Sample ID	DOM-2				FO-22				FO-J1				I10-2				JW-12				JW-13				JW-14				JW-6				JW-9			
Sample Date	09/19/01				09/18/01				09/18/01				09/19/01				09/18/01				09/19/01				09/18/01				09/19/01				09/18/01			
Sample Type	N1				N1				N1				N1				N1				N1				N1				N1				N1			
Lab ID	AP22339				AP22328				AP22329				AP22340				AP22331				AP22334				AP22332				AP22335				AP22330			
	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL
SW8260 (UG/L)																																				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	**0.47**	F	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	**7.90**		1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	**0.10**	F	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
SURR 1,2-Dichloroethane-D4	87.2				81.7				82.60				93.10				84.57				87.40				83.82				86.50				84.10			
SURR 4-Bromofluorobenzene	94.2				92.2				92.60				98.20				93.05				94.20				91.93				93.40				92.50			
SURR Dibromofluoromethane	92.6				89.5				89.60				95.00				90.59				91.90				90.92				91.10				91.50			
SURR Toluene-D8	90.2				90.6				90.40				96.50				90.04				90.40				90.75				89.60				89.90			
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	**0.12**	F	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	0.14	U	1	1.	0.14	U	1	1.0	0.14	U	1	1.0	**0.23**	F	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1

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

Tables present all laboratory results for analytes detected above the method detection limit.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 36416

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

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

Sample ID	LS-1				LS-2				LS-3				LS-4				LS-6 POST				LS-6 PRE				LS-6 PRE				LS-7 POST				LS-7 PRE			
Sample Date	09/17/01				09/27/01				09/27/01				09/17/01				09/19/01				09/19/01				09/19/01				09/19/01				09/19/01			
Sample Type	N1				N1				N1				N1				N1				FD1				N1				N1				FD1			
Lab ID	AP22236				AP22601				AP22602				AP22235				AP22343				AP22342				AP22341				AP22346				AP22345			
	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL
SW8260 (UG/L)																																				
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	**0.2**	F	1	0.3	**0.20**	F	1	0.3	**0.13**	F	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	**0.11**	F	1	0.3	**0.10**	F	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.1	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.1	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.1	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.1	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	**0.77**	F	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.1	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
SURR 1,2-Dichloroethane-D4					85.90				84.10								92.60				94.20				92.80				93.80				92.90			
SURR 4-Bromofluorobenzene					94.20				95.90								98.60				100.00				98.30				101.00				98.40			
SURR Dibromofluoromethane					87.60				86.90								94.90				95.30				95.20				94.00				95.70			
SURR Toluene-D8					90.00				91.10								97.00				96.90				96.90				96.60				97.30			
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.2	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.2	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	**0.47**	F	1	1.4	**4.60**		1	1.4	**4.80**		1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	**7.20**		1	1.4	**6.70**		1	1.4	0.11	U	1	1.4	**4.60**		1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**0.37**	F	1	1.0	**0.17**	F	1	1.0	**0.43**	F	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	**0.39**	F	1	1.0	**0.35**	F	1	1.0	0.14	U	1	1.0	**0.83**	F	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1

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

Tables present all laboratory results for analytes detected above the method detection limit.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 36416

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

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

Sample ID	LS-7 PRE				RFR-10				RFR-10 POST H				RFR-6				RFR-7				RFR-8				RFR-9			
Sample Date	09/19/01				09/19/01				10/12/01				09/19/01				09/19/01				09/18/01				09/19/01			
Sample Type	N1				N1				N1				N1				N1				N1				N1			
Lab ID	AP22344				AP22333				AP23211				AP22336				AP22338				AP22327				AP22337			
	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL	Results	Flags	Dilution	SQL
SW8260 (UG/L)																												
Benzene	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Bromobenzene	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Bromochloromethane	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4	0.16	U	1	0.4
Bromodichloromethane	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8	0.12	U	1	0.8
Bromoform	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2	0.14	U	1	1.2
Bromomethane	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1	0.36	U	1	1.1
Butylbenzene, N-	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Butylbenzene, sec-	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3	0.05	U	1	1.3
Butylbenzene, tert-	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4	0.05	U	1	1.4
Carbon tetrachloride	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1	0.09	U	1	2.1
Chlorobenzene	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4	0.09	U	1	0.4
Chloroethane	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0	0.26	U	1	1.0
Chloroform	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3	0.06	U	1	0.3
Chlorohexane, 1-	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6	0.12	U	1	0.6
Chloromethane	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3	0.41	U	1	1.3
Chlorotoluene, 2-	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4	0.12	U	1	0.4
Chlorotoluene, 4-	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6	0.09	U	1	0.6
Dibromo-3-chloropropane, 1,2-	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6	0.72	U	1	2.6
Dibromochloromethane	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5	0.09	U	1	0.5
Dibromomethane	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4	0.10	U	1	2.4
Dichlorobenzene, 1,2-	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3	0.08	U	1	0.3
Dichlorobenzene, 1,3-	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2	0.12	U	1	1.2
Dichlorobenzene, 1,4-	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3	0.09	U	1	0.3
Dichlorodifluoromethane	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0	0.24	U	1	1.0
Dichloroethane, 1,1-	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
Dichloroethane, 1,2-	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6	0.10	U	1	0.6
Dichloroethene, 1,1-	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2	0.16	U	1	1.2
Dichloroethene, cis-1,2-	0.11	U	1	1.2	**0.42**	F	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2	0.11	U	1	1.2
Dichloroethene, trans-1,2-	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6	0.14	U	1	0.6
Dichloropropane, 1,2-	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4	0.14	U	1	0.4
Dichloropropane, 1,3-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Dichloropropane, 2,2-	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5	0.53	U	1	3.5
Dichloropropene, 1,1-	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0	0.12	U	1	1.0
Dichloropropene, cis-1,3-	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0	0.09	U	1	1.0
Dichloropropene, trans-1,3-	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Ethylbenzene	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6	0.06	U	1	0.6
Ethylene dibromide	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6	0.11	U	1	0.6
Hexachlorobutadiene	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1	0.19	U	1	1.1
Isopropylbenzene	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Isopropyltoluene, 4- (Cymene, p-)	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2	0.06	U	1	1.2
Methylene chloride	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0	0.19	U	1	1.0
Naphthalene	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Propylbenzene, N-	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4	0.10	U	1	0.4
Styrene	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4	0.07	U	1	0.4
SURR 1,2-Dichloroethane-D4	93.20				85.20				102.00				87.20				88.80				85.30				86.20			
SURR 4-Bromofluorobenzene	98.40				92.80				109.00				92.30				94.30				92.20				95.10			
SURR Dibromofluoromethane	93.70				91.00				96.70				91.00				92.90				91.60				91.60			
SURR Toluene-D8	95.50				89.20				105.00				89.00				89.80				90.10				90.60			
Tetrachloroethane, 1,1,1,2-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Tetrachloroethane, 1,1,2,2-	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4	0.20	U	1	0.4
Tetrachloroethene	**4.50**		1	1.4	**14.00**		1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4	0.11	U	1	1.4
Toluene	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1	0.11	U	1	1.1
Trichlorobenzene, 1,2,3-	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5	0.13	U	1	0.5
Trichlorobenzene, 1,2,4-	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5	0.08	U	1	0.5
Trichloroethane, 1,1,1-	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8	0.08	U	1	0.8
Trichloroethane, 1,1,2-	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0	0.16	U	1	1.0
Trichloroethene	**0.81**	F	1	1.0	**7.50**		1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0	0.14	U	1	1.0
Trichlorofluoromethane	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8	0.09	U	1	0.8
Trichloropropane, 1,2,3-	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2	0.23	U	1	3.2
Trimethylbenzene, 1,2,4-	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3	0.07	U	1	1.3
Trimethylbenzene, 1,3,5-	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5	0.06	U	1	0.5
Vinyl chloride	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1	0.27	U	1	1.1
Xylene, m,p-	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5	0.14	U	1	0.5
Xylene, o-	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1	0.07	U	1	1.1

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

Tables present all laboratory results for analytes detected above the method detection limit.
All samples were analyzed by APPL Inc.
Referenced laboratory package number: APPL Inc.: 36416

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

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