Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File: VV022971.D

Acq On : 21 Oct 2021 22:01

Operator : SY/MD Sample : M4265-10MS

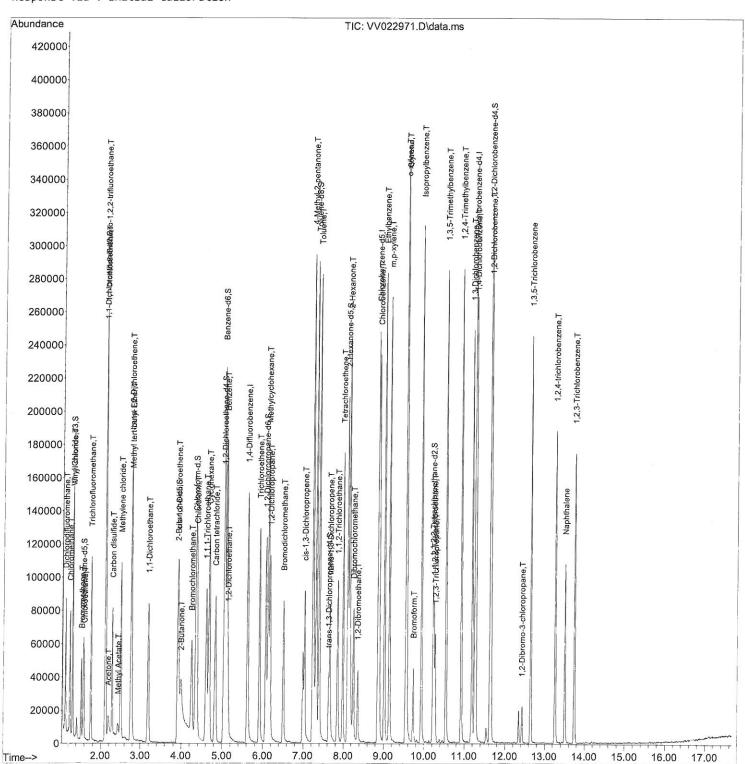
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Oct 22 04:58:17 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration Instrument:
MSVOA_V
ClientSampleId:
GB7J3MS

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File: VV022971.D

Acq On : 21 Oct 2021 22:01

Operator : SY/MD Sample : M4265-10MS

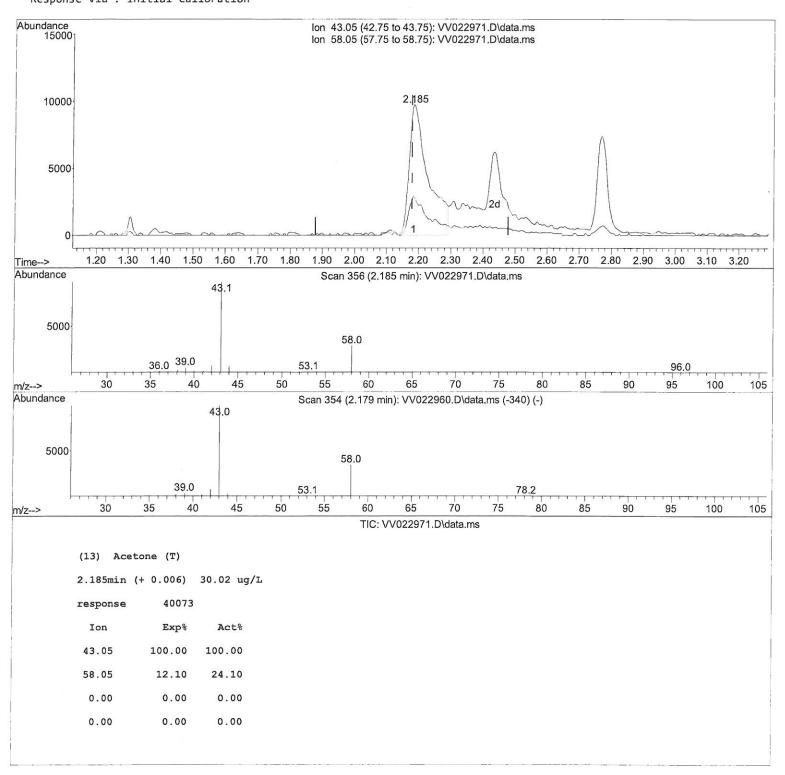
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Oct 22 04:58:17 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration Instrument: MSVOA_V ClientSampleld: GB7J3MS

Manual Integrations APPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File: VV022971.D

Acq On : 21 Oct 2021 22:01

Operator : SY/MD Sample : M4265-10MS

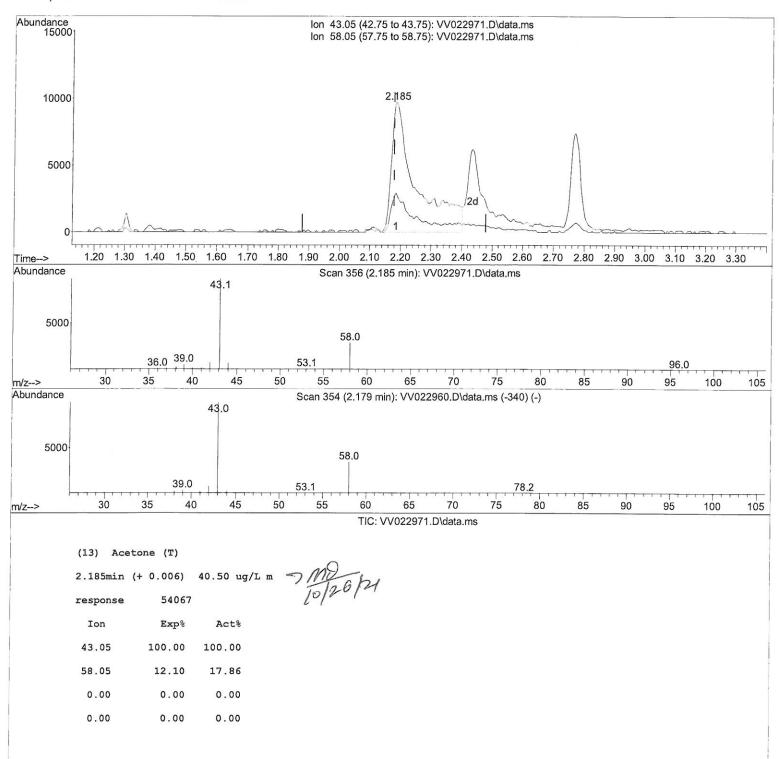
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Oct 22 04:58:17 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleld : GB7J3MS

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File : VV022971.D

Acq On : 21 Oct 2021 22:01 Operator : SY/MD

Sample

: M4265-10MS : 25.0mL/MSVOA_V/WATER Misc ALS Vial : 13 Sample Multiplier: 1

Quant Time: Oct 22 04:58:17 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration

Instrument : MSVOA_V ClientSampleId : GB7J3MS

Manual IntegrationsAPPROVED

Compound				Conc Units Dev		
Tutuural Ctandanda						
<pre>Internal Standards 1) 1,4-Difluorobenzene</pre>	5.619	114	132693	5.000 ug/L	0.00	a
28) Chlorobenzene-d5	8.853		134656	5.000 ug/L	0.00	
58) 1,4-Dichlorobenzene-d4	11.249		70262	5.000 ug/L	0.00	
remarks • Government of the control						
System Monitoring Compounds						
Vinyl Chloride-d3	1.304			4.738 ug/L	0.00	
Spiked Amount 5.000	Range 40			ry = 94.800%		
7) Chloroethane-d5	1.568	69		4.253 ug/L	0.00	
Spiked Amount 5.000 11) 1,1-Dichloroethene-d2	Range 65 2.108	63	Recover 67771			
Spiked Amount 5.000	Range 60		Recover		0.00	
20) 2-Butanone-d5	3.905	46		33.859 ug/L	0.02	
Spiked Amount 50.000	Range 40			ry = 67.720%		
24) Chloroform-d	4.349		104781	5.569 ug/L	0.00	
Spiked Amount 5.000	Range 70	- 125	Recover	ry = 111.400%	1	
26) 1,2-Dichloroethane-d4	5.034	65	45399	4.972 ug/L	0.00	
Spiked Amount 5.000	Range 70	- 130		ry = 99.400%		
32) Benzene-d6		84	200421	5.000 ug/L	0.00	
Spiked Amount 5.000	Range 70			ry = 100.000%		
36) 1,2-Dichloropropane-d6	6.072	67		5.180 ug/L	0.00	
Spiked Amount 5.000	Range 60			ry = 103.600%		
41) Toluene-d8	7.317	98	188573	5.442 ug/L	0.00	
Spiked Amount 5.000	Range 70		Recover			
43) trans-1,3-Dichloroprop Spiked Amount 5.000	Range 55		Recover	4.887 ug/L °V = 97.800%	0.00	
46) 2-Hexanone-d5	8.095	63		56.892 ug/L	0.00	
Spiked Amount 50.000	Range 45			y = 113.780%		
56) 1,1,2,2-Tetrachloroeth	. 10.217	84		5.563 ug/L	0.00	
	Range 65			y = 111.200%		
66) 1,2-Dichlorobenzene-d4				5.274 ug/L	0.00	
Spiked Amount 5.000	Range 80	- 120	Recover	y = 105.400%		
Target Compounds				Qva	lue	
Dichlorodifluoromethane	1.127	85	43607	4.509 ug/L	98	
Chloromethane	1.240	50	48269	4.966 ug/L	99	
Vinyl chloride	1.310	62	51367	5.115 ug/L	98	
Bromomethane	1.523	94	19525	3.119 ug/L	99	
8) Chloroethane	1.584	64	25318	4.075 ug/L	98	
Trichlorofluoromethane	1.754	101	64279	4.370 ug/L	97	
10) 1,1,2-Trichloro-1,2,2			33034	3.860 ug/L	95	
12) 1,1-Dichloroethene	2.117	96	30032	3.747 ug/L	93	MO 5/21
13) Acetone14) Carbon disulfide	2.185 2.291	43 76	54067m 83809	40.505 ug/L	100	10/26/21
15) Methyl Acetate	2.436	43	10933	3.720 ug/L 2.863 ug/L	94	<i>C</i> /
16) Methylene chloride	2.503	84	45121	4.090 ug/L	89	
17) Methyl tert-butyl Ether	2.770	73	85774	4.422 ug/L	96	
18) trans-1,2-Dichloroethene		96	45464	5.216 ug/L	97	
19) 1,1-Dichloroethane	3.188	63	85705	5.459 ug/L	97	
21) 2-Butanone	3.992	43	72497	33.171 ug/L	99	
22) cis-1,2-Dichloroethene	3.912	96	47009	5.136 ug/L	100	
23) Bromochloromethane	4.249	128	21386	5.259 ug/L	96	

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\

Data File : VV022971.D

Acq On : 21 Oct 2021 22:01 Operator : SY/MD

: M4265-10MS Sample

: 25.0mL/MSVOA_V/WATER Misc ALS Vial : 13 Sample Multiplier: 1

Quant Time: Oct 22 04:58:17 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Oct 22 04:55:17 2021 Response via : Initial Calibration

Instrument : MSVOA_V ClientSampleId : GB7J3MS

Manual IntegrationsAPPROVED

Compound	рт	QIon	Response	Conc Units Dev	(Min)				
	N.I.		Response	conc onits bev	(MTII)				
25) Chloroform	4.375	83	87771	4.823 ug/L	100				
27) 1,2-Dichloroethane	5.133	62	44665	4.917 ug/L	100				
29) 1,1,1-Trichloroethane	4.606	97	75618	4.918 ug/L	100				
30) Cyclohexane	4.674	56	62514	4.406 ug/L	99				
31) Carbon tetrachloride	4.828	117	67050	5.024 ug/L	98				
33) Benzene	5.101	78	184849	5.097 ug/L	100				
34) Trichloroethene	5.915	95	45474	4.611 ug/L	96				
35) Methylcyclohexane	6.130	83	65250	4.722 ug/L	99				
37) 1,2-Dichloropropane	6.175	63	42951	4.924 ug/L	98				
38) Bromodichloromethane	6.513	83	57122	5.315 ug/L	100				
39) cis-1,3-Dichloropropene	7.030	75	57789	5.005 ug/L	99				
40) 4-Methyl-2-pentanone	7.230	43	233061	49.729 ug/L	97				
42) Toluene	7.387	91	205859	5.591 ug/L	98				
44) trans-1,3-Dichloropropene	7.654	75	49720	5.282 ug/L	95				
45) 1,1,2-Trichloroethane	7.844	97	33199	5.129 ug/L	98				
47) Tetrachloroethene	7.976	164	39901	5.003 ug/L	98				
48) 2-Hexanone	8.143	43	174207	50.838 ug/L	97				
49) Dibromochloromethane	8.246	129	39857	5.260 ug/L	90				
50) 1,2-Dibromoethane	8.355	107	30099	4.980 ug/L	100				
51) Chlorobenzene	8.882	112	124501	5.147 ug/L	97				
52) Ethylbenzene	9.014	91	193502	5.177 ug/L	99				
53) m,p-xylene	9.140	106	77917	5.312 ug/L	98				
54) o-xylene	9.545	106	73484	5.325 ug/L	94				
55) Styrene	9.561	104	126930	5.338 ug/L	98				
57) 1,1,2,2-Tetrachloroethane	10.242	83	37494	5.726 ug/L	97				
59) Bromoform	9.734	173	21261	5.402 ug/L	98				
60) Isopropylbenzene	9.934	105	197119	5.523 ug/L	99				
61) 1,2,3-Trichloropropane	10.275	75	26656	5.460 ug/L	98				
62) 1,3,5-Trimethylbenzene	10.541	105	149417	5.182 ug/L	98				
63) 1,2,4-Trimethylbenzene	10.914	105	157223	5.448 ug/L	98				
64) 1,3-Dichlorobenzene	11.181	146	101938	5.442 ug/L	97				
65) 1,4-Dichlorobenzene	11.275	146	99408	5.227 ug/L	97				
67) 1,2-Dichlorobenzene	11.644	146	93867	5.342 ug/L	99				
68) 1,2-Dibromo-3-chloropr	12.429	75	5327	5.995 ug/L	86				
69) 1,3,5-Trichlorobenzene	12.644	180	77802	5.514 ug/L	97				
70) 1,2,4-trichlorobenzene	13.262	180	57326	5.318 ug/L	98				
71) Naphthalene	13.503	128	86402	5.043 ug/L	98				
72) 1,2,3-Trichlorobenzene	13.744	180	53600	5.274 ug/L	100				

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