Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110521\

Data File: VV023238.D

: 05 Nov 2021 17:34

Operator | : SY/MD Sample : M4535-04MS

Misc : 25.0mL/MSVOA V/WATER ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 09 03:24:34 2021

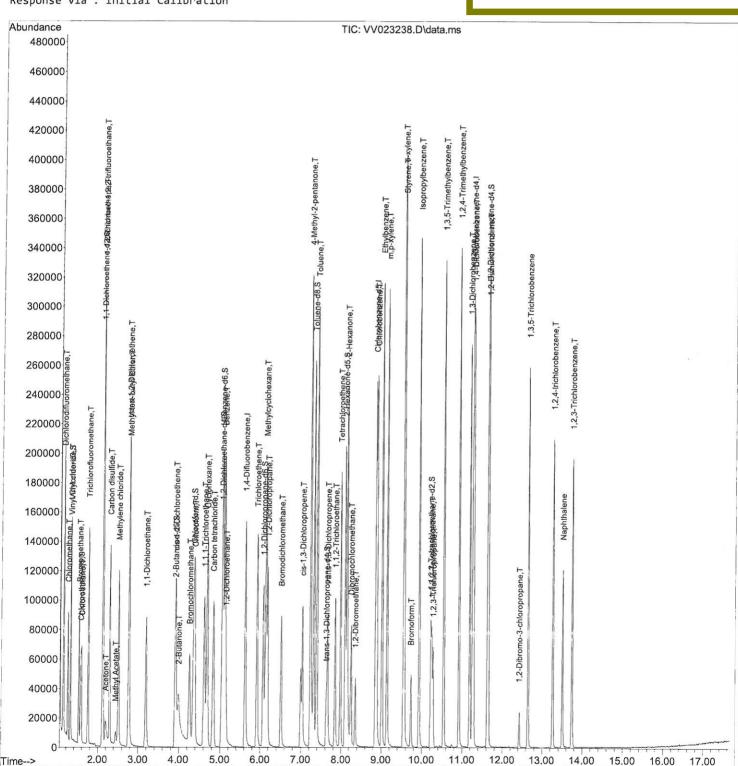
Quant Method: Z:\voasrv\HPCHEM1\MSVOA V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 09 02:04:24 2021 Response via: Initial Calibration

Instrument: MSVOA_V ClientSampleId:

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 11/10/2021 Supervised By : Mahesh Dadoda 11/10/2021



Time-->

Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110521\

Data File: VV023238.D

Acq On : 05 Nov 2021 17:34

Operator : SY/MD Sample : M4535-04MS

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

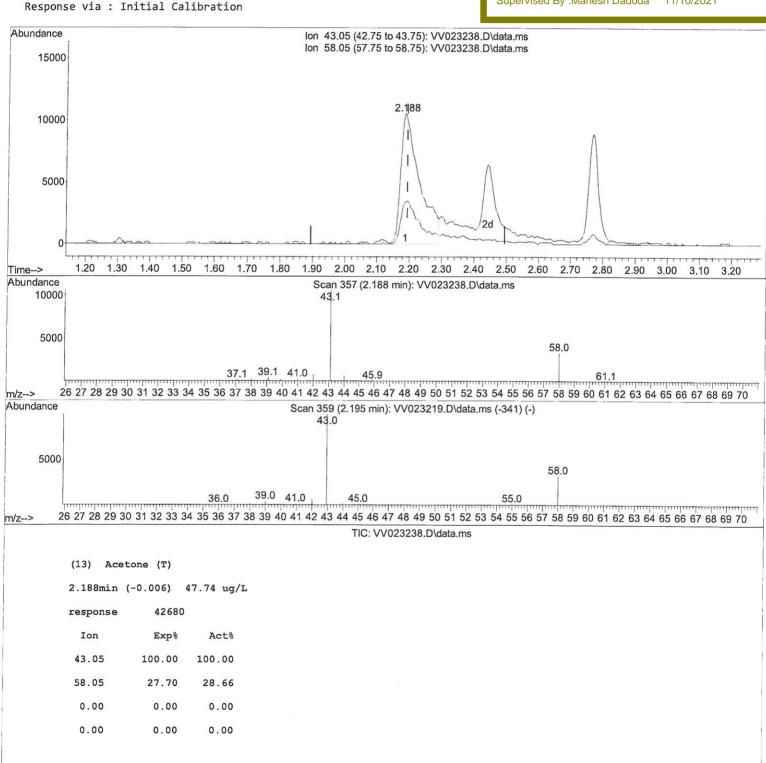
Quant Time: Nov 09 03:24:34 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 09 02:04:24 2021 Response via : Initial Calibration Instrument: MSVOA_V ClientSampleId: H4621MS

Manual Integrations APPROVED

Reviewed By :John Carlone 11/10/2021 Supervised By :Mahesh Dadoda 11/10/2021



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110521\

Data File: VV023238.D

: 05 Nov 2021 17:34 Acq On

: SY/MD Operator : M4535-04MS Sample

: 25.0mL/MSVOA_V/WATER Misc ALS Vial : 21 Sample Multiplier: 1

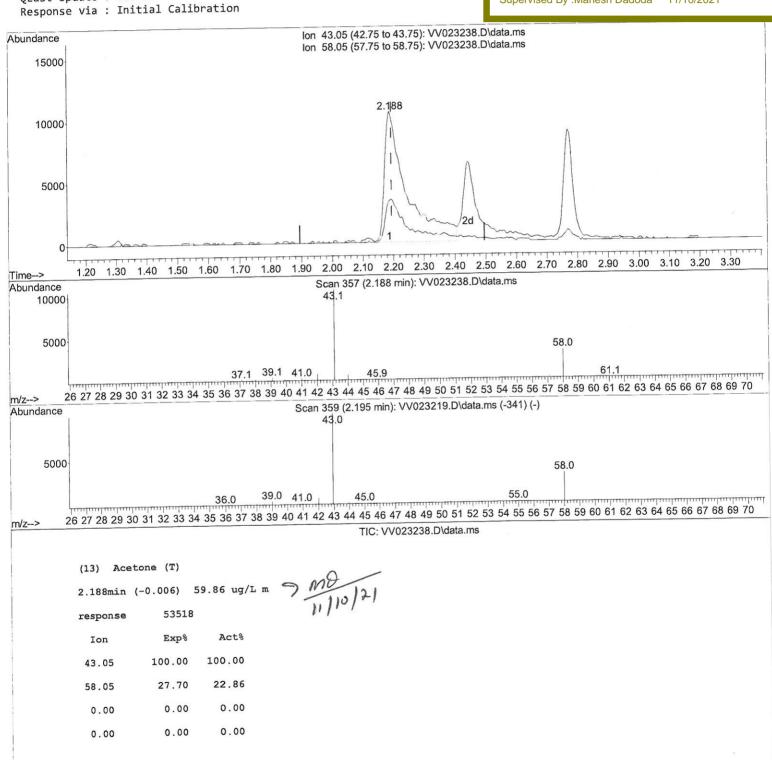
Quant Time: Nov 09 03:24:34 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_v\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 09 02:04:24 2021 Instrument: MSVOA_V ClientSampleId: 14621MS

Manual Integrations APPROVED

Reviewed By :John Carlone 11/10/2021 Supervised By :Mahesh Dadoda 11/10/2021



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV110521\

Data File : VV023238.D

Acq On : 05 Nov 2021 17:34

Operator : SY/MD Sample : M4535-04MS

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 09 03:24:34 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 09 02:04:24 2021 Response via : Initial Calibration Instrument: MSVOA_V ClientSampleld: H4621MS

Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc Units Dev(Min)	17
Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	135562	5.000 ug/L	0.00)
28) Chlorobenzene-d5	8.854		133745	5.000 ug/L	0.00	
58) 1,4-Dichlorobenzene-d4	11.249		73767	5.000 ug/L	0.00)
30) 1,4 Dienzor obenizent 1.						
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	41057	4.835 ug/L	0.00	
	Range 40	- 130	Recover			
7) Chloroethane-d5	1.568	69	34302	4.956 ug/L	0.00	
	Range 65		Recover			
<pre>11) 1,1-Dichloroethene-d2</pre>	2.108	63	78546	4.941 ug/L	0.00	
	Range 60		Recover	1.57	0 00	
20) 2-Butanone-d5	3.902	46		52.498 ug/L	0.00	
	Range 40		Recover	TOTAL CONTRACTOR OF SAME	0 00	
24) Chloroform-d	4.352		94995	5.249 ug/L	0.00	
T	Range 70		Recover	ry = 105.000% 5.362 ug/L	0.00	
26) 1,2-Dichloroethane-d4	5.037 Range 70		43639 Recover		0.00	
	5.050	84		5.211 ug/L	0.00	
<pre>32) Benzene-d6 Spiked Amount 5.000</pre>	Range 70		Recover		0.00	
36) 1,2-Dichloropropane-d6	6.072	67	53030	5.250 ug/L	0.00	
	Range 60		Recover		53,025	
41) Toluene-d8	7.317		171429	5.331 ug/L	0.00	
Spiked Amount 5.000	Range 70		Recover			
43) trans-1,3-Dichloroprop			19421	5.070 ug/L	0.00	
	Range 55		Recover	ry = 101.400%		
46) 2-Hexanone-d5	8.092	63	79448	56.373 ug/L	0.00	
Spiked Amount 50.000	Range 45	- 130	Recover	ry = 112.740%		
56) 1,1,2,2-Tetrachloroeth	. 10.217		39265	5.405 ug/L	0.00	
	Range 65	- 120	Recover			
00) 1,2 22011201 00011	11.625		63524	5.172 ug/L	0.00	
Spiked Amount 5.000	Range 80	- 120	Recover	ry = 103.400%		
Target Compounds				0va	lue	
2) Dichlorodifluoromethane	1.127	85	66727	5.048 ug/L	99	
3) Chloromethane	1.240	50	57358	5.104 ug/L	95	
5) Vinyl chloride	1.311	62	56727	5.054 ug/L	98	
6) Bromomethane	1.523	94	35852	4.997 ug/L	99	
8) Chloroethane	1.584	64	32874	5.075 ug/L	97	
Trichlorofluoromethane	1.754	101	84919	5.035 ug/L	99	
10) 1,1,2-Trichloro-1,2,2	. 2.118	101	41543	4.893 ug/L	99	9
12) 1,1-Dichloroethene	2.118	96	40457	5.005 ug/L	95	M/10/21
13) Acetone	2.188	43	53518m	59.864 ug/L	/	11/10/-1
14) Carbon disulfide	2.294	76	150556	4.935 ug/L	99	•
15) Methyl Acetate	2.442	43	14869	5.877 ug/L #	89	
16) Methylene chloride	2.507	84	50301	4.264 ug/L	95	
17) Methyl tert-butyl Ether	2.770	73	95835	5.385 ug/L	95	
18) trans-1,2-Dichloroethene		96	50237	5.055 ug/L	97	
19) 1,1-Dichloroethane	3.191	63	88763	5.290 ug/L	96 98	
21) 2-Butanone	3.986	43	71843	49.706 ug/L	90	
22) cis-1,2-Dichloroethene	3.915	96 128	50744 22415	5.306 ug/L # 5.083 ug/L	88	
23) Bromochloromethane	4.253	128	22413	J.005 ug/L	00	

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Quant Title : TRACE VOA SFAM1.0 QLast Update : Tue Nov 09 02:04:24 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleld : H4621MS

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 11/10/2021 Supervised By :Mahesh Dadoda 11/10/2021

певро	.,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,									
	Compound	R.T.	QIon	Response	Conc Unit	s Dev((Min)			
25)	Chloroform	4.378	83	94970	5.310 u	g/L	98			
	1,2-Dichloroethane	5.133	62	48747	5.124 u		98			
	1,1,1-Trichloroethane	4.609	97	82377	5.071 u	-	100			
	Cyclohexane	4.677	56	72343	4.970 u	g/L	99			
	Carbon tetrachloride	4.828	117	75345	5.165 u		98			
200000000000000000000000000000000000000	Benzene	5.101	78	199207	5.329 u	g/L	100			
34)	Trichloroethene	5.915	95	49837	5.013 u	g/L	99			
35)	Methylcyclohexane	6.133	83	74714	4.762 u	g/L	96			
	1,2-Dichloropropane	6.175	63	46584	5.338 u	g/L	99			
38)	Bromodichloromethane	6.513	83	60312	5.157 u	g/L	99			
39)	cis-1,3-Dichloropropene	7.031	75	57393	4.572 u	g/L	100			
40)	4-Methyl-2-pentanone	7.230	43	235246	58.121 u	g/L	98			
42)	Toluene	7.391	91	220485	5.514 u	g/L	98			
44)	trans-1,3-Dichloropropene	7.654	75	51958	4.989 u	g/L	99			
45)	1,1,2-Trichloroethane	7.841	97	32697	5.214 u	g/L	98			
47)	Tetrachloroethene	7.976	164	43350	5.032 u	g/L	98			
48)	2-Hexanone	8.143	43	170010	59.944 uį	g/L	98			
49)	Dibromochloromethane	8.246	129	41965	5.282 u	g/L	98			
50)	1,2-Dibromoethane	8.355	107	30631	5.271 u	g/L	97			
51)	Chlorobenzene	8.883	112	136831	5.149 u	g/L	98			
52)	Ethylbenzene	9.014	91	217551	5.159 ug	g/L	98			
53)	m,p-xylene	9.140	106	84874	5.128 ug	g/L	95			
54)	o-xylene	9.545	106	81678	5.261 u	g/L	97			
55)	Styrene	9.561	104	139345	5.239 ug	g/L	95			
57)	1,1,2,2-Tetrachloroethane	10.243	83	37264	5.424 ug	g/L	98			
59)	Bromoform	9.731	173	22641	5.139 ug		97			
	Isopropylbenzene	9.934	105	219341	5.182 u	g/L	99			
	1,2,3-Trichloropropane	10.275	75	26472	5.402 ug	<u> </u>	98			
	1,3,5-Trimethylbenzene	10.538	105	175836	5.010 ug		100			
63)	1,2,4-Trimethylbenzene	10.915	105	182725	5.230 ug		98			
64)	1,3-Dichlorobenzene	11.181	146	111255	5.144 uş	-	96			
	1,4-Dichlorobenzene	11.272	146	110945	5.023 ug	_	99			
	1,2-Dichlorobenzene	11.644	146	101702	5.255 ug	7.0	99			
68)	1,2-Dibromo-3-chloropr	12.429	75	5618	5.381 սք		93			
	1,3,5-Trichlorobenzene	12.644	180	81757	4.828 ug		99			
70)	1,2,4-trichlorobenzene	13.262	180	64242	4.737 ug		98			
200-200-200-200	Naphthalene	13.503	128	95680	4.785 ug		99			
72)	1,2,3-Trichlorobenzene	13.744	180	59324	5.000 ug	g/L	98			

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