Volatile Organic Compound Analysis Results for Samples Collected in Nuiqsut, Alaska

Sample Location: Nuiqsut Ambient Air Quality Monitoring Station

Date Sample Collected: 5/9/2022

Analysis Conducted by: Atmospheric Analysis & Consulting, Inc.

Analysis Method: EPA Method TO-15

Laboratory Analysis Report

CLIENT : SLR International Corporation
PROJECT NO : 221010
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/11/2022
DATE REPORTED : 05/12/2022
ANALYST : MB/DL

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	NUI-Primary			Sample Reporting Limit (SRL) (MRLxDF's)	NUI-DUP			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
AAC ID	221010-31123				221010-31124				
Date Sampled	05/09/2022				05/09/2022				
Date Analyzed	05/11/2022				05/11/2022				
Can Dilution Factor	1.68				1.72				
Compound	Result	Qualifier	Analysis DF		Result	Qualifier	Analysis DF		
Chlorodifluoromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Propene	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Dichlorodifluoromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Chloromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Dichlorotetrafluoroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Vinyl Chloride	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Methanol	<SRL	U	1	8.41	<SRL	U	1	8.61	5.00
1,3-Butadiene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Bromomethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Chloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Dichlorofluoromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Ethanol	6.04		1	3.37	25.3		1	3.44	2.00
Vinyl Bromide	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Acetone	<SRL	U	1	3.37	<SRL	U	1	3.44	2.00
Trichlorofluoromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
2-Propanol (IPA)	<SRL	U	1	3.37	<SRL	U	1	3.44	2.00
Acrylonitrile	<SRL	U	1	3.37	<SRL	U	1	3.44	2.00
1,1-Dichloroethene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Methylene Chloride (DCM)	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Allyl Chloride	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Carbon Disulfide	<SRL	U	1	3.37	<SRL	U	1	3.44	2.00
Trichlorotrifluoroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
trans-1,2-Dichloroethene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,1-Dichloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Methyl Tert Butyl Ether (MTBE)	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Vinyl Acetate	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
2-Butanone (MEK)	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
cis-1,2-Dichloroethene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Hexane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Chloroform	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Ethyl Acetate	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Tetrahydrofuran	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2-Dichloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,1,1-Trichloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Benzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50

Laboratory Analysis Report

CLIENT : SLR International Corporation
PROJECT NO : 221010
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 05/11/2022
DATE REPORTED : 05/12/2022
ANALYST : MB/DL

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	NUI-Primary			Sample Reporting Limit (SRL) (MRLxDF's)	NUI-DUP			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
AAC ID	221010-31123				221010-31124				
Date Sampled	05/09/2022				05/09/2022				
Date Analyzed	05/11/2022				05/11/2022				
Can Dilution Factor	1.68				1.72				
Compound	Result	Qualifier	Analysis DF		Result	Qualifier	Analysis DF		
Carbon Tetrachloride	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Cyclohexane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2-Dichloropropane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Bromodichloromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,4-Dioxane	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Trichloroethene (TCE)	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
2,2,4-Trimethylpentane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Heptane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
cis-1,3-Dichloropropene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
4-Methyl-2-pentanone (MiBK)	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
trans-1,3-Dichloropropene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,1,2-Trichloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Toluene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
2-Hexanone (MBK)	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Dibromochloromethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2-Dibromoethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Tetrachloroethene (PCE)	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Chlorobenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Ethylbenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
m & p-Xylene	<SRL	U	1	1.68	<SRL	U	1	1.72	1.00
Bromoform	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Styrene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,1,2,2-Tetrachloroethane	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
o-Xylene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
4-Ethyltoluene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,3,5-Trimethylbenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2,4-Trimethylbenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
Benzyl Chloride (a-Chlorotoluene)	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,3-Dichlorobenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,4-Dichlorobenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2-Dichlorobenzene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
1,2,4-Trichlorobenzene	<SRL	U	1	3.37	<SRL	U	1	3.44	2.00
Hexachlorobutadiene	<SRL	U	1	0.84	<SRL	U	1	0.86	0.50
BFB-Surrogate Std. % Recovery		89%				87%			70-130%

U - Compound was not detected at or above the SRL.

Atmospheric Analysis & Consulting, Inc.

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 05/11/2022 | **INSTRUMENT ID : GC/MS-02**
MATRIX : High Purity N_2 | **CALIBRATION STD ID : MS1-040622-01**
UNITS : PPB (v/v) | **ANALYST : MB/DL**

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Continuing Calibration Verification of the 04/22/2022 Calibration

Analyte Compounds	*Source*[1]	*CCV*[2]	*% Recovery*[3]
4-BFB (surrogate standard)	10.00	9.95	100
Chlorodifluoromethane	10.40	10.20	98.1
Propene	10.60	10.57	100
Dichlorodifluoromethane	10.40	10.04	97
Dimethyl Ether	10.20	10.18	100
Chloromethane	10.40	11.05	106
Dichlorotetrafluoroethane	10.30	9.51	92
Vinyl Chloride	10.50	11.66	111
Acetaldehyde	21.10	21.69	103
Methanol	18.80	19.79	105
1,3-Butadiene	10.60	11.17	105
Bromomethane	10.40	10.09	97
Chloroethane	10.30	10.73	104
Dichlorofluoromethane	10.20	10.56	104
Ethanol	11.20	11.25	100
Vinyl Bromide	10.10	9.51	94
Acrolein	11.10	11.83	107
Acetone	10.60	10.55	100
Trichlorofluoromethane	10.50	9.32	89
2-Propanol (IPA)	11.00	11.13	101
Acrylonitrile	11.20	11.95	107
1,1-Dichloroethene	10.40	9.58	92
Methylene Chloride (DCM)	10.50	10.26	98
TertButanol (TBA)	11.10	12.78	115
Allyl Chloride	10.20	9.62	94
Carbon Disulfide	10.50	11.22	107
Trichlorotrifluoroethane	10.40	9.73	94
trans-1,2-Dichloroethene	10.60	10.00	94
1,1-Dichloroethane	10.50	10.27	98
Methyl Tert Butyl Ether (MTBE)	10.50	9.60	91
Vinyl Acetate	11.00	11.18	102
2-Butanone (MEK)	10.60	10.38	98
cis-1,2-Dichloroethene	10.50	10.29	98
Hexane	10.70	9.81	92
Chloroform	10.60	10.11	95
Ethyl Acetate	10.60	10.73	101
Tetrahydrofuran	10.20	9.11	89
1,2-Dichloroethane	10.50	9.07	86
1,1,1-Trichloroethane	10.40	8.92	86
Benzene	10.60	9.62	91
Carbon Tetrachloride	10.20	8.01	79
Cyclohexane	10.50	8.83	84

Analyte Compounds (Continued)	*Source*[1]	*CCV*[2]	*% Recovery*[3]
1,2-Dichloropropane	10.50	9.52	91
Bromodichloromethane	10.40	8.68	83
1,4-Dioxane	10.40	10.29	99
Trichloroethene (TCE)	10.40	8.43	81
2,2,4-Trimethylpentane	10.00	9.20	92
Methyl Methacrylate	11.00	9.63	88
Heptane	10.50	9.45	90
cis-1,3-Dichloropropene	10.40	8.87	85
4-Methyl-2-pentanone (MiBK)	10.40	10.05	97
trans-1,3-Dichloropropene	10.50	9.26	88
1,1,2-Trichloroethane	10.50	9.31	89
Toluene	10.60	9.50	90
2-Hexanone (MBK)	10.50	10.39	99
Dibromochloromethane	10.30	8.34	81
1,2-Dibromoethane	10.60	9.34	88
Tetrachloroethene (PCE)	10.40	8.57	82
Chlorobenzene	10.60	9.23	87
Ethylbenzene	10.50	9.69	92
m & p-Xylene	21.00	19.20	91
Bromoform	10.50	8.92	85
Styrene	10.50	9.59	91
1,1,2,2-Tetrachloroethane	10.50	10.44	99
o-Xylene	10.50	9.85	94
1,2,3-Trichloropropane	11.00	9.83	89
Isopropylbenzene (Cumene)	10.30	9.83	95
α-Pinene	10.70	9.14	85
2-Chlorotoluene	10.30	8.83	86
n-Propylbenzene	10.10	9.56	95
4-Ethyltoluene	10.30	10.29	100
1,3,5-Trimethylbenzene	10.30	9.82	95
β-Pinene LR	11.00	6.51	59
1,2,4-Trimethylbenzene	10.30	9.95	97
Benzyl Chloride (a-Chlorotoluene)	10.40	9.24	89
1,3-Dichlorobenzene	10.40	9.23	89
1,4-Dichlorobenzene	10.30	9.40	91
Sec-ButylBenzene	10.10	10.01	99
1,2-Dichlorobenzene	10.60	9.92	94
n-ButylBenzene	10.20	10.31	101
1,2-Dibromo-3-Chloropropane	10.10	8.81	87
1,2,4-Trichlorobenzene	11.00	8.25	75
Naphthalene	11.50	9.28	81
Hexachlorobutadiene	11.00	9.33	85

[1] Concentration of analyte compound in certified source standard.

[2] Measured result from daily Continuing Calibration Verification (CCV).

[3] The acceptable range for analyte recovery is 100±30%.

LR - Recovery for this compound was low. Results should be considered estimated.

2225 Sperry Ave., Ventura, CA 93003 www.aaclab.com • (805) 650-1642

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 05/11/2022
MATRIX : High Purity N_2
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-02
CALIBRATION STD ID : MS1-040622-01
ANALYST : MB/DL

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Laboratory Control Spike Analysis

System Monitoring Compounds	*Sample Concentration*	*Spike Added*	*LCS[1] Recovery*	*LCSD[1] Recovery*	*LCS[1] % Recovery[2]*	*LCSD[1] % Recovery[2]*	*RPD[3]*
4-BFB (surrogate standard)	0.0	9.80	9.95	9.73	101.5	99.3	2.2
1,1-Dichloroethene	0.0	10.40	9.58	9.66	92	93	0.8
Methylene Chloride (DCM)	0.0	10.50	10.26	10.90	98	104	6.0
Benzene	0.0	10.60	9.62	9.90	91	93	2.9
Trichloroethene (TCE)	0.0	10.40	8.43	8.80	81	85	4.3
Toluene	0.0	10.60	9.50	9.66	90	91	1.7
Tetrachloroethene (PCE)	0.0	10.40	8.57	8.62	82	83	0.6
Chlorobenzene	0.0	10.60	9.23	9.32	87	88	1.0
Ethylbenzene	0.0	10.50	9.69	9.66	92	92	0.3
m & p-Xylene	0.0	21.00	19.20	18.72	91	89	2.5
o-Xylene	0.0	10.50	9.85	9.98	94	95	1.3

[1] Laboratory Control Spike (LCS) / Laboratory Control Spike Duplicate (LCSD)

[2] The acceptable range for analyte recovery is 100±30%.

[3] Relative Percent Difference (RPD) between LCS recovery and LCSD recovery (acceptable range is <25%).

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QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 05/11/2022 | **INSTRUMENT ID : GC/MS-02**
MATRIX : High Purity He or N_2 | **ANALYST : MB/DL**
UNITS : PPB (v/v)

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Method Blank Analysis

Analyte Compounds	*MB 051122*	*Reporting Limit (RL)*
4-BFB (surrogate standard)	90%	100±30%
Chlorodifluoromethane	<RL	0.5
Propene	<RL	1.0
Dichlorodifluoromethane	<RL	0.5
Dimethyl Ether	<RL	0.5
Chloromethane	<RL	0.5
Dichlorotetrafluoroethane	<RL	0.5
Vinyl Chloride	<RL	0.5
Acetaldehyde	<RL	5.0
Methanol	<RL	5.0
1,3-Butadiene	<RL	0.5
Bromomethane	<RL	0.5
Chloroethane	<RL	0.5
Dichlorofluoromethane	<RL	0.5
Ethanol	<RL	2.0
Vinyl Bromide	<RL	0.5
Acrolein	<RL	1.0
Acetone	<RL	2.0
Trichlorofluoromethane	<RL	0.5
2-Propanol (IPA)	<RL	2.0
Acrylonitrile	<RL	2.0
1,1-Dichloroethene	<RL	0.5
Methylene Chloride (DCM)	<RL	1.0
TertButanol (TBA)	<RL	0.5
Allyl Chloride	<RL	1.0
Carbon Disulfide	<RL	2.0
Trichlorotrifluoroethane	<RL	0.5
trans-1,2-Dichloroethene	<RL	0.5
1,1-Dichloroethane	<RL	0.5
Methyl Tert Butyl Ether (MTBE)	<RL	0.5
Vinyl Acetate	<RL	1.0
2-Butanone (MEK)	<RL	1.0
cis-1,2-Dichloroethene	<RL	0.5
Hexane	<RL	0.5
Chloroform	<RL	0.5
Ethyl Acetate	<RL	0.5
Tetrahydrofuran	<RL	0.5
1,2-Dichloroethane	<RL	0.5
1,1,1-Trichloroethane	<RL	0.5
Benzene	<RL	0.5
Carbon Tetrachloride	<RL	0.5
Cyclohexane	<RL	0.5

Analyte Compounds (Continued)	*MB 051122*	*Reporting Limit (RL)*
1,2-Dichloropropane	<RL	0.5
Bromodichloromethane	<RL	0.5
1,4-Dioxane	<RL	1.0
Trichloroethene (TCE)	<RL	0.5
2,2,4-Trimethylpentane	<RL	0.5
Methyl Methacrylate	<RL	0.5
Heptane	<RL	0.5
cis-1,3-Dichloropropene	<RL	0.5
4-Methyl-2-pentanone (MiBK)	<RL	0.5
trans-1,3-Dichloropropene	<RL	0.5
1,1,2-Trichloroethane	<RL	0.5
Toluene	<RL	0.5
2-Hexanone (MBK)	<RL	1.0
Dibromochloromethane	<RL	0.5
1,2-Dibromoethane	<RL	0.5
Tetrachloroethene (PCE)	<RL	0.5
Chlorobenzene	<RL	0.5
Ethylbenzene	<RL	0.5
m & p-Xylene	<RL	1.0
Bromoform	<RL	0.5
Styrene	<RL	0.5
1,1,2,2-Tetrachloroethane	<RL	0.5
o-Xylene	<RL	0.5
1,2,3-Trichloropropane	<RL	0.5
Isopropylbenzene (Cumene)	<RL	0.5
α-Pinene	<RL	0.5
2-Chlorotoluene	<RL	0.5
n-Propylbenzene	<RL	0.5
4-Ethyltoluene	<RL	0.5
1,3,5-Trimethylbenzene	<RL	0.5
β-Pinene	<RL	0.5
1,2,4-Trimethylbenzene	<RL	0.5
Benzyl Chloride (a-Chlorotoluene)	<RL	0.5
1,3-Dichlorobenzene	<RL	0.5
1,4-Dichlorobenzene	<RL	0.5
Sec-ButylBenzene	<RL	0.5
1,2-Dichlorobenzene	<RL	0.5
n-ButylBenzene	<RL	0.5
1,2-Dibromo-3-Chloropropane	<RL	0.5
1,2,4-Trichlorobenzene	<RL	2.0
Naphthalene	<RL	2.0
Hexachlorobutadiene	<RL	0.5

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 05/11/2022
MATRIX : Air
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-02
ANALYST : MB/DL
DILUTION FACTOR[1] : x14.17

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Duplicate Analysis of AAC Sample ID: 221003-31110

Analyte Compounds	*Sample*	*Duplicate*	*RPD[2]*
4-BFB (surrogate standard)	8.99	9.28	3.2
Chlorodifluoromethane	<SRL	<SRL	*NA*
Propene	619	613	0.9
Dichlorodifluoromethane	<SRL	<SRL	*NA*
Dimethyl Ether	<SRL	<SRL	*NA*
Chloromethane	<SRL	<SRL	*NA*
Dichlorotetrafluoroethane	<SRL	<SRL	*NA*
Vinyl Chloride	<SRL	<SRL	*NA*
Acetaldehyde	172	166	3.6
Methanol	<SRL	<SRL	*NA*
1,3-Butadiene	<SRL	<SRL	*NA*
Bromomethane	<SRL	<SRL	*NA*
Chloroethane	<SRL	<SRL	*NA*
Dichlorofluoromethane	<SRL	<SRL	*NA*
Ethanol	35.9	36.8	2.7
Vinyl Bromide	<SRL	<SRL	*NA*
Acrolein	<SRL	<SRL	*NA*
Acetone	45.1	42.4	6.2
Trichlorofluoromethane	<SRL	<SRL	*NA*
2-Propanol (IPA)	<SRL	<SRL	*NA*
Acrylonitrile	<SRL	<SRL	*NA*
1,1-Dichloroethene	<SRL	<SRL	*NA*
Methylene Chloride (DCM)	<SRL	<SRL	*NA*
TertButanol (TBA)	<SRL	<SRL	*NA*
Allyl Chloride	<SRL	<SRL	*NA*
Carbon Disulfide	<SRL	<SRL	*NA*
Trichlorotrifluoroethane	<SRL	<SRL	*NA*
trans-1,2-Dichloroethene	<SRL	<SRL	*NA*
1,1-Dichloroethane	<SRL	<SRL	*NA*
Methyl Tert Butyl Ether (MTBE)	<SRL	<SRL	*NA*
Vinyl Acetate	<SRL	<SRL	*NA*
2-Butanone (MEK)	24.4	24.0	1.8
cis-1,2-Dichloroethene	<SRL	<SRL	*NA*
Hexane	13.6	13.7	1.0
Chloroform	<SRL	<SRL	*NA*
Ethyl Acetate	<SRL	<SRL	*NA*
Tetrahydrofuran	<SRL	<SRL	*NA*
1,2-Dichloroethane	<SRL	<SRL	*NA*
1,1,1-Trichloroethane	<SRL	<SRL	*NA*
Benzene	<SRL	<SRL	*NA*
Carbon Tetrachloride	<SRL	<SRL	*NA*
Cyclohexane	<SRL	<SRL	*NA*

Analyte Compounds (Continued)	*Sample*	*Duplicate*	*RPD[2]*
1,2-Dichloropropane	<SRL	<SRL	*NA*
Bromodichloromethane	<SRL	<SRL	*NA*
1,4-Dioxane	<SRL	<SRL	*NA*
Trichloroethene (TCE)	<SRL	<SRL	*NA*
2,2,4-Trimethylpentane	<SRL	<SRL	*NA*
Methyl Methacrylate	<SRL	<SRL	*NA*
Heptane	51.4	46.9	9.2
cis-1,3-Dichloropropene	<SRL	<SRL	*NA*
4-Methyl-2-pentanone (MiBK)	36.3	35.3	2.8
trans-1,3-Dichloropropene	<SRL	<SRL	*NA*
1,1,2-Trichloroethane	<SRL	<SRL	*NA*
Toluene	329	314	4.5
2-Hexanone (MBK)	<SRL	<SRL	*NA*
Dibromochloromethane	<SRL	<SRL	*NA*
1,2-Dibromoethane	<SRL	<SRL	*NA*
Tetrachloroethene (PCE)	<SRL	<SRL	*NA*
Chlorobenzene	<SRL	<SRL	*NA*
Ethylbenzene	33.3	32.9	1.3
m & p-Xylene	<SRL	<SRL	*NA*
Bromoform	<SRL	<SRL	*NA*
Styrene	<SRL	<SRL	*NA*
1,1,2,2-Tetrachloroethane	<SRL	<SRL	*NA*
o-Xylene	<SRL	<SRL	*NA*
1,2,3-Trichloropropane	<SRL	<SRL	*NA*
Isopropylbenzene (Cumene)	<SRL	<SRL	*NA*
α-Pinene	107	97.7	8.9
2-Chlorotoluene	<SRL	<SRL	*NA*
n-Propylbenzene	<SRL	<SRL	*NA*
4-Ethyltoluene	11.5	12.6	9.4
1,3,5-Trimethylbenzene	8.50	8.22	3.4
β-Pinene	122	107	12.6
1,2,4-Trimethylbenzene	<SRL	<SRL	*NA*
Benzyl Chloride (a-Chlorotoluene)	<SRL	<SRL	*NA*
1,3-Dichlorobenzene	<SRL	<SRL	*NA*
1,4-Dichlorobenzene	12.6	11.8	7.0
Sec-ButylBenzene	<SRL	<SRL	*NA*
1,2-Dichlorobenzene	<SRL	<SRL	*NA*
n-ButylBenzene	<SRL	<SRL	*NA*
1,2-Dibromo-3-Chloropropane	<SRL	<SRL	*NA*
1,2,4-Trichlorobenzene	<SRL	<SRL	*NA*
Naphthalene	<SRL	<SRL	*NA*
Hexachlorobutadiene	<SRL	<SRL	*NA*

[1] Dilution factor is the product of the Canister Dilution Factor and the Analysis Dilution Factor.

[2] Relative Percent Difference (RPD) between Sample analysis and Duplicate analysis (acceptable range is <25%).

SRL - Sample Reporting Limit (minimum)