Atmospheric Analysis & Consulting, Inc.

CLIENT	: Yorke Engineering
PROJECT NAME	: Odor Sampling & Analysis
PROJECT NO.	: 0357-007-01
AAC PROJECT NO.	: 202141
REPORT DATE	: 12/4/2020

On November 24, 2020, Atmospheric Analysis & Consulting, Inc. received one (1) six-Liter Silonite Canister for Volatile Organic Compounds analysis by EPA Method TO-15. Upon receipt, the sample was assigned a unique Laboratory ID number as follows:

Client ID	Lab ID	Return Pressure (mmHga)
Outside AAA	202141-14807	565.6

This analysis is accredited under the laboratory's ISO/IEC 17025:2017 accreditation issued by the ANSI National Accreditation Board. Refer to certificate and scope of accreditation AT-1908. For detailed information pertaining to specific EPA, NCASI, ASTM and SCAQMD accreditations (Methods & Analytes), please visit our website at www.aaclab.com.

I certify that this data is technically accurate, complete, and in compliance with the terms and conditions of the contract. No problems were encountered during receiving, preparation, and/or analysis of this sample.

The Technical Director or his designee, as verified by the following signature, has authorized release of the data contained in this hardcopy report.

If you have any questions or require further explanation of data results, please contact the undersigned.

Sucha Parmar, Ph.D.
Technical Director

This report consists of 8 pages.

Laboratory Analysis Report

CLIENT : Yorke Engineering
PROJECT NO : 202141
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 11/24/2020
DATE REPORTED : 12/04/2020
ANALYST : MB/RC

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	Outside AAA			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
AAC ID	202141-14807				
Date Sampled	11/19/2020				
Date Analyzed	12/03/2020				
Canister Dilution Factor	1.80				
Compound	Result	Qualifier	Analysis DF		
Chlorodifluoromethane	<SRL	U	1	0.90	0.5
Propene	<SRL	U	1	1.80	1.0
Dichlorodifluoromethane	<SRL	U	1	0.90	0.5
Dimethyl Ether	<SRL	U	1	0.90	0.5
Chloromethane	<SRL	U	1	0.90	0.5
Dichlorotetrafluoroethane	<SRL	U	1	0.90	0.5
Vinyl Chloride	<SRL	U	1	0.90	0.5
Acetaldehyde	8.07		1	3.60	2.0
Methanol	10.8		1	9.00	5.0
1,3-Butadiene	<SRL	U	1	0.90	0.5
Bromomethane	0.94		1	0.90	0.5
Chloroethane	<SRL	U	1	0.90	0.5
Dichlorofluoromethane	<SRL	U	1	0.90	0.5
Ethanol	14.0		1	3.60	2.0
Vinyl Bromide	<SRL	U	1	0.90	0.5
Acrolein	<SRL	U	1	1.80	1.0
Acetone	9.34		1	3.60	2.0
Trichlorofluoromethane	<SRL	U	1	0.90	0.5
2-Propanol (IPA)	<SRL	U	1	3.60	2.0
Acrylonitrile	<SRL	U	1	3.60	2.0
1,1-Dichloroethene	<SRL	U	1	0.90	0.5
Methylene Chloride (DCM)	<SRL	U	1	1.80	1.0
TertButanol (TBA)	<SRL	U	1	0.90	0.5
Allyl Chloride	<SRL	U	1	1.80	1.0
Carbon Disulfide	<SRL	U	1	3.60	2.0
Trichlorotrifluoroethane	<SRL	U	1	0.90	0.5
trans-1,2-Dichloroethene	<SRL	U	1	0.90	0.5
1,1-Dichloroethane	<SRL	U	1	0.90	0.5
Methyl Tert Butyl Ether (MTBE)	<SRL	U	1	0.90	0.5
Vinyl Acetate	<SRL	U	1	1.80	1.0
2-Butanone (MEK)	<SRL	U	1	1.80	1.0
cis-1,2-Dichloroethene	<SRL	U	1	0.90	0.5
Hexane	<SRL	U	1	0.90	0.5
Chloroform	<SRL	U	1	0.90	0.5
Ethyl Acetate	<SRL	U	1	0.90	0.5
Tetrahydrofuran	<SRL	U	1	0.90	0.5
1,2-Dichloroethane	<SRL	U	1	0.90	0.5

Laboratory Analysis Report

CLIENT : Yorke Engineering
PROJECT NO : 202141
MATRIX : AIR
UNITS : PPB (v/v)

DATE RECEIVED : 11/24/2020
DATE REPORTED : 12/04/2020
ANALYST : MB/RC

VOLATILE ORGANIC COMPOUNDS BY EPA TO-15

Client ID	Outside AAA			Sample Reporting Limit (SRL) (MRLxDF's)	Method Reporting Limit (MRL)
AAC ID	202141-14807				
Date Sampled	11/19/2020				
Date Analyzed	12/03/2020				
Canister Dilution Factor	1.80				
Compound	Result	Qualifier	Analysis DF		
1,1,1-Trichloroethane	<SRL	U	1	0.90	0.5
Benzene	0.97		1	0.90	0.5
Carbon Tetrachloride	<SRL	U	1	0.90	0.5
Cyclohexane	<SRL	U	1	0.90	0.5
1,2-Dichloropropane	<SRL	U	1	0.90	0.5
Bromodichloromethane	<SRL	U	1	0.90	0.5
1,4-Dioxane	<SRL	U	1	1.80	1.0
Trichloroethene (TCE)	<SRL	U	1	0.90	0.5
2,2,4-Trimethylpentane	<SRL	U	1	0.90	0.5
Methyl Methacrylate	<SRL	U	1	0.90	0.5
Heptane	<SRL	U	1	0.90	0.5
cis-1,3-Dichloropropene	<SRL	U	1	0.90	0.5
4-Methyl-2-pentanone (MiBK)	<SRL	U	1	0.90	0.5
trans-1,3-Dichloropropene	<SRL	U	1	0.90	0.5
1,1,2-Trichloroethane	<SRL	U	1	0.90	0.5
Toluene	<SRL	U	1	0.90	0.5
2-Hexanone (MBK)	<SRL	U	1	1.80	1.0
Dibromochloromethane	<SRL	U	1	0.90	0.5
1,2-Dibromoethane	<SRL	U	1	0.90	0.5
Tetrachloroethene (PCE)	<SRL	U	1	0.90	0.5
Chlorobenzene	<SRL	U	1	0.90	0.5
Ethylbenzene	<SRL	U	1	0.90	0.5
m & p-Xylene	<SRL	U	1	1.80	1.0
Bromoform	<SRL	U	1	0.90	0.5
Styrene	<SRL	U	1	0.90	0.5
1,1,2,2-Tetrachloroethane	<SRL	U	1	0.90	0.5
o-Xylene	<SRL	U	1	0.90	0.5
1,2,3-Trichloropropane	<SRL	U	1	0.90	0.5
Isopropylbenzene (Cumene)	<SRL	U	1	0.90	0.5
α-Pinene	<SRL	U	1	0.90	0.5
2-Chlorotoluene	<SRL	U	1	0.90	0.5
n-Propylbenzene	<SRL	U	1	0.90	0.5
4-Ethyltoluene	<SRL	U	1	0.90	0.5
1,3,5-Trimethylbenzene	<SRL	U	1	0.90	0.5
β-Pinene	<SRL	U	1	0.90	0.5
1,2,4-Trimethylbenzene	<SRL	U	1	0.90	0.5
Benzyl Chloride (a-Chlorotoluene)	<SRL	U	1	0.90	0.5
1,3-Dichlorobenzene	<SRL	U	1	0.90	0.5
1,4-Dichlorobenzene	<SRL	U	1	0.90	0.5
Sec-ButylBenzene	<SRL	U	1	0.90	0.5
1,2-Dichlorobenzene	<SRL	U	1	0.90	0.5
n-ButylBenzene	<SRL	U	1	1.80	1.0
1,2-Dibromo-3-Chloropropane	<SRL	U	1	0.90	0.5
1,2,4-Trichlorobenzene	<SRL	U	1	0.90	0.5
Naphthalene	<SRL	U	1	1.80	1.0
Hexachlorobutadiene	<SRL	U	1	0.90	0.5
BFB-Surrogate Std. % Recovery		101%			70-130%

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

Sucha Parmar, Ph.D.
Technical Director

Atmospheric Analysis & Consulting, Inc.

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 12/03/2020
MATRIX : High Purity N_2
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-03
CALIBRATION STD ID : PS101520-02
ANALYST : MB/RC

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Continuing Calibration Verification of the 11/04/2020 Calibration

Analyte Compounds	*Source*[1]	*CCV*[2]	*% Recovery*[3]
4-BFB (surrogate standard)	10.00	10.81	108
Chlorodifluoromethane	10.70	9.53	89
Propene LR	11.00	7.69	70
Dichlorodifluoromethane	10.30	9.57	93
Dimethyl Ether	10.70	9.24	86
Chloromethane LR	10.60	5.36	51
Dichlorotetrafluoroethane	10.20	9.78	96
Vinyl Chloride	10.50	9.95	95
Acetaldehyde	19.80	20.25	102
Methanol	16.80	15.62	93
1,3-Butadiene	10.90	10.27	94
Bromomethane	10.60	11.74	111
Chloroethane	10.20	9.63	94
Dichlorofluoromethane	10.40	10.37	100
Ethanol	10.20	9.33	91
Vinyl Bromide	10.60	10.53	99
Acrolein	11.00	9.49	86
Acetone	10.50	8.75	83
Trichlorofluoromethane	10.50	10.18	97
2-Propanol (IPA)	9.80	8.19	84
Acrylonitrile	11.50	10.59	92
1,1-Dichloroethene	10.80	10.53	98
Methylene Chloride (DCM)	10.90	10.03	92
TertButanol (TBA)	10.50	9.65	92
Allyl Chloride	10.60	10.00	94
Carbon Disulfide	10.20	10.51	103
Trichlorotrifluoroethane	10.90	10.46	96
trans-1,2-Dichloroethene	10.30	10.20	99
1,1-Dichloroethane	10.40	9.71	93
Methyl Tert Butyl Ether (MTBE)	11.00	10.43	95
Vinyl Acetate	11.00	10.22	93
2-Butanone (MEK)	10.60	10.53	99
cis-1,2-Dichloroethene	10.70	10.58	99
Hexane	11.00	10.21	93
Chloroform	10.80	10.21	95
Ethyl Acetate	10.80	10.26	95
Tetrahydrofuran	10.30	9.20	89
1,2-Dichloroethane	10.80	9.89	92
1,1,1-Trichloroethane	10.70	10.04	94
Benzene	10.80	9.19	85
Carbon Tetrachloride	10.80	9.94	92
Cyclohexane	10.80	10.03	93

Analyte Compounds (Continued)	*Source*[1]	*CCV*[2]	*% Recovery*[3]
1,2-Dichloropropane	10.80	10.50	97
Bromodichloromethane	9.90	9.66	98
1,4-Dioxane	9.90	8.54	86
Trichloroethene (TCE)	9.90	9.49	96
2,2,4-Trimethylpentane	10.70	10.24	96
Methyl Methacrylate	10.60	10.38	98
Heptane	10.80	10.50	97
cis-1,3-Dichloropropene	10.40	10.04	97
4-Methyl-2-pentanone (MiBK)	10.40	10.39	100
trans-1,3-Dichloropropene	10.20	9.84	96
1,1,2-Trichloroethane	10.90	10.42	96
Toluene	11.00	10.00	91
2-Hexanone (MBK)	10.10	9.99	99
Dibromochloromethane	10.40	10.73	103
1,2-Dibromoethane	10.90	10.80	99
Tetrachloroethene (PCE)	10.60	10.28	97
Chlorobenzene	10.80	10.23	95
Ethylbenzene	10.90	10.41	96
m & p-Xylene	21.20	21.98	104
Bromoform	10.60	10.50	99
Styrene	10.80	11.03	102
1,1,2,2-Tetrachloroethane	10.70	10.66	100
o-Xylene	10.70	10.59	99
1,2,3-Trichloropropane	10.70	10.80	101
Isopropylbenzene (Cumene)	10.70	10.84	101
α-Pinene	11.60	11.44	99
2-Chlorotoluene	10.70	10.51	98
n-Propylbenzene	10.10	10.97	109
4-Ethyltoluene	10.70	11.31	106
1,3,5-Trimethylbenzene	10.60	10.70	101
β-Pinene HR	9.30	12.84	138
1,2,4-Trimethylbenzene	10.50	10.30	98
Benzyl Chloride (a-Chlorotoluene)	10.20	11.22	110
1,3-Dichlorobenzene	10.20	11.16	109
1,4-Dichlorobenzene	10.60	11.54	109
Sec-ButylBenzene	10.70	11.25	105
1,2-Dichlorobenzene	10.60	11.35	107
n-ButylBenzene	10.30	11.71	114
1,2-Dibromo-3-Chloropropane	10.20	10.46	103
1,2,4-Trichlorobenzene	10.90	9.98	92
Naphthalene	11.00	10.55	96
Hexachlorobutadiene	10.90	12.21	112

[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%.

HR - Recovery for this compound was high. Results should be consider biased high.

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

Sucha Parmar, Ph.D.
Technical Director

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 12/03/2020
MATRIX : High Purity N_2
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-03
CALIBRATION STD ID : PS101520-02
ANALYST : MB/RC

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	10.00	10.81	11.00	108.1	110	1.7
1,1-Dichloroethene	0.0	10.80	10.53	10.79	98	100	2.4
Methylene Chloride (DCM)	0.0	10.90	10.03	10.15	92	93	1.2
Benzene	0.0	10.80	9.19	9.86	85	91	7.0
Trichloroethene (TCE)	0.0	9.90	9.49	9.99	96	101	5.1
Toluene	0.0	11.00	10.00	10.88	91	99	8.4
Tetrachloroethene (PCE)	0.0	10.60	10.28	10.94	97	103	6.2
Chlorobenzene	0.0	10.80	10.23	10.35	95	96	1.2
Ethylbenzene	0.0	10.90	10.41	10.48	96	96	0.7
m & p-Xylene	0.0	21.20	21.98	21.24	104	100	3.4
o-Xylene	0.0	10.70	10.59	10.49	99	98	0.9

[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%).

Sucha Parmar, Ph.D.
Technical Director

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 12/03/2020
MATRIX : High Purity He or N_2
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-03
ANALYST : MB/RC

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Method Blank Analysis

Analyte Compounds	*MB 120320*	*Reporting Limit (RL)*
4-BFB (surrogate standard)	94%	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	2.0
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 120320*	*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	1.0
1,2-Dibromo-3-Chloropropane	<RL	0.5
1,2,4-Trichlorobenzene	<RL	0.5
Naphthalene	<RL	1.0
Hexachlorobutadiene	<RL	0.5

Sucha Parmar, Ph.D.
Technical Director

Atmospheric Analysis & Consulting, Inc.

QUALITY CONTROL / QUALITY ASSURANCE REPORT

ANALYSIS DATE : 12/03/2020
MATRIX : Air
UNITS : PPB (v/v)

INSTRUMENT ID : GC/MS-03
ANALYST : MB/RC
DILUTION FACTOR[1] : x23.41

VOLATILE ORGANIC COMPOUNDS BY EPA METHOD TO-15

Duplicate Analysis of AAC Sample ID: 202153-14898

Analyte Compounds	*Sample*	*Duplicate*	*RPD[2]*
4-BFB (surrogate standard)	9.30	9.82	5.4
Chlorodifluoromethane	<SRL	<SRL	*NA*
Propene	1570.00	1490.00	5.2
Dichlorodifluoromethane	<SRL	<SRL	*NA*
Dimethyl Ether	<SRL	<SRL	*NA*
Chloromethane	<SRL	<SRL	*NA*
Dichlorotetrafluoroethane	<SRL	<SRL	*NA*
Vinyl Chloride	<SRL	<SRL	*NA*
Acetaldehyde	<SRL	<SRL	*NA*
Methanol	<SRL	<SRL	*NA*
1,3-Butadiene	<SRL	<SRL	*NA*
Bromomethane	<SRL	<SRL	*NA*
Chloroethane	<SRL	<SRL	*NA*
Dichlorofluoromethane	<SRL	<SRL	*NA*
Ethanol	<SRL	<SRL	*NA*
Vinyl Bromide	<SRL	<SRL	*NA*
Acrolein	<SRL	<SRL	*NA*
Acetone	<SRL	<SRL	*NA*
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)	<SRL	<SRL	*NA*
cis-1,2-Dichloroethene	<SRL	<SRL	*NA*
Hexane	<SRL	<SRL	*NA*
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	<SRL	<SRL	*NA*
cis-1,3-Dichloropropene	<SRL	<SRL	*NA*
4-Methyl-2-pentanone (MiBK)	<SRL	<SRL	*NA*
trans-1,3-Dichloropropene	<SRL	<SRL	*NA*
1,1,2-Trichloroethane	<SRL	<SRL	*NA*
Toluene	<SRL	<SRL	*NA*
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	<SRL	<SRL	*NA*
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	<SRL	<SRL	*NA*
2-Chlorotoluene	<SRL	<SRL	*NA*
n-Propylbenzene	<SRL	<SRL	*NA*
4-Ethyltoluene	<SRL	<SRL	*NA*
1,3,5-Trimethylbenzene	<SRL	<SRL	*NA*
β-Pinene	<SRL	<SRL	*NA*
1,2,4-Trimethylbenzene	<SRL	<SRL	*NA*
Benzyl Chloride (a-Chlorotoluene)	<SRL	<SRL	*NA*
1,3-Dichlorobenzene	<SRL	<SRL	*NA*
1,4-Dichlorobenzene	<SRL	<SRL	*NA*
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)

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CHAIN OF CUSTODY/ ANALYSIS REQUEST FORM

Client Name Ninyo & Moore	Project Name Odor Sampling & Analysis					Analysis Requested				Send report:
Project Mgr. (Print Name) Keith Gilbert	Project Number 0357-007-01					EPA TO-15	SCAQMD 307-91			
Sampler's Name (Print Name) Bipul K. Saraf	Sampler's Signature Bipul K. Saraf									Attn: Bipul Saraf E-Mail – Bsaraf@YorkeEngr.com Phone#: 949-444-8063 Fax#_____
AAC Sample No.	Date Sampled	Time Sampled	Sample Type	Client Sample ID/Description	Type/No. of Containers					
141807	11/19-11/20	16:55	Canister ID 1156	Outside AAA	1	X	X			Send invoice to: Attn: Accounting@YorkeEngr.com P.O. #
										Turnaround Time 24-Hr ____ 48-Hr ____ 5 Day ____ Normal _X_ Other (Specify) ____
										Special Instructions/remarks: Please report all compounds including aldehyde per TO-15. Include all compounds per SCAQMD 307-91.
Relinquished by (Signature): Bipul K. Saraf	Print Name: Bipul Saraf				Date/Time 11/23/2020 12:30 PM	Received by (signature):				Print Name
Relinquished by (Signature):	Print Name:				Date/Time 11/24/20 1034	Received by (signature):				Print Name Gabriel Rojas

1x can + 1x Fritech