Data File : VX025229.D

Acq On : 18 Nov 2021 23:05

Operator : JC/MD Sample : M4677-08MSD

Misc : 5.0mL/MSVOA\_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Nov 19 05:31:58 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM111121WMA.M

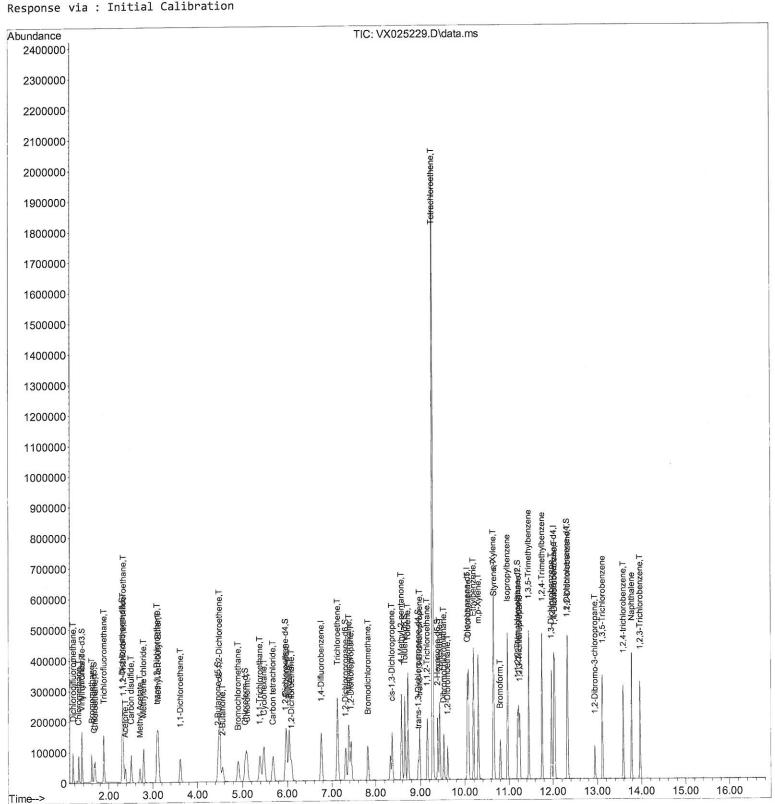
Quant Title : VOC Analysis

QLast Update : Fri Nov 19 05:25:45 2021

Instrument :
MSVOA\_X
ClientSampleId :
H0AB5MSD

# **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/19/2021 Supervised By :Mahesh Dadoda 11/22/2021



Data File : VX025229.D

Acq On : 18 Nov 2021 23:05

Operator : JC/MD

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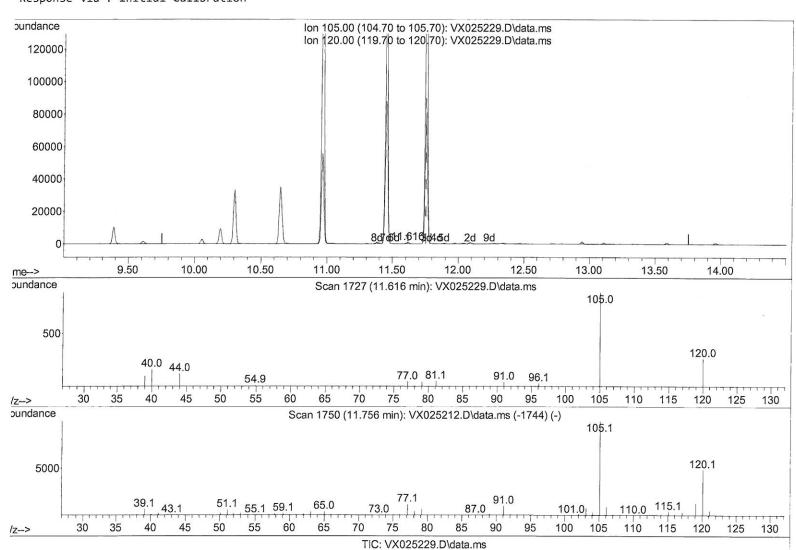
Quant Title : VOC Analysis

QLast Update : Fri Nov 19 05:25:45 2021 Response via : Initial Calibration



### **Manual Integrations APPROVED**

Reviewed By :John Carlone 11/19/2021 Supervised By :Mahesh Dadoda 11/22/2021



### (63) 1,2,4-Trimethylbenzene

11.616min (-0.140) 0.22 ug/L

response	1089	
Ion	Exp%	Act%
105.00	100.00	100.00
120.00	38.80	32.14
0.00	0.00	0.00
0.00	0.00	0.00

Data File: VX025229.D

Acq On : 18 Nov 2021 23:05

Operator : JC/MD Sample : M4677-08MSD

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Quant Time: Nov 19 05:31:58 2021

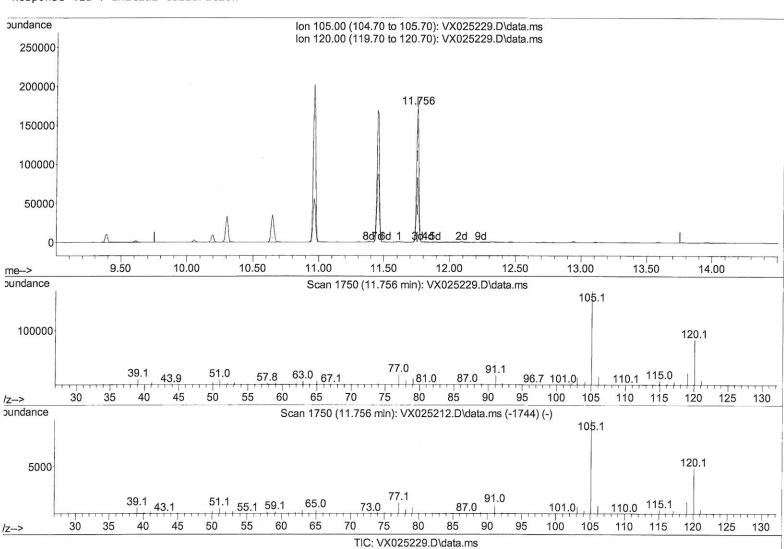
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Fri Nov 19 05:25:45 2021 Response via : Initial Calibration Instrument : MSVOA\_X ClientSampleId : H0AB5MSD

#### **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/19/2021 Supervised By :Mahesh Dadoda 11/22/2021



#### (63) 1,2,4-Trimethylbenzene

11.756min (+ 0.000) 43.88 ug/L m response 217779 Ion Exp% Act% 105.00 100.00 100.00 120.00 38.80 0.16# 0.00 0.00 0.00 0.00 0.00 0.00

Data File : VX025229.D

Acq On : 18 Nov 2021 23:05

Dperator : JC/MD
Sample : M4677-08MSD

fisc : 5.0mL/MSVOA\_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Nov 19 05:31:58 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM111121WMA.M

R.T. QIon Response Conc Units Dev(Min)

Quant Title : VOC Analysis

Compound

QLast Update : Fri Nov 19 05:25:45 2021 Response via : Initial Calibration Instrument: MSVOA\_X ClientSampleId: H0AB5MSD

# **Manual IntegrationsAPPROVED**

Reviewed By :John Carlone 11/19/2021 Supervised By :Mahesh Dadoda 11/22/2021

Internal Standards						
<ol> <li>1,4-Difluorobenzene</li> </ol>	6.763	114	172958	50.000	ug/L	0.00
28) Chlorobenzene-d5	10.055	117	160005	50.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	12.024	152	81543	50.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.368	65	44837	38.372	•	0.00
Spiked Amount 50.000	Range 60		Recove		76.740%	
<ol><li>7) Chloroethane-d5</li></ol>	1.660	69	37655	56.544	ug/L	0.00
Spiked Amount 50.000	Range 70	- 130	Recove	-	113.080%	
11) 1,1-Dichloroethene-d2	2.306	63	77822	38.724	ug/L	0.00
Spiked Amount 50.000	_	- 125	Recove		77.440%	
21) 2-Butanone-d5	4.459	46	77923	88.248	ug/L	0.00
Spiked Amount 100.000	Range 40	- 130	Recove	ry =	88.250%	
24) Chloroform-d	5.062	84	89685	43.630	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recove	ry =	87.260%	
26) 1,2-Dichloroethane-d4	5.958	65	53848	43.218	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recove	ry =	86.440%	
32) Benzene-d6	5.977	84	178069	40.776	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recove	ry =	81.560%	
36) 1,2-Dichloropropane-d6	7.312	67	54906	41.177	ug/L	0.00
Spiked Amount 50.000	Range 70	- 120	Recove	ry =	82.360%	
41) Toluene-d8	8.653	98	164664	39.484	ug/L	0.00
Spiked Amount 50.000	Range 80	- 120	Recove	ry =	78.960%	#
43) trans-1,3-Dichloroprop	. 8.952	79	28172	38.920	ug/L	0.00
Spiked Amount 50.000		- 125	Recove	ry =	77.840%	
47) 2-Hexanone-d5	9.384	63	61292	85.599	ug/L	0.00
Spiked Amount 100.000	Range 45	- 130	Recove		85.600%	
56) 1,1,2,2-Tetrachloroeth	. 11.195	84	82660	43.040	ug/L	0.00
		- 120	Recove		86.080%	
66) 1,2-Dichlorobenzene-d4	12.323	152	69974	43.288		0.00
	Range 80	- 120	Recove		86.580%	
*	<del>-</del>			153		
Target Compounds					Qva.	lue
<ol><li>Dichlorodifluoromethane</li></ol>	1.166	85	49967	36.965	1000	97
<ol><li>Chloromethane</li></ol>	1.294	50	50591	34.570		90
<ol><li>Vinyl chloride</li></ol>	1.374	62	60305	39.961		97
<ol><li>6) Bromomethane</li></ol>	1.599	94	36335	62.744	ug/L	99
<ol><li>8) Chloroethane</li></ol>	1.685	64	38896	51.326	ug/L	98
<ol><li>Trichlorofluoromethane</li></ol>	1.886	101	93237	42.327	ug/L	99
10) 1,1,2-Trichloro-1,2,2		101	47663	41.756		95
12) 1,1-Dichloroethene	2.319	96	45015	40.820	ug/L #	83
13) Acetone	2.380	43	46577	56.232	ug/L	97
14) Carbon disulfide	2.508	76	114207	33.613	ug/L	99
15) Methyl Acetate	2.709	43	53049	39.338	ug/L #	81
16) Methylene chloride	2.788	84	51874	42.524	ug/L	86
17) trans-1,2-Dichloroethene	3.093	96	50730	42.359		91
18) Methyl tert-butyl Ether	3.117	73	164166	43.855		90
19) 1,1-Dichloroethane	3.611	63	86061	42.481		93
20) cis-1,2-Dichloroethene	4.489	96	140531	106.053		98
22) 2-Butanone	4.562	43	81631	74.036	•	86
23) Bromochloromethane	4.904	128	30532	44.754		74
55.					示2	
1MYI M111121WMA M Eri Nov 19 06	-31-52 202	1				

Data File : VX025229.D

Acq On : 18 Nov 2021 23:05

Dperator : JC/MD
Sample : M4677-08MSD

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Quant Title : VOC Analysis

¿Last Update : Fri Nov 19 05:25:45 2021
Response via : Initial Calibration

Instrument: MSVOA\_X ClientSampleId: H0AB5MSD

### **Manual Integrations APPROVED**

Reviewed By :John Carlone 11/19/2021 Supervised By :Mahesh Dadoda 11/22/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)	
25) Chloroform	5.099	83	94383	45.455 ug/L 95	
27) 1,2-Dichloroethane	6.092	62	69106	45.392 ug/L # 89	
29) Cyclohexane	5.477	56	79084	38.687 ug/L 88	
30) 1,1,1-Trichloroethane	5.385	97	84133	42.790 ug/L # 94	
31) Carbon tetrachloride	5.678	117	75589	43.538 ug/L 99	
33) Benzene	6.044	78	206178	41.941 ug/L 100	
34) Trichloroethene	7.129	95	102368	79.688 ug/L 81	
35) Methylcyclohexane	7.385	83	88113	40.562 ug/L 94	
37) 1,2-Dichloropropane	7.434	63	51053	41.457 ug/L 100	
38) Bromodichloromethane	7.824	83	71786	43.076 ug/L 96	
39) cis-1,3-Dichloropropene	8.366	75	81184	40.149 ug/L 98	
40) 4-Methyl-2-pentanone	8.574	43	167368	85.516 ug/L # 84	
42) Toluene	8.720	91	227133	42.538 ug/L 96	
44) trans-1,3-Dichloropropene	8.982	75	79243	40.268 ug/L 97	
45) 1,1,2-Trichloroethane	9.153	97	55284	44.458 ug/L 98	
46) Tetrachloroethene	9.275	164	394772	368.332 ug/L 90	
48) 2-Hexanone	9.433	43	127131	79.247 ug/L # 85	
49) Dibromochloromethane	9.525	129	62952	44.348 ug/L 100	
50) 1,2-Dibromoethane	9.610	107	58692	44.102 ug/L # 96	
51) Chlorobenzene	10.079	112	160945	46.912 ug/L 97	
52) Ethylbenzene	10.195	91	245945	42.851 ug/L 93	
53) m,p-Xylene	10.305	106	99663	42.949 ug/L 80	
54) o-Xylene	10.646	106	99788	43.579 ug/L 79	
55) Styrene	10.659	104	166692	42.871 ug/L 81	
57) 1,1,2,2-Tetrachloroethane		83	85974	43.547 ug/L 97	
59) Bromoform	10.799	173	48342	43.867 ug/L # 95	
60) Isopropylbenzene	10.963	105	253773	43.523 ug/L 94	
61) 1,2,3-Trichloropropane	11.244	75	68719	44.755 ug/L 96	
62) 1,3,5-Trimethylbenzene	11.451	105	213020	43.060 ug/L 87	$\cap$
63) 1,2,4-Trimethylbenzene	11.756	105	217779m	43.878 ug/L 7/	70
64) 1,3-Dichlorobenzene	11.969	146	119641	45.012 ug/L 94 -	11/23/21
65) 1,4-Dichlorobenzene	12.043	146	118487	44.618 ug/L 93	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,
67) 1,2-Dichlorobenzene	12.335	146	119270	45.179 ug/L 94	
68) 1,2-Dibromo-3-chloropr	12.945	75	18954	42.681 ug/L # 62	
69) 1,3,5-Trichlorobenzene	13.116	180	86560	45.014 ug/L 97	
70) 1,2,4-trichlorobenzene	13.591		75782	45.179 ug/L 97	
71) Naphthalene	13.780		263310	46.292 ug/L 99	
72) 1,2,3-Trichlorobenzene	13.963	180	76395	45.941 ug/L 95	

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