Copco1 Reservoir - Sediment - Additional Analytes

Analyte	Sampling Sites																	
	CDH-S-009A(0.0-4.6)	CDH-S-010(0.0-5.0)	CDH-S-010(5.0-8.0)	CDH-S-011(0.0-1.3)	CDH-S-012(0.0-5.4)	CDH-S-013(0.0-5.7)	CDH-S-014(0.0-5.3)	CDH-S-015A(0.0-5.0)	CDH-S-015A(5.0-9.7)	CDH-S-015A(0.0-9.7)	CDH-S-016(0.0-5.0)	CDH-S-016(5.0-7.5)	CDH-S-017(0.0-1.2)	CDH-S-018(0.0-5.0)	CDH-S-018(5.0-8.9)	CDH-S-019(0.0-4.8)	CDH-S-020(0.0-5.0)	CDH-S-020(5.0-7.0)
Organics																		
Pesticides/Herbicides: Organophosphate Pesticides (ug/kg dry weight, Method 8141A unless otherwise noted)																		
0,0,0-Triethylphosphorothioate	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Azinphosmethyl	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Bolstar	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Chlorpyrifos (see footnote for methods)	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Coumaphos	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Demeton	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Demeton-O	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Demeton-S	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Diazinon (Method MLA-047-Rev 03)	-	-	-	-	-	-	<0.018	-	-	<0.016	-	-	-	-	-	-	-	-
Dichlorvos	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Dimethoate	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Disulfoton	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
EPN	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Ethoprop	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Famphur	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Fensulfothion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Fenthion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Malathion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Methyl parathion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Mevinphos	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Parathion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Phorate	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Ronnel	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Stirophos	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Sulfotep	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Thionazin	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Tokuthion	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Trichloronate or Tirchloronat	-	-	-	-	-	-	<350	-	-	<330	-	-	-	-	-	-	-	-
Organics																		
Pesticides/Herbicides: Carbamate Pesticides (ug/kg dry weight, Method MLA-047 Rev 03 unless otherwise noted)																		
3-Hydroxycarbofuran	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Aldicarb	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Aldicarb Sulfone	-	-	-	-	-	-	<0.37	-	-	<0.33	-	-	-	-	-	-	-	-
Aldicarb Sulfoxide	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Aminocarb	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Bendiocarb	-	-	-	-	-	-	<0.37	-	-	<0.32	-	-	-	-	-	-	-	-
Carbaryl	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Carbazole (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Carbofuran	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Dioxacarb	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Imidacloprid	-	-	-	-	-	-	<0.37	-	-	<0.32	-	-	-	-	-	-	-	-
Methiocarb	-	-	-	-	-	-	<0.37	-	-	<0.33	-	-	-	-	-	-	-	-
Methomyl	-	-	-	-	-	-	<0.18 L	-	-	<0.16 L	-	-	-	-	-	-	-	-
Mexacarbate	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Oxamyl	-	-	-	-	-	-	<0.19 L	-	-	<0.17 L	-	-	-	-	-	-	-	-
Piperonyl butoxide	-	-	-	-	-	-	0.040	-	-	0.035	-	-	-	-	-	-	-	-
Pirimicarb	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Promecarb	-	-	-	-	-	-	<0.37	-	-	<0.32	-	-	-	-	-	-	-	-
Propoxur	-	-	-	-	-	-	<0.18	-	-	<0.16	-	-	-	-	-	-	-	-
Organics																		
Pesticides/Herbicides: Pyrethroids (Insecticides) (ug/kg dry weight, Method GCMS-NCI-SIM)																		
Allethrin	-	-	-	-	-	-	<0.33	-	-	<0.33 L	-	-	-	-	-	-	-	-
Bifenthrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Cyfluthrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Cypermethrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Deltamethrin: tralomethrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Esfenvalerate: fenvalerate	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Fenpropathrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Lambda-cyhalothrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Permethrin (Total)	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Tau-Fluvalinate	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-
Tetramethrin	-	-	-	-	-	-	<0.33	-	-	<0.33	-	-	-	-	-	-	-	-

Preliminary Data - Subject to Revision

Copco1 Reservoir - Sediment - Additional Analytes

Analyte	Sampling Sites																	
	CDH-S-009A(0.0-4.6)	CDH-S-010(0.0-5.0)	CDH-S-010(5.0-8.0)	CDH-S-011(0.0-1.3)	CDH-S-012(0.0-5.4)	CDH-S-013(0.0-5.7)	CDH-S-014(0.0-5.3)	CDH-S-015A(0.0-5.0)	CDH-S-015A(5.0-9.7)	CDH-S-015A(0.0-9.7)	CDH-S-016(0.0-5.0)	CDH-S-016(5.0-7.5)	CDH-S-017(0.0-1.2)	CDH-S-018(0.0-5.0)	CDH-S-018(5.0-8.9)	CDH-S-019(0.0-4.8)	CDH-S-020(0.0-5.0)	CDH-S-020(5.0-7.0)
Organics																		
Phthalates (ug/kg dry weight, Method 8270D)																		
Di-N-butyl phthalate	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Diethyl phthalate	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Dimethyl phthalate	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Bis(2-ethylhexyl) phthalate	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Butyl benzyl phthalate	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Di-N-octyl phthalate	<710 V	<580 V	<730 V	<580	<620 V	<700 V	<660	<680 V	<620 V	-	<650 V	<620 V	<700	<710 V	<650 V	<620	<620	<590
Organics																		
VOCs (ug/kg dry weight, Method 8260C, unless otherwise noted)																		
1,2,4-Trimethylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2-Dibromoethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,3,5-Trimethylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
2,4-Dinitrotoluene (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,6-Dinitrotoluene (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2-Butanone	29	<18	<22	<18	<19	36	<20	26	37	-	35	<19	<21	53	29	<19	21	27
2-Hexanone	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
4-Methyl-2-pentanone	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Acetone	360	63	92	40	81	480	90	390	510	-	460	270	100	730	440	120	210	280
Allyl chloride	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Benzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bromobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bromochloromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bromodichloromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bromoform	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bromomethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Carbon disulfide	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Carbon tetrachloride	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Chlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Chloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Chloroform	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Chloromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
cis-1,2-Dichloroethene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
cis-1,3-Dichloropropene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Cyclohexane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Dibromochloromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Dibromomethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Dichlorodifluoromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Dichlorofluoromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Diesel Range Organics (mg/kg, Method 8015 DRO)	<17	<14	<18	65	<15	<17	42	<16	<15	-	<16	<15	47	<17	<16	26	49	40
Ethyl acetate	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Ethyl ether	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Ethyl methacrylate	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Ethylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Freon 113	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Iodomethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Isopropylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
meta, para-Xylene	<43	<36	<45	<36	<38	<42	<40	<41	<38	-	<39	<38	<42	<43	<40	<37	<38	<36
Methyl acetate	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Methylcyclohexane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Methylene chloride	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
MTBE	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
N-Butylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
N-Propylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
ortho-Xylene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Pentachloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
p-Isopropyltoluene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Residual Range Organics (mg/kg, Method 8015 RRO)	<170	<140	<180	320	<150	<170	180	<160	<150	-	<160	<150	220 T	<170	<160	<150 T	240 T	210 T
sec-Butylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Styrene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
tert-Butylbenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Tetrachloroethene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Toluene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18

Preliminary Data - Subject to Revision

Analyte	Copco1 Reservoir - Sediment - Additional Analytes																	
	Sampling Sites																	
	CDH-S-009A(0.0-4.6)	CDH-S-010(0.0-5.0)	CDH-S-010(5.0-8.0)	CDH-S-011(0.0-1.3)	CDH-S-012(0.0-5.4)	CDH-S-013(0.0-5.7)	CDH-S-014(0.0-5.3)	CDH-S-015A(0.0-5.0)	CDH-S-015A(5.0-9.7)	CDH-S-015A(0.0-9.7)	CDH-S-016(0.0-5.0)	CDH-S-016(5.0-7.5)	CDH-S-017(0.0-1.2)	CDH-S-018(0.0-5.0)	CDH-S-018(5.0-8.9)	CDH-S-019(0.0-4.8)	CDH-S-020(0.0-5.0)	CDH-S-020(5.0-7.0)
Organics																		
SVOCs: Phenols (ug/kg dry weight, Method 8270D unless otherwise noted)																		
2,4,5-Trichlorophenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,4,6-Trichlorophenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,4-Dichlorophenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,4-Dimethylphenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2-Chlorophenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2-Methylphenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Chloro-3-methylphenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Methylphenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Nitrophenol	<2,900	<2,300	<2,900	<2,300	<620	<2,800	<2,700	<2,700	<2,500	-	<2,600	<2,500	<2,800	<2,800	<2,600	<2,500	<2,500	<2,400
Pentachlorophenol (Method 8151A)	<14	<12	<15	<12	<13	<14	<13	<14	<13	-	<13	<13	<14	<14	<13	<12	<13	<12
Phenol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2-Nitrophenol	<710	<580	<730	<580	<2,500	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,4-Dinitrophenol	<2,900	<2,300	<2,900	<2,300	<2,500	<2,800	<2,700	<2,700	<2,500	-	<2,600	<2,500	<2,800	<2,800	<2,600	<2,500	<2,500	<2,400
4,6-Dinitro-2-methylphenol	<2,900	<2,300	<2,900	<2,300	<2,500	<2,800	<2,700	<2,700	<2,500	-	<2,600	<2,500	<2,800	<2,800	<2,600	<2,500	<2,500	<2,400
Organics																		
SVOCs: Chlorinated hydrocarbons (ug/kg dry weight, Method 8260C, unless otherwise noted)																		
1,1,1,2-Tetrachloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1,1-Trichloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1,2,2-Tetrachloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1,2-Trichloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1-Dichloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1-Dichloroethene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,1-Dichloropropene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2,3-Trichlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2,3-Trichloropropane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2,4-Trichlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2-Dibromo-3-chloropropane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2-Dichlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2-Dichloroethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,2-Dichloropropane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,3-Dichlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,3-Dichloropropane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1,4-Dichlorobenzene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
1-Chlorohexane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
2-Chloronaphthalene (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
2,2-Dichloropropane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
2-Chlorotoluene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
3,3'-Dichlorobenzidine (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Chlorophenyl phenyl ether (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Chlorotoluene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Bis(2-chloroethoxy) methane (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Bis(2-chloroethyl) ether (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Bis(2-chloroisopropyl) ether (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Hexachlorobenzene (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Hexachlorocyclopentadiene (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Hexachlorobutadiene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Hexachloroethane (Method 8270D)	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
trans-1,2-Dichloroethene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
trans-1,3-Dichloropropene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
trans-1,4-Dichloro-2-butene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Trichloroethene	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18
Trichlorofluoromethane	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18

Preliminary Data - Subject to Revision

Analyte	Copco1 Reservoir - Sediment - Additional Analytes																	
	Sampling Sites																	
	CDH-S-009A(0.0-4.6)	CDH-S-010(0.0-5.0)	CDH-S-010(5.0-8.0)	CDH-S-011(0.0-1.3)	CDH-S-012(0.0-5.4)	CDH-S-013(0.0-5.7)	CDH-S-014(0.0-5.3)	CDH-S-015A(0.0-5.0)	CDH-S-015A(5.0-9.7)	CDH-S-015A(0.0-9.7)	CDH-S-016(0.0-5.0)	CDH-S-016(5.0-7.5)	CDH-S-017(0.0-1.2)	CDH-S-018(0.0-5.0)	CDH-S-018(5.0-8.9)	CDH-S-019(0.0-4.8)	CDH-S-020(0.0-5.0)	CDH-S-020(5.0-7.0)
Organics																		
SVOCs: Other SVOCs (ug/kg dry weight, Method 8270D unless otherwise noted)																		
2-Nitroaniline	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
3-Nitroaniline	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Chloroaniline	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
4-Nitroaniline	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Benzoic acid	<2,900	<2,300	<2,900	<2,300	<2,500	<2,800	<2,700	<2,700	<2,500	-	<2,600	<2,500	<2,800	<2,800	<2,600	<2,500	<2,500	<2,400
Benzyl alcohol	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Dibenzofuran	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Isophorone	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Nitrobenzene	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
N-Nitrosodi-N-propylamine	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
N-Nitrosodiphenylamine	<710	<580	<730	<580	<620	<700	<660	<680	<620	-	<650	<620	<700	<710	<650	<620	<620	<590
Pyridine	<710 L	<580 L	<730 L	<580 L	<620 L	<700 L	<660 L	<680 L	<620 L	-	<650 L	<620 L	<700 L	<710 L	<650 L	<620 L	<620 L	<590 L
Tetrahydrofuran (Method 8260C)	<21	<18	<22	<18	<19	<21	<20	<20	<19	-	<20	<19	<21	<21	<20	<19	<19	<18

Qualifiers:
V: result may vary excessively from the true value
H: result may have a high bias
L: result may have a low bias
T: result obtained past the holding time
U: result determined to be an outlier at the time of data validation
J: result is between the reporting limit and lowest calibration level
- : no data
<: not detected at reporting limit shown
Chlorpyrifos: CDH-S-014(0.0-5.3) by Method GCMS-NCI-SIM and CDH-S-015A(0.0-0.9) by Method MLA-047 Rev 03

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