(QT Reviewed) Quantitation Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025132.D

: 11 Nov 2021 15:36 Acq On

: JC/MD Operator : VSTD00535 Sample

: 5.0mL/MSVOA_X/WATER Misc Sample Multiplier: 1 ALS Vial : 8

Quant Time: Nov 11 16:12:44 2021

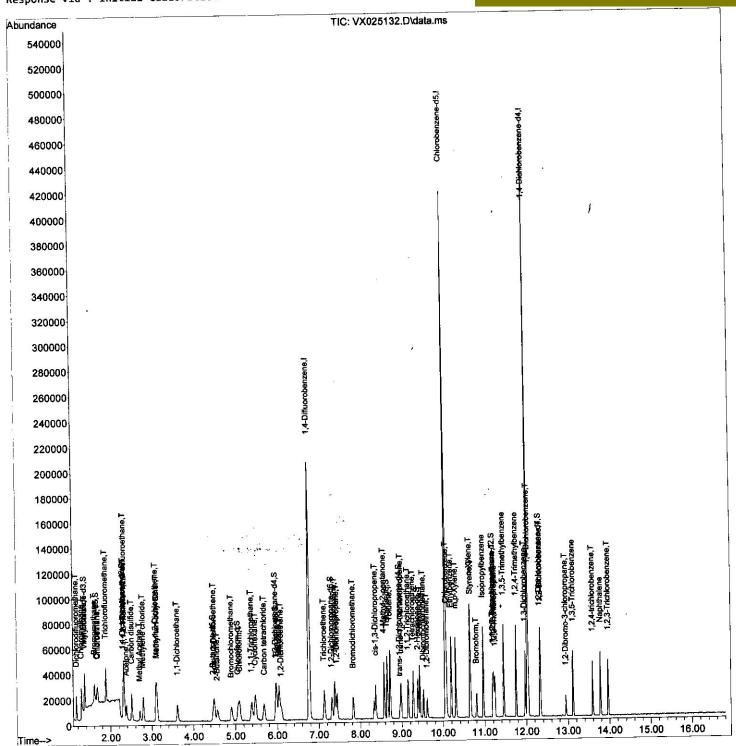
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 14:40:35 2021 Response via : Initial Calibration

Instrument: MSVOA_X **ClientSampleld**: STD005635

Manual IntegrationsAPPROVED



SFAMXLM111121WMA.M Thu Nov 18 03:41:32 2021

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025132.D

Acq On : 11 Nov 2021 15:36

Operator : JC/MD Sample : VSTD00535

Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

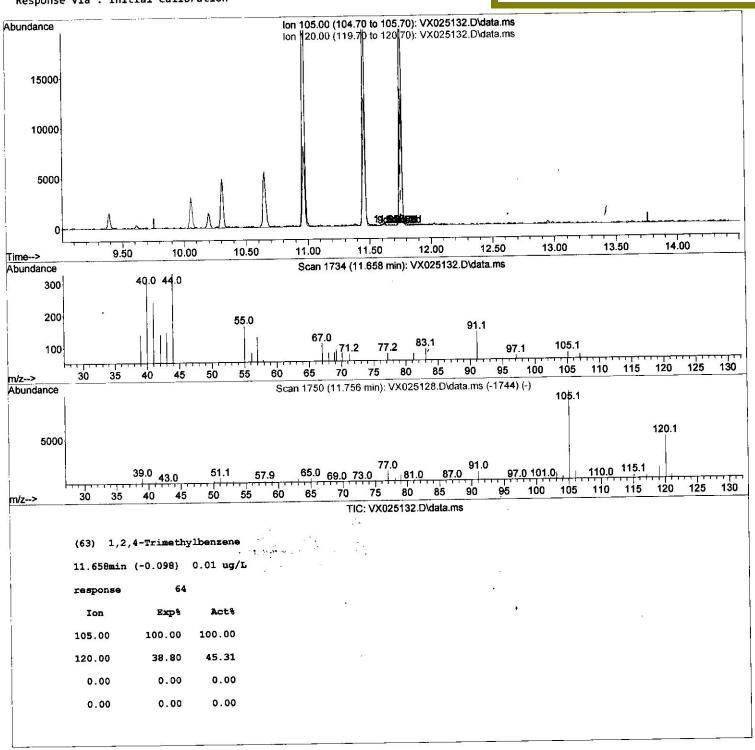
Quant Time: Nov 18 03:42:24 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis OLast Update : Fri Nov 12 12:01:23

QLast Update : Fri Nov 12 12:01:23 2021 Response via : Initial Calibration Instrument: MSVOA_X ClientSampleId: VSTD005635

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File: VX025132.D

Acq On : 11 Nov 2021 15:36

Operator : JC/MD Sample : VSTD00535

Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

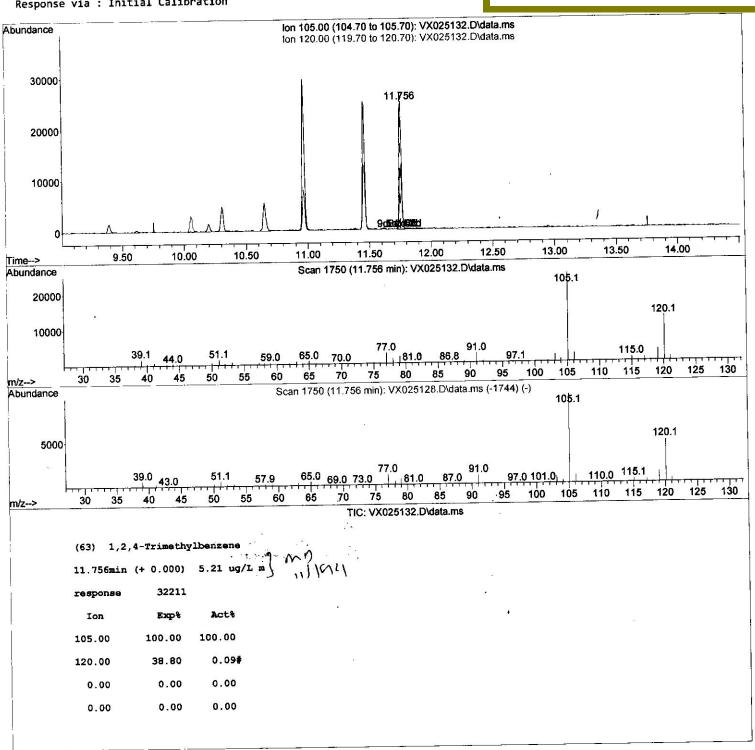
Quant Time: Nov 11 16:12:44 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 14:40:35 2021 Response via : Initial Calibration Instrument: MSVOA_X ClientSampleId: VSTD005635

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File: VX025132.D

Acq On : 11 Nov 2021 15:36

Operator : JC/MD Sample : VSTD00535

Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

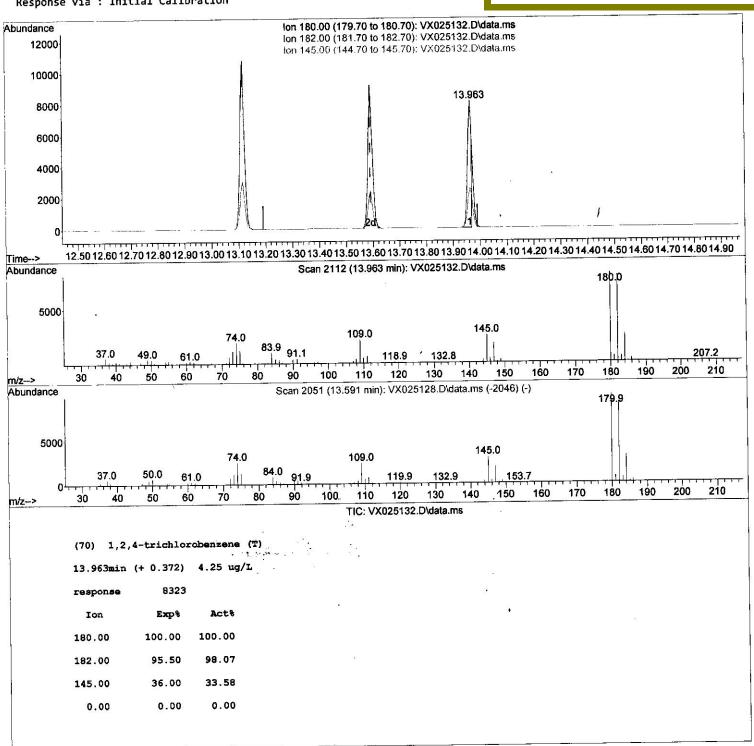
Quant Time: Nov 18 03:42:24 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Fri Nov 12 12:01:23 2021 Response via : Initial Calibration Instrument:
MSVOA_X
ClientSampleId:
VSTD005635

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025132.D

: 11 Nov 2021 15:36 Acq On

Operator : JC/MD : VSTD00535 Sample

: 5.0mL/MSVOA_X/WATER Misc Sample Multiplier: 1 : 8 ALS Vial

Quant Time: Nov 11 16:12:44 2021

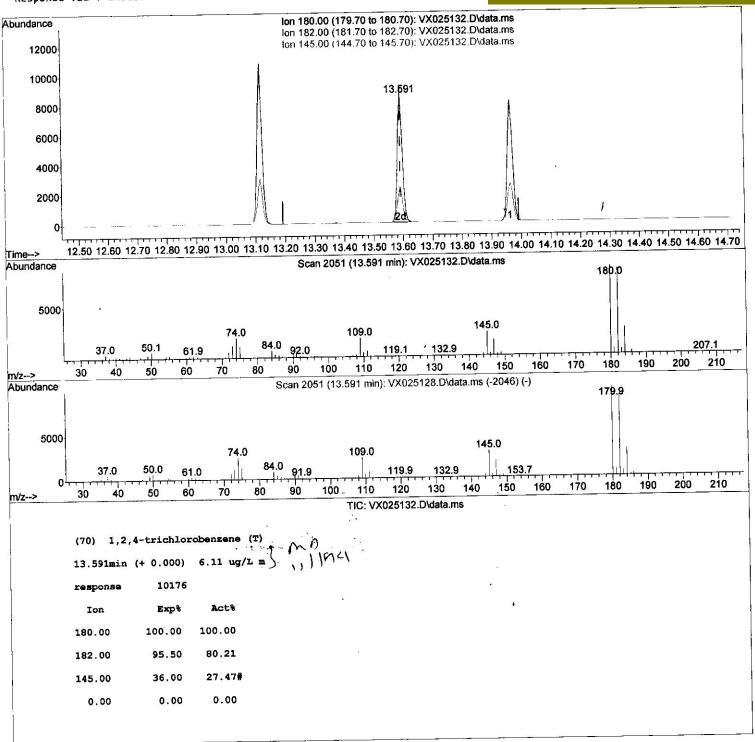
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 14:40:35 2021 Response via : Initial Calibration

Instrument: MSVOA_X ClientSampleId: /STD00<u>5635</u>

Manual IntegrationsAPPROVED



(QT Reviewed) Quantitation Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025132.D

: 11 Nov 2021 15:36 Acq On

Operator : JC/MD Sample : VSTD00535

: 5.0mL/MSVOA_X/WATER Misc ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 11 16:12:44 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis QLast Update : Thu Nov 11 14:40:35 2021 Response via : Initial Calibration

nstrument :
ISVOA_X
lientSampleId :
STD005635

Manual IntegrationsAPPROVED

Compound	R.T. ([]Ion	Response	Conc Units De	v(Min)
<pre>Internal Standards 1) 1,4-Difluorobenzene</pre>	6.769	114	221023	50.00 ug/L	0.00
28) Chlorobenzene-d5	10.055	117	197310	50.00 ug/L	0.00
58) 1,4-Dichlorobenzene-d4	12.024	152	95130	50.00 ug/L	0.00
58) 1,4-DICHIO ODERZENC 44				 20	
System Monitoring Compounds					
4) Vinyl Chloride-d3	1.368	65	8391	4.43 ug/L	0.00
7) Chloroethane-d5	1.666	69	6490	5.85 ug/L	0.00
11) 1,1-Dichloroethene-d2	2.306	63	14133	3.91 ug/L	0.00
21) 2-Butanone-d5	4.471	46	12147	9.15 ug/L	0.00
24) Chloroform-d	5.062	84	14175	3.86 ug/L	0.00
26) 1,2-Dichloroethane-d4	5.964	65	9027	3.81 ug/L	0.00
32) Benzene-d6	5.983	84	29486	4.49 ug/L	0.00 0.00
36) 1,2-Dichloropropane-d6	7.312	67	8949	4.43 ug/L	0.00
41) Toluene-d8	8.653	98	28450	4.89 ug/L	0.00
43) trans-1,3-Dichloroprop	8.952	79	5021	4.79 ug/L 10.70 ug/L	0.00
47) 2-Hexanone-d5	9.391	63	9337	4.43 ug/L	0.00
56) 1,1,2,2-Tetrachloroeth	11.195	84	12600	5.96 ug/L	0.00
66) 1,2-Dichlorobenzene-d4	12.323	152	10991	3.90 ug/L	0.00
				İ	Qvalue
Target Compounds	1.166	85	9552	4.39 ug/L	100
2) Dichlorodifluoromethane	1.288	50	10393	6.29 ug/L	89
3) Chloromethane	1.374	62	10822	5.74 ug/L	96
5) Vinyl chloride	1.611	94	5168	4.69 ug/L	96
6) Bromomethane	1.691	64	6322	5.99 ug/L	99
8) Chloroethane9) Trichlorofluoromethane	1.886	101	15282	4.57 ug/L	97
10) 1,1,2-Trichloro-1,2,2	2.331	101	7803	4.51 ug/L	95
12) 1,1-Dichloroethene	2.319	96	7396	4.80 ug/L	91
13) Acetone	2.386	43	11071	8.58 ug/L	95
14) Carbon disulfide	2.514	76	25783	6.26 ug/L	98
15) Methyl Acetate	2.709	43	9703	5.29 ug/L	
16) Methylene chloride	2.788	84	8869	5.15 ug/L	83
17) trans-1,2-Dichloroethene	3.093	96	8583	5.42 ug/L	89
18) Methyl tert-butyl Ether	3.117	73	25746	4.35 ug/L	
19) 1,1-Dichloroethane	3.611	63	13802		92
20) cis-1,2-Dichloroethene	4.489		9456	5.20 ug/L	88
22) 2-Butanone	4.562				86 # 72
23) Bromochloromethane	4.904		4766	5.48 ug/L	# /2 95 ·
25) Chloroform	5.099		14189	4.09 ug/L 3.88 ug/L	
27) 1,2-Dichloroethane	6.092		10263		90
29) Cyclohexane	5.477		13680	4.80 ug/L	94
30) 1,1,1-Trichloroethane	5.385		13875 11705	4.24 ug/L 4.44 ug/L	96
31) Carbon tetrachloride	5.690			4.71 ug/L	100
33) Benzene	6.044			(T).	82
34) Trichloroethene	7.129			5.09 ug/L	94
35) Methylcyclohexane	7.385			4.98 ug/L	
37) 1,2-Dichloropropane	7.434 7.830			4.76 ug/L	
38) Bromodichloromethane	8.372			4.92 ug/L	
39) cis-1,3-Dichloropropene	8.574			10.14 ug/L	
40) 4-Methyl-2-pentanone	8.720			5.02 ug/L	
42) Toluene	3.720				

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025132.D

Acq On : 11 Nov 2021 15:36

Operator : JC/MD Sample : VSTD00535

: 5.0mL/MSVOA_X/WATER Misc ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 11 16:12:44 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 14:40:35 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response (Conc Units Dev(Min)
44) trans-1,3-Dichloropropene	8.982	75	13013	4.93 ug/L	96
45) 1,1,2-Trichloroethane	9.153	97	8635	5.68 ug/L	97
46) Tetrachloroethene	9.275	164	7180	6.74 ug/L	88
48) 2-Hexanone	9.433	43	20844	9.84 ug/L #	85
49) Dibromochloromethane	9.525	129	9502	5.80 ug/L	99
50) 1,2-Dibromoethane	9.610	107	8897	5.46 ug/L #	95
51) Chlorobenzene	10.079	112	23035	5.63 ug/L	96
52) Ethylbenzene	10.195	91	37335	5.00 ug/L	91
53) m,p-Xylene	10.305	106	15042	5.43 ug/L	81
54) o-Xylene	10.646	106	15216	5.74 ug/L	81
55) Styrene	10.659	104	2492 9	5.44 ug/L	86
57) 1,1,2,2-Tetrachloroethane	11.213	83	13167	4.68 ug/L	93
59) Bromoform	10.805	173	7087	6.39 ug/L #	95
60) Isopropylbenzene	10.963	105	37801	5.08 ug/L	96
61) 1,2,3-Trichloropropane	11.244	75	10048	4.61 ug/L	94
62) 1,3,5-Trimethylbenzene	11.451	105	32231 7	5.08 ug/L	89
63) 1,2,4-Trimethylbenzene	11.756	105	32211m /	5.21 ug/L	
64) 1,3-Dichlorobenzene	11.969	146	17443	6.09 ug/L	97
65) 1,4-Dichlorobenzene	12.043	146	17195	5.90 ug/L	95
67) 1,2-Dichlorobenzene	12.335	146	16949	5.78 ug/L	92
68) 1,2-Dibromo-3-chloropr	12.945	75	2666	4.08 ug/L #	59
69) 1,3,5-Trichlorobenzene	13.116	180	12306	6.18 ug/L	96
70) 1,2,4-trichlorobenzene	13.591	180	10176m	6.11 ug/L	
71) Naphthalene	13.780	128	32139	/ 5.50 ug/L	99
72) 1,2,3-Trichlorobenzene	13.963	180	9887	6.10 ug/L	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

* 1 to 200 or 1 to

Instrument: MSVOA_X
ClientSampleId: VSTD005635

Manual IntegrationsAPPROVED