ARSENIC	TR	1.8	2	1	MG/KG	F	B	
				BARIUM	=	7.9	0.4	1	MG/KG			
				CADMIUM	TR	0.06	0.2	1	MG/KG	F	B	
				CHROMIUM	=	5	1	1	MG/KG			
				COPPER	=	2.7	2	1	MG/KG			
				NICKEL	=	4.9	2	1	MG/KG			
				LEAD	TR	1.4	2	1	MG/KG	F	B	
				ZINC	=	5.6	2	1	MG/KG			
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-9	9/17/2002	SW8260	NONE	CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 ET-9	9/17/2002	SW8260	NONE	METHYLENE CHLORIDE	TR	0.0014	0.005	1	MG/KG	F	J	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSB3 WT A-00	9/12/2002	SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	=	0.0538	0.045	1	MG/L			
				BARIUM	=	0.663	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.01	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag

Flag 2 = Laboratory data qualifier

Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-00	9/12/2002	SW6010	SW1311	NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.03	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0017	0.025	1	MG/L	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.011	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.014	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT A-1		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.777	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0082	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.053	0.02	1	MG/L			
				LEAD	=	0.933	0.05	1	MG/L			
				ANTIMONY	=	0.0454	0.045	1	MG/L			
				SELENIUM	TR	0.04	0.1	1	MG/L	F	B	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-1	9/12/2002	SW8260	SW1311	BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0016	0.025	1	MG/L	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.0034	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.012	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	M	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT A-2		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	1.49	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0168	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.056	0.02	1	MG/L			
				LEAD	=	1.47	0.05	1	MG/L			
				ANTIMONY	TR	0.041	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.001	0.025	1	MG/L	F	J	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-2	9/12/2002	SW8260	SW1311	METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0065	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.0053	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT A-3		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.832	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0083	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.063	0.02	1	MG/L			
				LEAD	=	2.43	0.05	1	MG/L			
				ANTIMONY	=	1.1	0.045	1	MG/L			
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.0013	0.025	1	MG/L	F	J	
				BENZENE	ND	0	0.25	10	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.25	10	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.25	10	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.25	10	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.25	10	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-3	9/12/2002	SW8260	SW1311	METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.5	10	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0033	0.025	1	MG/L	F	J	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.25	10	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	>	0.57	0.025	1	MG/L	J	E	
				TRICHLOROETHYLENE (TCE)	=	0.5	0.25	10	MG/L		D	
				CHLOROFORM	TR	0.0094	0.025	1	MG/L	F	J	
				CHLOROFORM	ND	0	0.25	10	MG/L	U	U	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
				VINYL CHLORIDE	ND	0	0.5	10	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT A-4		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	TR	0.0207	0.045	1	MG/L	F	B	
				BARIUM	=	0.625	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0133	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.071	0.02	1	MG/L			
				LEAD	=	6.7	0.05	1	MG/L			
				ANTIMONY	=	0.441	0.045	1	MG/L			
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.0024	0.025	1	MG/L	F	J	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-4	9/12/2002	SW8260	SW1311	METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	=	0.066	0.025	1	MG/L			
				TRICHLOROETHYLENE (TCE)	=	0.044	0.025	1	MG/L			
				CHLOROFORM	TR	0.012	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT A-5		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.195	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.01	0.02	1	MG/L	F	B	
				LEAD	TR	0.03	0.05	1	MG/L	F	B	
				ANTIMONY	TR	0.0291	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.013	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT A-5	9/12/2002	TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT B-1		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.598	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0095	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.019	0.02	1	MG/L	F	B	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0282	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0022	0.025	1	MG/L	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.0073	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.0092	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT B-2		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	TR	0.0231	0.045	1	MG/L	F	B	
				BARIUM	=	0.18	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT B-2	9/12/2002	SW6010	SW1311	CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0173	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	TR	0.0097	0.025	1	MG/L	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.0039	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.011	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT B-3		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.127	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.007	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.041	0.02	1	MG/L			
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0252	0.045	1	MG/L	F	B	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT B-3	9/12/2002	SW6010	SW1311	SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	=	0.061	0.025	1	MG/L			
				CHLOROFORM	TR	0.0096	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT C-1		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.412	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0011	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.015	0.02	1	MG/L	F	B	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0358	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT C-1	9/12/2002	SW8260	SW1311	1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.015	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	=	119	50	1	MG/KG			
CSB3 WT C-2		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.792	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0373	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.036	0.02	1	MG/L			
				LEAD	=	0.657	0.05	1	MG/L			
				ANTIMONY	=	0.0651	0.045	1	MG/L			
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0041	0.025	1	MG/L	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT C-2	9/12/2002	SW8260	SW1311	CHLOROFORM	TR	0.012	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT C-3		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.765	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0084	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.06	0.02	1	MG/L			
				LEAD	TR	0.049	0.05	1	MG/L	F	B	
				ANTIMONY	TR	0.0432	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.0013	0.025	1	MG/L	F	J	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	=	0.036	0.025	1	MG/L			
				CHLOROFORM	TR	0.0092	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT C-4		SW6010	SW1311	SILVER	TR	0.006	0.01	1	MG/L	F	B	
				ARSENIC	TR	0.0209	0.045	1	MG/L	F	B	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT C-4	9/12/2002	SW6010	SW1311	BARIUM	=	0.526	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.574	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.557	0.02	1	MG/L			
				LEAD	=	0.106	0.05	1	MG/L			
				ANTIMONY	TR	0.0372	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0029	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.014	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT C-5		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.337	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.037	0.02	1	MG/L			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT C-5	9/12/2002	SW6010	SW1311	LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0155	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.013	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT D-1		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.604	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0066	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.013	0.02	1	MG/L	F	B	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0215	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT D-1	9/12/2002	SW8260	SW1311	CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.015	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT D-2		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	TR	0.0141	0.045	1	MG/L	F	B	
				BARIUM	=	0.708	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0067	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.091	0.02	1	MG/L			
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0409	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT D-2	9/12/2002	SW8260	SW1311	TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0013	0.025	1	MG/L	F	J	
				CHLOROFORM	TR	0.012	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT D-3		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.397	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.033	0.02	1	MG/L			
				LEAD	TR	0.026	0.05	1	MG/L	F	B	
				ANTIMONY	TR	0.0292	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	TR	0.0002	0.0003	1	MG/L	F	B	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.012	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT D-4	9/12/2002	SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.413	0.005	1	MG/L		E	
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0011	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.019	0.02	1	MG/L	F	B	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0343	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	U	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	U	U	
				CHLOROFORM	TR	0.01	0.025	1	MG/L	F	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	U	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-1 0-2'	9/5/2002	SW6010		ARSENIC	ND	0	2	1	MG/KG	U	U	
				BARIUM	=	46	0.4	1	MG/KG			
				CADMIUM	=	1	0.2	1	MG/KG			
				CHROMIUM	=	11.3	1	1	MG/KG			
				COPPER	=	16	2	1	MG/KG			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW6010	NONE	NICKEL	=	10	2	1	MG/KG			
				LEAD	=	51.4	2	1	MG/KG			
				ZINC	=	57	2	1	MG/KG			
			SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.503	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0041	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0219	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	M	U	
		SW8260	NONE	ACETONE	ND	0	0.24	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.025	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.025	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.05	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.025	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.025	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW8260	NONE	SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.025	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				BENZENE	ND	0	0.01	1	MG/KG	U	U	
				TOLUENE	ND	0	0.025	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.025	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.025	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.025	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.025	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.025	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.029	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.05	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW8260	NONE	DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.025	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.025	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.025	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.029	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.025	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.025	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.025	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.015	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.025	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.025	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW8260	NONE	1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.025	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.025	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.025	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.025	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.025	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.015	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.025	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.1	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.1	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.025	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.025	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.025	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW8260	NONE	TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/KG	U	U	
				STYRENE	ND	0	0.025	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.029	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.1	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.025	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.025	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.025	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.025	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.01	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.025	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.025	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	SW8260	NONE	1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.025	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.025	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.025	1	MG/KG	U	U	
			SW1311	BENZENE	ND	0	0.025	1	MG/L	R	U	
				BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.016	0.025	1	MG/L	R	J	
				CHLOROFORM	TR	0.022	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 0-2'	9/5/2002	TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-1 2-8'		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.452	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0012	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.01	0.02	1	MG/L	F	B	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0176	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.016	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-1 8-16'		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.334	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-1 8-16'	9/5/2002	SW6010	SW1311	CADMIUM	ND	0	0.005	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.033	0.02	1	MG/L			
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.025	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.016	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	=	537	50	1	MG/KG			
CSB3 WT-2 0-2'		SW6010		ARSENIC	ND	0	2	1	MG/KG	U	U	
				BARIUM	=	38.7	0.4	1	MG/KG			
				CADMIUM	=	0.55	0.2	1	MG/KG			
				CHROMIUM	=	10.7	1	1	MG/KG			
				COPPER	=	10.2	2	1	MG/KG			
				NICKEL	=	7.9	2	1	MG/KG			
				LEAD	=	11	2	1	MG/KG			
				ZINC	=	29.5	2	1	MG/KG			
			SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW6010	SW1311	ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.41	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0019	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0181	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260	NONE	ACETONE	ND	0	0.24	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BROMOBENZENE	ND	0	0.024	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.05	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				t-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW8260	NONE	BENZENE	ND	0	0.01	1	MG/KG	U	U	
				TOLUENE	ND	0	0.024	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.024	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.024	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.024	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.029	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	M	U	
				DIBROMOCHLOROMETHANE	ND	0	0.015	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.05	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	M	U	
				DIBROMOMETHANE	ND	0	0.024	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW8260	NONE	1,2-DICHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2-DICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,3-DICHLOROBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	M	U	
				1,4-DICHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.029	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.024	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.024	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.024	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.015	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.024	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.024	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW8260	NONE	ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.024	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	M	U	
				HEXACHLOROBUTADIENE	ND	0	0.015	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.024	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	M	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.1	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.1	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.024	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	M	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.024	1	MG/KG	U	U	
				STYRENE	ND	0	0.024	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW8260	NONE	BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.029	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.1	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.024	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.01	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	M	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	M	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.024	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 0-2'	9/5/2002	SW8260	NONE	M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.024	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.024	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
			SW1311	BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.016	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-2 2-6'		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.551	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.109	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	TR	0.015	0.02	1	MG/L	F	B	
				LEAD	=	0.05	0.05	1	MG/L		B	
				ANTIMONY	TR	0.0295	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 2-6'	9/5/2002	SW8260	SW1311	CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	TR	0.006	0.025	1	MG/L	R	J	
				CHLOROFORM	TR	0.016	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-2 6-12'		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.54	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0879	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.033	0.02	1	MG/L			
				LEAD	TR	0.031	0.05	1	MG/L	F	B	
				ANTIMONY	TR	0.0362	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-2 6-12'	9/5/2002	SW8260	SW1311	TRICHLOROETHYLENE (TCE)	TR	0.0027	0.025	1	MG/L	R	J	
				CHLOROFORM	TR	0.013	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	=	57	50	1	MG/KG			
CSB3 WT-3 2-4'		SW6010		ARSENIC	TR	0.5	2	1	MG/KG	F	B	
				BARIUM	=	41.7	0.4	1	MG/KG			
				CADMIUM	=	0.66	0.2	1	MG/KG			
				CHROMIUM	=	10.5	1	1	MG/KG			
				COPPER	=	8.4	2	1	MG/KG			
				NICKEL	=	8.3	2	1	MG/KG			
				LEAD	=	12.8	2	1	MG/KG			
				ZINC	=	20.9	2	1	MG/KG			
			SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.381	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	TR	0.0034	0.005	1	MG/L	F	B	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.02	1	MG/L	U	U	
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0288	0.045	1	MG/L	F	B	
				SELENIUM	TR	0.05	0.1	1	MG/L	F	B	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260	NONE	ACETONE	ND	0	0.24	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 2-4'	9/5/2002	SW8260	NONE	BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.024	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.05	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				BENZENE	ND	0	0.01	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				TOLUENE	ND	0	0.024	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.024	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.024	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.024	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 2-4'	9/5/2002	SW8260	NONE	CHLOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.029	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.05	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.024	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.01	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.029	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.024	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.024	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 2-4'	9/5/2002	SW8260	NONE	1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.024	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.015	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.024	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.024	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.024	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.015	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.024	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.1	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.1	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 2-4'	9/5/2002	SW8260	NONE	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.024	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.024	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.024	1	MG/KG	U	U	
				STYRENE	ND	0	0.024	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.029	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.1	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.015	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.024	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.024	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 2-4'	9/5/2002	SW8260	NONE	TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.024	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.01	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.024	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.029	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.024	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.024	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.024	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.024	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
			SW1311	BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.012	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 4-6'	9/5/2002	SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.887	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.3	0.005	1	MG/L			
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.032	0.02	1	MG/L			
				LEAD	=	0.131	0.05	1	MG/L			
				ANTIMONY	TR	0.024	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	TR	0.00078	0.025	1	MG/L	R	J	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0033	0.025	1	MG/L	R	J	
				CHLOROFORM	TR	0.013	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSB3 WT-3 6-TD		SW6010	SW1311	SILVER	ND	0	0.01	1	MG/L	U	U	
				ARSENIC	ND	0	0.045	1	MG/L	U	U	
				BARIUM	=	0.508	0.005	1	MG/L			
				BERYLLIUM	ND	0	0.005	1	MG/L	U	U	
				CADMIUM	=	0.0992	0.005	1	MG/L			

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3 WT-3 6-TD	9/5/2002	SW6010	SW1311	CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				NICKEL	=	0.03	0.02	1	MG/L			
				LEAD	ND	0	0.05	1	MG/L	U	U	
				ANTIMONY	TR	0.0385	0.045	1	MG/L	F	B	
				SELENIUM	ND	0	0.1	1	MG/L	U	U	
		SW7470		MERCURY	ND	0	0.0003	1	MG/L	U	U	
		SW8260		BENZENE	ND	0	0.025	1	MG/L	R	U	
				CHLOROBENZENE	ND	0	0.025	1	MG/L	R	U	
				CARBON TETRACHLORIDE	ND	0	0.025	1	MG/L	R	U	
				1,2-DICHLOROETHANE	ND	0	0.025	1	MG/L	R	U	
				1,1-DICHLOROETHENE	ND	0	0.025	1	MG/L	R	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.05	1	MG/L	R	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.025	1	MG/L	R	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.025	1	MG/L	R	U	
				CHLOROFORM	TR	0.013	0.025	1	MG/L	R	J	
				VINYL CHLORIDE	ND	0	0.05	1	MG/L	R	U	
		TNRCC 1	NONE	TOTAL PETROLEUM HYDROCARBONS	ND	0	50	1	MG/KG	U	U	
CSCS-1	9/9/2002	SW6010		ARSENIC	=	3.4	2	1	MG/KG			
				BARIUM	=	36.9	0.4	1	MG/KG		E*	
				CADMIUM	=	0.63	0.2	1	MG/KG		N*	
				CHROMIUM	=	11	1	1	MG/KG			
				COPPER	=	11.9	2	1	MG/KG		E*	
				NICKEL	=	7	2	1	MG/KG		NE	
				LEAD	=	28.4	2	1	MG/KG		NE	
				ZINC	=	110	2	1	MG/KG		NE	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-1	9/9/2002	SW8260	NONE	BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-1	9/9/2002	SW8260	NONE	1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	TR	0.0011	0.005	1	MG/KG	F	J	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-1	9/9/2002	SW8260	NONE	TRICHLOROETHYLENE (TCE)	TR	0.0009	0.005	1	MG/KG	F	J	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSCS-10	10/11/2002	SW6010		ZINC	=	42.1	2	1	MG/KG			
		SW7471		MERCURY	ND	0	0.2	2	MG/KG	U	U	
CSCS-2	9/9/2002	SW6010		ARSENIC	=	3.5	2	1	MG/KG			
				BARIUM	=	57.6	0.4	1	MG/KG		E*	
				CADMIUM	=	0.39	0.2	1	MG/KG		N*	
				CHROMIUM	=	11.8	1	1	MG/KG			
				COPPER	=	8.2	2	1	MG/KG		E*	
				NICKEL	=	9	2	1	MG/KG		NE	
				LEAD	=	11.6	2	1	MG/KG		NE	
				ZINC	=	28.1	2	1	MG/KG		NE	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-2	9/9/2002	SW8260	NONE	BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-2	9/9/2002	SW8260	NONE	DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-2	9/9/2002	SW8260	NONE	trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	TR	0.001	0.005	1	MG/KG	F	J	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-2	9/9/2002	SW8260	NONE	n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-2	9/9/2002	SW8260	NONE	1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSCS-3		SW6010		ARSENIC	=	5.3	2	1	MG/KG			
				BARIUM	=	40.8	0.4	1	MG/KG		E*	
				CADMIUM	=	1.4	0.2	1	MG/KG		N*	
				CHROMIUM	=	9.6	1	1	MG/KG			
				COPPER	=	12.6	2	1	MG/KG		E*	
				NICKEL	=	7.5	2	1	MG/KG		NE	
				LEAD	=	29.1	2	1	MG/KG		NE	
				ZINC	=	88.3	2	1	MG/KG		NE	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-3	9/9/2002	SW8260	NONE	BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	M	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-3	9/9/2002	SW8260	NONE	P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	M	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	M	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	TR	0.002	0.005	1	MG/KG	F	J	
				cis-1,2-DICHLOROETHYLENE	TR	0.0009	0.005	1	MG/KG	F	J	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-3	9/9/2002	SW8260	NONE	trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	M	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	M	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	TR	0.0008	0.005	1	MG/KG	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-3	9/9/2002	SW8260	NONE	NAPHTHALENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	M	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	=	0.0072	0.005	1	MG/KG	J		
				TRICHLOROETHYLENE (TCE)	=	0.0096	0.005	1	MG/KG	J		
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-3	9/9/2002	SW8260	NONE	1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	M	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	M	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSCS-4		SW6010		ARSENIC	=	3.4	2	1	MG/KG			
				BARIUM	=	40.3	0.4	1	MG/KG	M	ME*	
				CADMIUM	=	0.4	0.2	1	MG/KG	M	N*	
				CHROMIUM	=	10.7	1	1	MG/KG			
				COPPER	=	9.5	2	1	MG/KG		E*	
				NICKEL	=	8	2	1	MG/KG	M	MNE	
				LEAD	=	19.5	2	1	MG/KG	M	MNE	
				ZINC	=	31.4	2	1	MG/KG	M	MNE	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	M	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-4	9/9/2002	SW8260	NONE	BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	M	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-4	9/9/2002	SW8260	NONE	P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	M	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	M	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	M	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	M	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	M	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	M	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	M	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	M	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	M	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-4	9/9/2002	SW8260	NONE	cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	M	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	M	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	M	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	M	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	M	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-4	9/9/2002	SW8260	NONE	NAPHTHALENE	ND	0	0.005	1	MG/KG	M	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	M	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	M	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				TRICHLOROETHYLENE (TCE)	TR	0.001	0.005	1	MG/KG	F	J	
				TRICHLOROETHYLENE (TCE)	ND	0	0.005	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-4	9/9/2002	SW8260	NONE	CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	M	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	M	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	M	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSCS-5		SW6010		ARSENIC	=	3.3	2	1	MG/KG			
				BARIUM	=	29.2	0.4	1	MG/KG		E*	
				CADMIUM	=	0.45	0.2	1	MG/KG		N*	
				CHROMIUM	=	8.3	1	1	MG/KG			
				COPPER	=	10.4	2	1	MG/KG		E*	
				NICKEL	=	6.5	2	1	MG/KG		NE	
				LEAD	=	12.6	2	1	MG/KG		NE	
				ZINC	=	186	2	1	MG/KG		NE	
		SW8260		ACETONE	ND	0	0.05	1	MG/KG	U	U	
				ACETONE	ND	0	0.05	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMODICHLOROMETHANE	ND	0	0.002	1	MG/KG	U	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BROMOBENZENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-5	9/9/2002	SW8260	NONE	BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOCHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				BROMOMETHANE	ND	0	0.01	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				SEC-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				t-BUTYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				BENZENE	ND	0	0.002	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				TOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				CHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				2-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	M	U	
				4-CHLOROTOLUENE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				1-CHLOROHEXANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CHLOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-5	9/9/2002	SW8260	NONE	CARBON TETRACHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	M	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	0.006	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	0.01	1	MG/KG	M	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DIBROMOMETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,3-DICHLOROBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	M	U	
				1,4-DICHLOROBENZENE	ND	0	0.002	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				1,1-DICHLOROETHENE	ND	0	0.006	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,1-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-5	9/9/2002	SW8260	NONE	cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.003	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.002	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				2,2-DICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				ETHYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	0.005	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.003	1	MG/KG	M	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	M	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	0.005	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	0.02	1	MG/KG	U	U	
				METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-5	9/9/2002	SW8260	NONE	METHYLENE CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	M	U	
				NAPHTHALENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	U	U	
				n-PROPYLBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	M	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				STYRENE	ND	0	0.005	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				BROMOFORM	ND	0	0.006	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				tert-BUTYL METHYL ETHER	ND	0	0.02	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.003	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.005	1	MG/KG	U	U	
				TRICHLOROETHYLENE (TCE)	TR	0.0022	0.005	1	MG/KG	F	J	
				TRICHLOROETHYLENE (TCE)	TR	0.0019	0.005	1	MG/KG	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-5	9/9/2002	SW8260	NONE	CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				CHLOROFORM	ND	0	0.002	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	0.005	1	MG/KG	M	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	M	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	0.006	1	MG/KG	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	M	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				VINYL CHLORIDE	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	0.005	1	MG/KG	U	U	
CSCS-6	10/11/2002	SW6010		ZINC	=	86.6	2	1	MG/KG			
		SW7471		MERCURY	TR	0.1	0.2	2	MG/KG	F	B	
CSCS-7		SW6010		ZINC	=	35.3	2	1	MG/KG			
		SW7471		MERCURY	ND	0	0.2	2	MG/KG	U	U	
CSCS-8		SW6010		ZINC	=	57.1	2	1	MG/KG			
		SW7471		MERCURY	ND	0	0.2	2	MG/KG	U	U	
CSCS-9		SW6010		ZINC	=	22.2	2	1	MG/KG			
		SW7471		MERCURY	ND	0	0.2	2	MG/KG	U	U	
CSCS-COMPOSITE	9/9/2002			MERCURY	=	0.93	0.1	1	MG/KG	M	N*	
		SW8270		ACENAPHTHENE	ND	0	0.7	1	MG/KG	U	U	
				ACENAPHTHYLENE	ND	0	0.7	1	MG/KG	U	U	
				ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-COMPOSITE	9/9/2002	SW8270	NONE	bis(2-CHLOROETHOXY) METHANE	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	0.7	1	MG/KG	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	0.7	1	MG/KG	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	TR	0.07	0.7	1	MG/KG	F	J	
				4-BROMOPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				BENZO(a)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				BENZOIC ACID	ND	0	5	1	MG/KG	U	U	
				BENZO(a)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(b)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	0.7	1	MG/KG	U	U	
				BENZO(k)FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				BENZYL ALCOHOL	ND	0	1.3	1	MG/KG	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	1.3	1	MG/KG	U	U	
				CHRYSENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROANILINE	ND	0	1.3	1	MG/KG	R	U	
				2-CHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				2-CHLORONAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	0.7	1	MG/KG	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	0.7	1	MG/KG	U	U	
				DIBENZOFURAN	ND	0	0.7	1	MG/KG	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	1.3	1	MG/KG	U	U	
				1,2-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,3-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				DIETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DIMETHYLPHENOL	ND	0	0.3	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-COMPOSITE	9/9/2002	SW8270	NONE	DIMETHYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	3.3	1	MG/KG	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	0.7	1	MG/KG	U	U	
				2,4-DINITROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				2,6-DINITROTOLUENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORENE	ND	0	0.7	1	MG/KG	U	U	
				FLUORANTHENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				HEXACHLOROETHANE	ND	0	0.7	1	MG/KG	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	0.7	1	MG/KG	U	U	
				ISOPHORONE	ND	0	0.7	1	MG/KG	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	0.3	1	MG/KG	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	2	1	MG/KG	U	U	
				2-METHYLNAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				NAPHTHALENE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				3-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				4-NITROANILINE	ND	0	3.3	1	MG/KG	U	U	
				NITROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2-NITROPHENOL	ND	0	0.3	1	MG/KG	U	U	
				4-NITROPHENOL	ND	0	1.6	1	MG/KG	U	U	
				PENTACHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSCS-COMPOSITE	9/9/2002	SW8270	NONE	PHENANTHRENE	ND	0	0.7	1	MG/KG	U	U	
				PHENOL	ND	0	0.3	1	MG/KG	U	U	
				PYRENE	ND	0	0.7	1	MG/KG	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	0.7	1	MG/KG	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	3.3	1	MG/KG	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	0.3	1	MG/KG	U	U	
CSB3ET-EB-1	9/17/2002	SW6010		ARSENIC	ND	0	0.01	1	MG/L	U	U	
				BARIUM	ND	0	0.002	1	MG/L	U	U	
				CADMIUM	ND	0	0.001	1	MG/L	U	U	
				CHROMIUM	ND	0	0.01	1	MG/L	U	U	
				COPPER	ND	0	0.01	1	MG/L	U	U	
				NICKEL	ND	0	0.01	1	MG/L	U	U	
				LEAD	ND	0	0.01	1	MG/L	U	U	
				ZINC	ND	0	0.02	1	MG/L	U	U	
		SW8260		ACETONE	TR	7.3	10	1	UG/L	R	JR	F
				BROMODICHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				BROMOBENZENE	ND	0	1	1	UG/L	U	U	
				BROMOCHLOROMETHANE	ND	0	1	1	UG/L	U	U	
				BROMOMETHANE	ND	0	3	1	UG/L	U	U	
				n-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				SEC-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				t-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				BENZENE	ND	0	0.4	1	UG/L	U	U	
				TOLUENE	ND	0	1	1	UG/L	U	U	
				CHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				2-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	
				4-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3ET-EB-1	9/17/2002	SW8260	NONE	CHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1-CHLOROHEXANE	ND	0	1	1	UG/L	U	U	
				CHLOROMETHANE	TR	0.25	1	1	UG/L	F	J	
				CARBON TETRACHLORIDE	ND	0	1	1	UG/L	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	1	1	UG/L	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	2	1	UG/L	U	U	
				DIBROMOMETHANE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.5	1	UG/L	R	UR	R
				1,2-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	1	1	UG/L	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.5	1	UG/L	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.4	1	UG/L	U	U	
				2,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				ETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	1	1	UG/L	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.6	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3ET-EB-1	9/17/2002	SW8260	NONE	ISOPROPYLBENZENE (CUMENE)	ND	0	1	1	UG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	10	1	UG/L	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	10	1	UG/L	U	U	
				METHYLENE CHLORIDE	ND	0	1	1	UG/L	U	U	
				NAPHTHALENE	ND	0	1	1	UG/L	U	U	
				n-PROPYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	1	1	UG/L	U	U	
				STYRENE	ND	0	1	1	UG/L	U	U	
				BROMOFORM	ND	0	1	1	UG/L	U	U	
				tert-BUTYL METHYL ETHER	ND	0	5	1	UG/L	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	1	1	UG/L	U	U	
				CHLOROFORM	ND	0	0.3	1	UG/L	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	1	1	UG/L	U	U	
				VINYL CHLORIDE	ND	0	1	1	UG/L	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	2	1	UG/L	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	1	1	UG/L	U	U	
		SW8270		ACENAPHTHENE	ND	0	10	1	UG/L	U	U	
				ACENAPHTHYLENE	ND	0	10	1	UG/L	U	U	
				ANTHRACENE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3ET-EB-1	9/17/2002	SW8270	NONE	BENZYL BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHOXY) METHANE	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ND	0	10	1	UG/L	U	U	
				bis(2-CHLOROISOPROPYL) ETHER	ND	0	10	1	UG/L	U	U	
				bis(2-ETHYLHEXYL) PHTHALATE	ND	0	10	1	UG/L	U	U	
				4-BROMOPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				BENZO(a)ANTHRACENE	ND	0	10	1	UG/L	U	U	
				BENZOIC ACID	ND	0	100	1	UG/L	U	U	
				BENZO(a)PYRENE	ND	0	10	1	UG/L	U	U	
				BENZO(b)FLUORANTHENE	ND	0	10	1	UG/L	U	U	
				BENZO(g,h,i)PERYLENE	ND	0	10	1	UG/L	U	U	
				BENZO(k)FLUORANTHENE	ND	0	10	1	UG/L	U	U	
				BENZYL ALCOHOL	ND	0	20	1	UG/L	U	U	
				4-CHLORO-3-METHYLPHENOL	ND	0	20	1	UG/L	U	U	
				CHRYSENE	ND	0	10	1	UG/L	U	U	
				4-CHLOROANILINE	ND	0	20	1	UG/L	U	U	
				2-CHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				2-CHLORONAPHTHALENE	ND	0	10	1	UG/L	U	U	
				4-CHLOROPHENYL PHENYL ETHER	ND	0	10	1	UG/L	U	U	
				DIBENZ(a,h)ANTHRACENE	ND	0	10	1	UG/L	U	U	
				DIBENZOFURAN	ND	0	10	1	UG/L	U	U	
				3,3'-DICHLOROBENZIDINE	ND	0	20	1	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				2,4-DICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
				DIETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3ET-EB-1	9/17/2002	SW8270	NONE	2,4-DIMETHYLPHENOL	ND	0	10	1	UG/L	U	U	
				DIMETHYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				4,6-DINITRO-2-METHYLPHENOL	ND	0	50	1	UG/L	U	U	
				DI-n-BUTYL PHTHALATE	ND	0	10	1	UG/L	U	U	
				DI-n-OCTYLPHTHALATE	ND	0	10	1	UG/L	U	U	
				2,4-DINITROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				2,6-DINITROTOLUENE	ND	0	10	1	UG/L	U	U	
				FLUORENE	ND	0	10	1	UG/L	U	U	
				FLUORANTHENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				HEXACHLOROETHANE	ND	0	10	1	UG/L	U	U	
				INDENO(1,2,3-c,d)PYRENE	ND	0	10	1	UG/L	U	U	
				ISOPHORONE	ND	0	10	1	UG/L	U	U	
				2-METHYLPHENOL (o-CRESOL)	ND	0	10	1	UG/L	U	U	
				4-METHYLPHENOL (p-CRESOL)	ND	0	50	1	UG/L	U	U	
				2-METHYLNAPHTHALENE	ND	0	10	1	UG/L	U	U	
				NAPHTHALENE	ND	0	10	1	UG/L	U	U	
				N-NITROSODIPHENYLAMINE	ND	0	10	1	UG/L	U	U	
				N-NITROSODI-n-PROPYLAMINE	ND	0	10	1	UG/L	U	U	
				2-NITROANILINE	ND	0	50	1	UG/L	U	U	
				3-NITROANILINE	ND	0	50	1	UG/L	U	U	
				4-NITROANILINE	ND	0	50	1	UG/L	U	U	
				NITROBENZENE	ND	0	10	1	UG/L	U	U	
				2-NITROPHENOL	ND	0	10	1	UG/L	U	U	
				4-NITROPHENOL	ND	0	50	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
CSB3ET-EB-1	9/17/2002	SW8270	NONE	PENTACHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				PHENANTHRENE	ND	0	10	1	UG/L	U	U	
				PHENOL	ND	0	10	1	UG/L	U	U	
				PYRENE	ND	0	10	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	10	1	UG/L	U	U	
				2,4,5-TRICHLOROPHENOL	ND	0	50	1	UG/L	U	U	
				2,4,6-TRICHLOROPHENOL	ND	0	10	1	UG/L	U	U	
TB-1_08282002	8/28/2002	SW8260		ACETONE	TR	2.9	10	1	UG/L	F	J	
				BROMODICHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				BROMOBENZENE	ND	0	1	1	UG/L	U	U	
				BROMOCHLOROMETHANE	ND	0	1	1	UG/L	U	U	
				BROMOMETHANE	ND	0	3	1	UG/L	R	UR	
				n-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				SEC-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				t-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				BENZENE	ND	0	0.4	1	UG/L	U	U	
				TOLUENE	TR	0.4	1	1	UG/L	F	J	
				CHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				2-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	
				4-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	
				CHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1-CHLOROHEXANE	ND	0	1	1	UG/L	U	U	
				CHLOROMETHANE	ND	0	1	1	UG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	1	1	UG/L	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	1	1	UG/L	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	2	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
TB-1_08282002	8/28/2002	SW8260	NONE	DIBROMOMETHANE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				1,2-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	1	1	UG/L	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.5	1	UG/L	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.4	1	UG/L	U	U	
				2,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				ETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	1	1	UG/L	U	U	
				TRICHLOROFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.6	1	UG/L	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	1	1	UG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	10	1	UG/L	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	10	1	UG/L	U	U	
				METHYLENE CHLORIDE	ND	0	1	1	UG/L	U	U	
				NAPHTHALENE	ND	0	1	1	UG/L	U	U	
				n-PROPYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
TB-1_08282002	8/28/2002	SW8260	NONE	TETRACHLOROETHYLENE(PCE)	ND	0	1	1	UG/L	U	U	
				STYRENE	ND	0	1	1	UG/L	U	U	
				BROMOFORM	ND	0	1	1	UG/L	U	U	
				tert-BUTYL METHYL ETHER	ND	0	5	1	UG/L	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	1	1	UG/L	U	U	
				CHLOROFORM	ND	0	0.3	1	UG/L	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	1	1	UG/L	U	U	
				VINYL CHLORIDE	ND	0	1	1	UG/L	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	2	1	UG/L	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	1	1	UG/L	U	U	
TRIP BLANK	9/17/2002			ACETONE	TR	5.4	10	1	UG/L	R	R	F
				BROMODICHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				BROMOBENZENE	ND	0	1	1	UG/L	U	U	
				BROMOCHLOROMETHANE	ND	0	1	1	UG/L	U	U	
				BROMOMETHANE	ND	0	3	1	UG/L	U	U	
				n-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				SEC-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				t-BUTYLBENZENE	ND	0	1	1	UG/L	U	U	
				BENZENE	ND	0	0.4	1	UG/L	U	U	
				TOLUENE	TR	0.29	1	1	UG/L	F	J	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
TRIP BLANK	9/17/2002	SW8260	NONE	CHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				2-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	
				4-CHLOROTOLUENE	ND	0	1	1	UG/L	U	U	
				CHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1-CHLOROHEXANE	ND	0	1	1	UG/L	U	U	
				CHLOROMETHANE	ND	0	1	1	UG/L	U	U	
				CARBON TETRACHLORIDE	ND	0	1	1	UG/L	U	U	
				P-CYMENE (p-ISOPROPYLTOLUENE)	ND	0	1	1	UG/L	U	U	
				DIBROMOCHLOROMETHANE	ND	0	0.5	1	UG/L	U	U	
				1,2-DIBROMO-3-CHLOROPROPANE	ND	0	2	1	UG/L	U	U	
				DIBROMOMETHANE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROETHANE	=	0.5	0.5	1	UG/L	R	R	R
				1,2-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,4-DICHLOROBENZENE	ND	0	0.5	1	UG/L	U	U	
				1,1-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				cis-1,2-DICHLOROETHYLENE	ND	0	1	1	UG/L	U	U	
				trans-1,2-DICHLOROETHENE	ND	0	1	1	UG/L	U	U	
				1,1-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				cis-1,3-DICHLOROPROPENE	ND	0	0.5	1	UG/L	U	U	
				trans-1,3-DICHLOROPROPENE	ND	0	1	1	UG/L	U	U	
				1,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,3-DICHLOROPROPANE	ND	0	0.4	1	UG/L	U	U	
				2,2-DICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				ETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ND	0	1	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier

Camp Stanley Storage Activity
Interim Remedial Action
Analytical Results
August - October 2002

Location	Sample Date	Method	Extraction	Parameter	Boolean	Value	RL	Dilution	Units	Flag 1	Flag 2	Flag 3
TRIP BLANK	9/17/2002	SW8260	NONE	TRICHLOROFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				DICHLORODIFLUOROMETHANE	ND	0	1	1	UG/L	U	U	
				HEXACHLOROBUTADIENE	ND	0	0.6	1	UG/L	U	U	
				ISOPROPYLBENZENE (CUMENE)	ND	0	1	1	UG/L	U	U	
				METHYL ETHYL KETONE (2-BUTANONE)	ND	0	10	1	UG/L	U	U	
				METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ND	0	10	1	UG/L	U	U	
				METHYLENE CHLORIDE	ND	0	1	1	UG/L	U	U	
				NAPHTHALENE	ND	0	1	1	UG/L	U	U	
				n-PROPYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,1,2,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				TETRACHLOROETHYLENE(PCE)	ND	0	1	1	UG/L	U	U	
				STYRENE	ND	0	1	1	UG/L	U	U	
				BROMOFORM	ND	0	1	1	UG/L	U	U	
				tert-BUTYL METHYL ETHER	ND	0	5	1	UG/L	U	U	
				1,1,1,2-TETRACHLOROETHANE	ND	0	0.5	1	UG/L	U	U	
				1,1,1-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,1,2-TRICHLOROETHANE	ND	0	1	1	UG/L	U	U	
				1,2,3-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRICHLOROBENZENE	ND	0	1	1	UG/L	U	U	
				TRICHLOROETHYLENE (TCE)	ND	0	1	1	UG/L	U	U	
				CHLOROFORM	ND	0	0.3	1	UG/L	U	U	
				1,2,3-TRICHLOROPROPANE	ND	0	1	1	UG/L	U	U	
				1,2,4-TRIMETHYLBENZENE	ND	0	1	1	UG/L	U	U	
				1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ND	0	1	1	UG/L	U	U	
				VINYL CHLORIDE	ND	0	1	1	UG/L	U	U	
				M,P-XYLENE (SUM OF ISOMERS)	ND	0	2	1	UG/L	U	U	
				O-XYLENE (1,2-DIMETHYLBENZENE)	ND	0	1	1	UG/L	U	U	

Flag 1 = AFCEE QAPP data flag
Flag 2 = Laboratory data qualifier
Flag 3 = EPA data qualifier