Data File: VX025228.D

: 18 Nov 2021 22:41 Acq On

Operator : JC/MD Sample : M4677-07MS

: 5.0mL/MSVOA\_X/WATER Misc ALS Vial : 18 Sample Multiplier: 1

Quant Time: Nov 19 05:31:36 2021

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

Quant Title : VOC Analysis

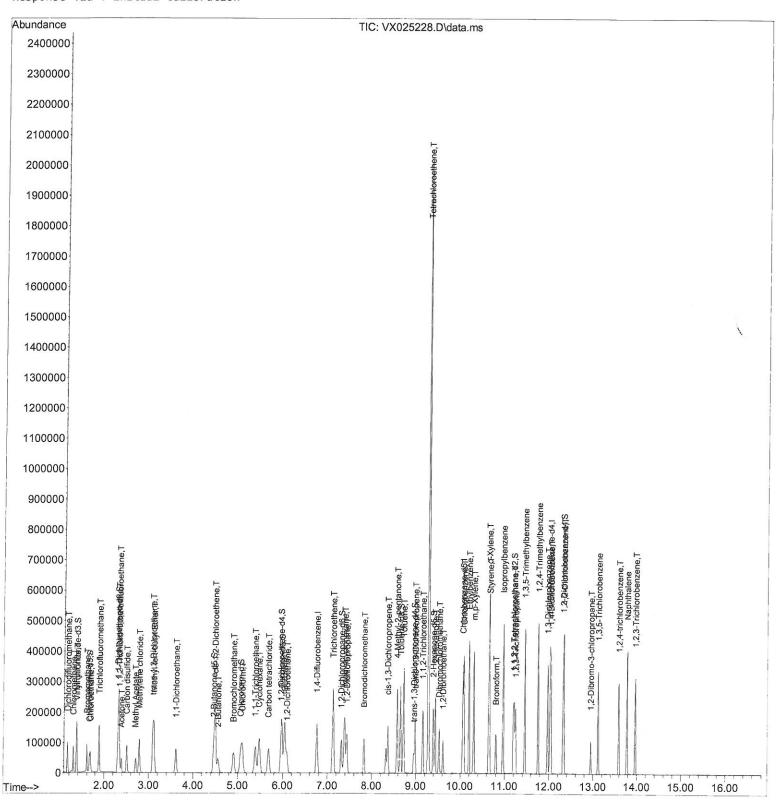
QLast Update : Fri Nov 19 05:25:45 2021

Response via: Initial Calibration



## **Manual IntegrationsAPPROVED**

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



Data File: VX025228.D

Acg On : 18 Nov 2021 22:41

: JC/MD Operator

: M4677-07MS

Sample Misc

: 5.0mL/MSVOA\_X/WATER ALS Vial : 18 Sample Multiplier: 1

Quant Time: Nov 19 05:31:36 2021

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

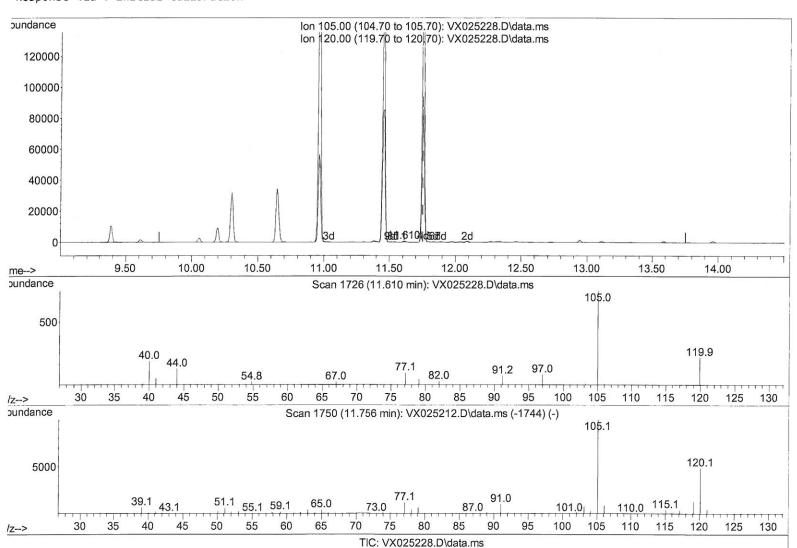
Quant Title : VOC Analysis

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



## **Manual Integrations APPROVED**

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



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

11.610min (-0.146) 0.20 ug/L

response	996			
Ion	Exp%	Act%		
105.00	100.00	100.00		
120.00	38.80	38.76		
0.00	0.00	0.00		
0.00	0.00	0.00		

Data File: VX025228.D

Acq On : 18 Nov 2021 22:41

Operator : JC/MD Sample : M4677-07MS

Misc : 5.0mL/MSVOA\_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Nov 19 05:31:36 2021

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

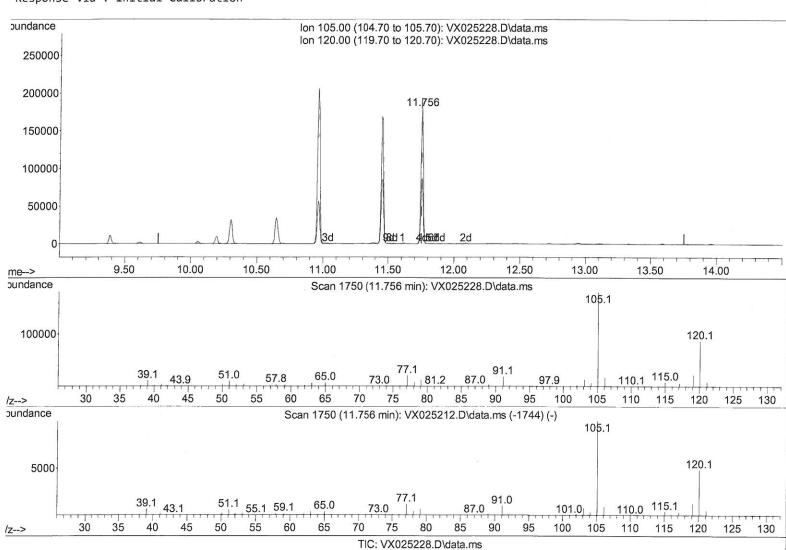
Quant Title : VOC Analysis

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

Instrument : MSVOA\_X ClientSampleId : H0AB5MS

### **Manual Integrations APPROVED**

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



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

11.	756min (	+ 0.000)	43.82 ug/L m	7 mg
res	ponse	217711		11/23/21
I	on	Exp%	Act%	
105	.00	100.00	100.00	
120	.00	38.80	0.18#	
0	.00	0.00	0.00	
0	.00	0.00	0.00	

Data File : VX025228.D

Acq On : 18 Nov 2021 22:41

Dperator : JC/MD
Sample : M4677-07MS

Misc : 5.0mL/MSVOA\_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Nov 19 05:31:36 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: H0AB5MS

# **Manual IntegrationsAPPROVED**

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

Compound		QIon	Response C	onc Ur	nits Dev(	Min)
Tatanana Ctandanda						
<pre>Internal Standards 1) 1,4-Difluorobenzene</pre>	6.763	114	175202	EQ 000	/1	0 00
28) Chlorobenzene-d5	10.055	117	161587	50.000 50.000		0.00 0.00
58) 1,4-Dichlorobenzene-d4	12.024	152	81616	50.000		0.00
50) 134 Dienzon Obenzene u4	12.024	132	01010	30.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.368	65	44812	37.860	ug/L	0.00
Spiked Amount 50.000	Range 60		Recovery		75.720%	
7) Chloroethane-d5	1.660	69		57.826		0.00
Spiked Amount 50.000	Range 70	- 130	Recovery		115.660%	
<pre>11) 1,1-Dichloroethene-d2</pre>	2.307	63	79553	39.078	ug/L	0.00
Spiked Amount 50.000	Range 60	- 125	Recovery	=	78.160%	
21) 2-Butanone-d5	4.459	46	79420	88.791	ug/L	0.00
Spiked Amount 100.000	Range 40	- 130	Recovery		88.790%	
24) Chloroform-d	5.062	84	90880	43.645	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recovery	=	87.280%	
26) 1,2-Dichloroethane-d4	5.958	65		43.982	Color State Colors	0.00
Spiked Amount 50.000	Range 70		Recovery		87.960%	
32) Benzene-d6	5.977	84		41.229	_	0.00
Spiked Amount 50.000		- 125	Recovery			
36) 1,2-Dichloropropane-d6	7.312	67	12 CO 10 CO 10	41.646	12 10 TO 100 TO	0.00
Spiked Amount 50.000	3. Table 1.	- 120	Recovery		83.300%	
41) Toluene-d8	8.653	98		40.361		0.00
Spiked Amount 50.000	•	- 120	Recovery	=	80.720%	0.00
43) trans-1,3-Dichloroprop.		79 125		39.224		0.00
Spiked Amount 50.000	Range 60 9.385	- 125 63	Recovery	=	78.440%	0.00
47) 2-Hexanone-d5 Spiked Amount 100.000		- 130		35.723	ug/L 85.720%	0.00
56) 1,1,2,2-Tetrachloroeth.	-	84	Recovery 84292	= 43.460		0.00
Spiked Amount 50.000	Range 65		Recovery		86.920%	0.00
66) 1,2-Dichlorobenzene-d4	12.323	152	-	13.946		0.00
Spiked Amount 50.000		- 120	Recovery	=	87.900%	0.00
			,		0, 1,200,0	
Target Compounds					Qva]	lue
2) Dichlorodifluoromethane	1.166	85	50731	37.049	-	98
<ol><li>Chloromethane</li></ol>	1.295	50	51205	34.542	ug/L	89
<ol><li>Vinyl chloride</li></ol>	1.374	62	61070	39.949	ug/L	98
<ol><li>Bromomethane</li></ol>	1.599	94	33115	6.451	ug/L	97
<ol><li>8) Chloroethane</li></ol>	1.685	64	38658	0.359	ug/L	97
<ol><li>Trichlorofluoromethane</li></ol>	1.886	101		12.402		100
10) 1,1,2-Trichloro-1,2,2		101		1.413		96
12) 1,1-Dichloroethene	2.319	96			ug/L #	83
13) Acetone	2.386	43		6.767		99
14) Carbon disulfide	2.508	76		32.791		99
15) Methyl Acetate	2.703	43			ug/L #	81
16) Methylene chloride	2.788	84		2.205		82
17) trans-1,2-Dichloroethene		96		2.154		88
18) Methyl tert-butyl Ether	3.117	73			ug/L #	90
19) 1,1-Dichloroethane	3.611	63		2.589		95
20) cis-1,2-Dichloroethene	4.489	96		4.425	-	100
<ul><li>22) 2-Butanone</li><li>23) Bromochloromethane</li></ul>	4.562 4.898	43 128		3.271		85 81
25) Di Omocnitoi omechane	4.070	140	JU1+1 4	.J • 013	ug/L #	OT

Data File : VX025228.D

Acq On : 18 Nov 2021 22:41

Dperator : JC/MD
Sample : M4677-07MS

disc : 5.0mL/MSVOA\_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Nov 19 05:31:36 2021

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

Quant Title : VOC Analysis

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

Instrument: MSVOA\_X ClientSampleId: H0AB5MS

# **Manual IntegrationsAPPROVED**

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	5.093	83	95434	45.372 ug/L	100
27) 1,2-Dichloroethane	6.092	62	70072	45.437 ug/L #	89
29) Cyclohexane	5.471	56	79451	38.486 ug/L	87
30) 1,1,1-Trichloroethane	5.385	97	85196	42.906 ug/L #	93
31) Carbon tetrachloride	5.678	117	75359	42.981 ug/L	98
33) Benzene	6.044	78	205279	41.349 ug/L	100
34) Trichloroethene	7.129	95	102825	79.260 ug/L	84
35) Methylcyclohexane	7.385	83	88206	40.207 ug/L	94
37) 1,2-Dichloropropane	7.434	63	51806	41.657 ug/L	100
38) Bromodichloromethane	7.824	83	71835	42.683 ug/L	98
39) cis-1,3-Dichloropropene	8.366	75	82715	40.506 ug/L	97
40) 4-Methyl-2-pentanone	8.574	43	165556	83.762 ug/L #	85
42) Toluene	8.720	91	227458	42.182 ug/L	96
44) trans-1,3-Dichloropropene	8.982	75	79372	39.938 ug/L	94
45) 1,1,2-Trichloroethane	9.153	97	55383	44.102 ug/L	98
46) Tetrachloroethene	9.275	164	393710	363.744 ug/L	90
48) 2-Hexanone	9.433	43	125981	77.762 ug/L #	86
49) Dibromochloromethane	9.525	129	64863	45.246 ug/L	99
50) 1,2-Dibromoethane	9.610	107	59204	44.051 ug/L #	96
51) Chlorobenzene	10.080	112	169516	48.926 ug/L	96
52) Ethylbenzene	10.195	91	248555	42.882 ug/L	93
53) m,p-Xylene	10.305	106	99100	42.288 ug/L	77
54) o-Xylene	10.646	106	98458	42.577 ug/L	80
55) Styrene	10.659	104	168456	42.900 ug/L	80
57) 1,1,2,2-Tetrachloroethane	11.213	83	86186	43.227 ug/L	97
59) Bromoform	10.805	173	48297	43.787 ug/L #	96
60) Isopropylbenzene	10.964	105	255508	43.781 ug/L	94
61) 1,2,3-Trichloropropane	11.244	75	67754	44.087 ug/L	97
62) 1,3,5-Trimethylbenzene	11.451	105	214357	43.291 ug/L	87 mg
63) 1,2,4-Trimethylbenzene	11.756	105	217711m	43.825 ug/L	87 mo 123 21 95 11 123 21
64) 1,3-Dichlorobenzene	11.969	146	118626	44.590 ug/L	
65) 1,4-Dichlorobenzene	12.043	146	118164	44.457 ug/L	92 ′
67) 1,2-Dichlorobenzene	12.335	146	119128	45.085 ug/L	93
68) 1,2-Dibromo-3-chloropr	12.945	75	19043	42.843 ug/L #	62
69) 1,3,5-Trichlorobenzene	13.116	180	82974	43.111 ug/L	96
70) 1,2,4-trichlorobenzene	13.591	180	73865	43.997 ug/L	98
71) Naphthalene	13.780		260589	45.773 ug/L	99
72) 1,2,3-Trichlorobenzene	13.963	180	75990	45.657 ug/L	96

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