Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File: VV023554.D

Acq On : 16 Nov 2021 23:33

Operator : SY/MD Sample : M4627-04MSD

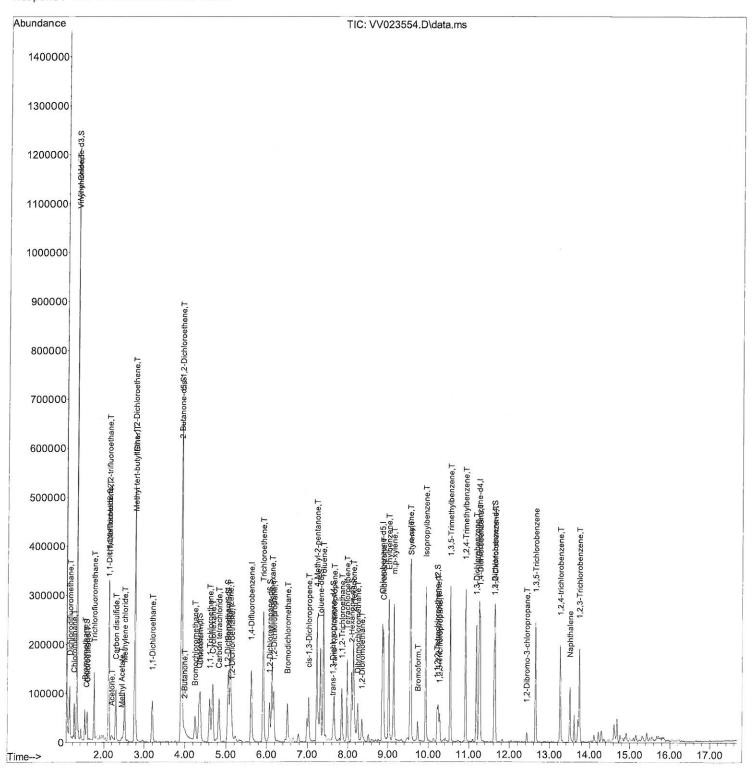
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Nov 17 03:41:01 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 02:49:39 2021 Response via : Initial Calibration Instrument:
MSVOA\_V
ClientSampleId:
H4637MSD

## **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File: VV023554.D

Acq On : 16 Nov 2021 23:33

Operator : SY/MD Sample : M4627-04MSD

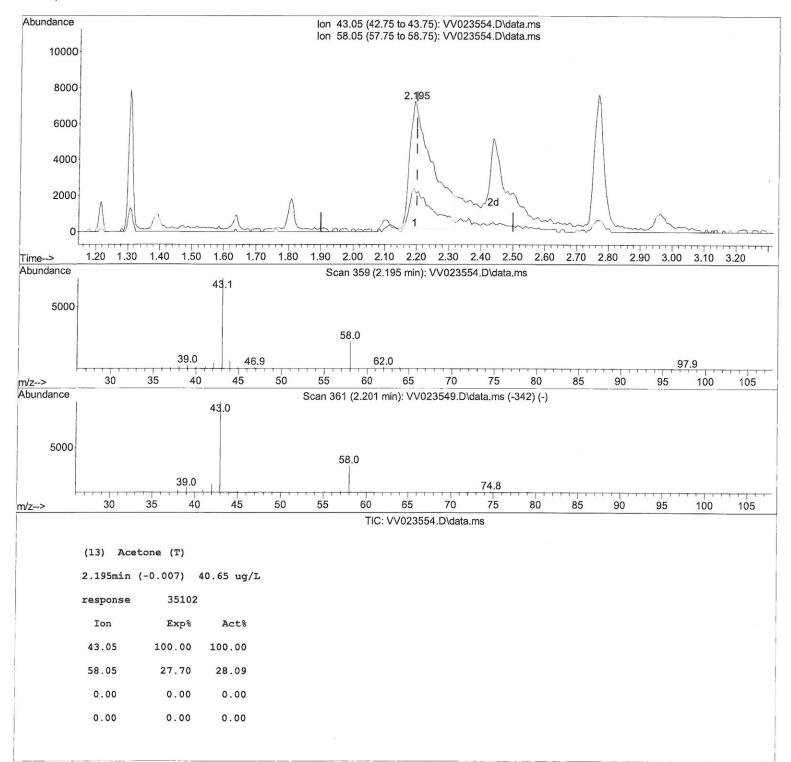
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Nov 17 03:41:01 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 02:49:39 2021 Response via : Initial Calibration Instrument : MSVOA\_V ClientSampleId : H4637MSD

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Data File : VV023554.D

Acq On : 16 Nov 2021 23:33

Operator : SY/MD Sample : M4627-04MSD

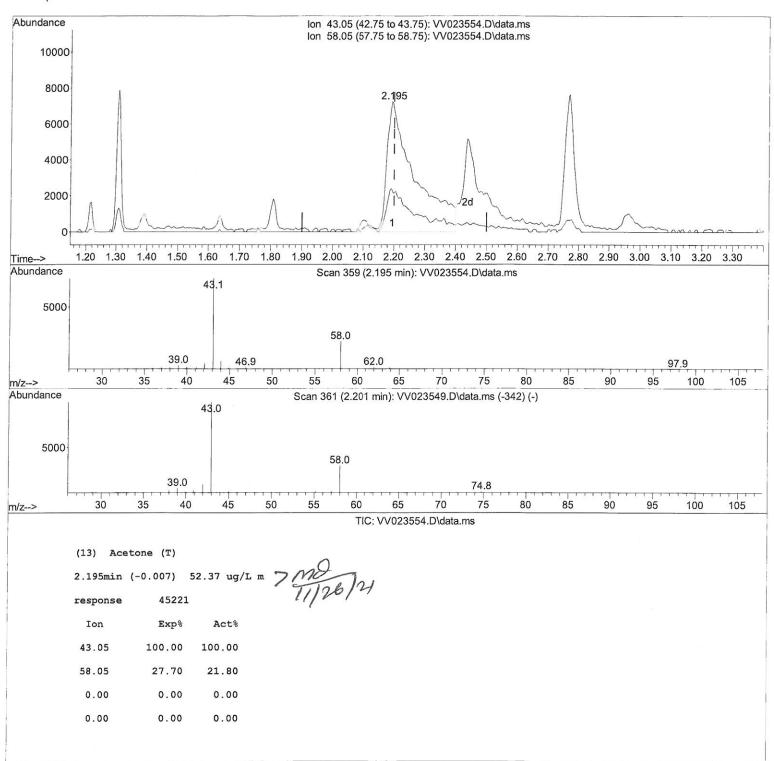
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Nov 17 03:41:01 2021

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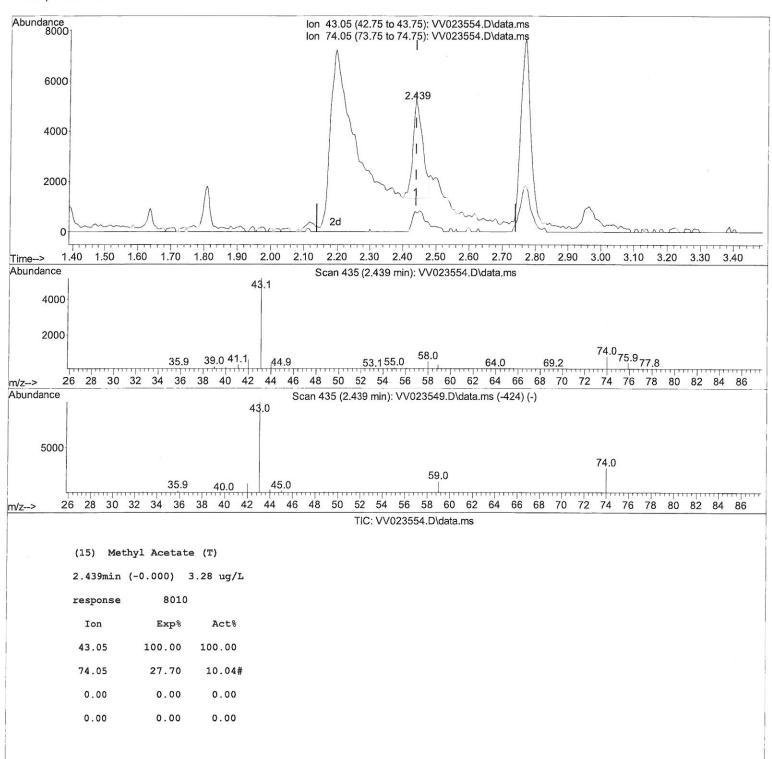
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Operator : SY/MD Sample : M4627-04MSD

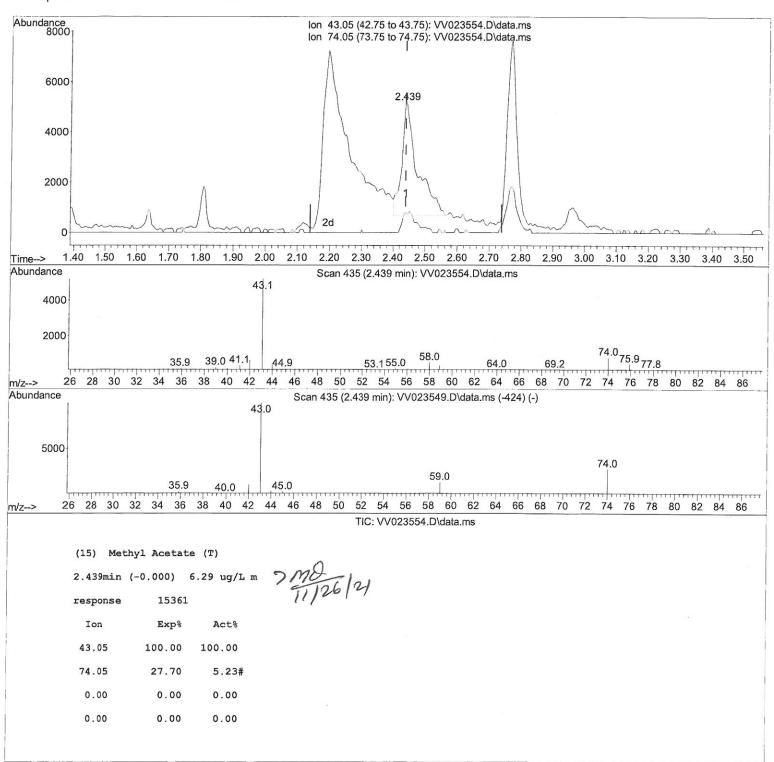
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 $\label{thm:local_var_def} Quant \ \mbox{Method} : \ \mbox{Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR110421WMA.M}$ 

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 02:49:39 2021 Response via : Initial Calibration Instrument : MSVOA\_V ClientSampleld : H4637MSD

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Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

Data File : VV023554.D

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Operator : SY/MD Sample : M4627-04MSD

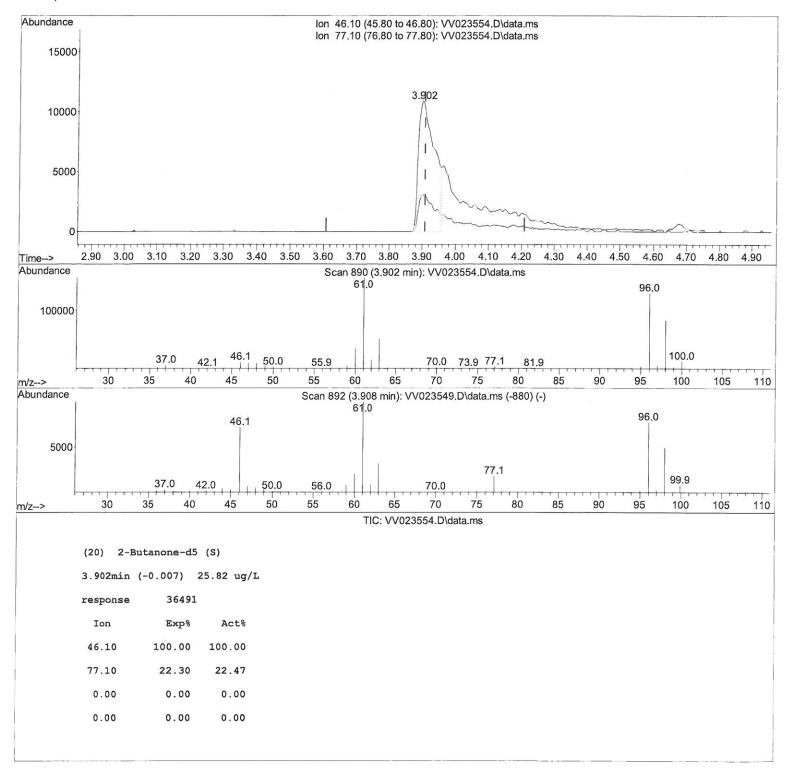
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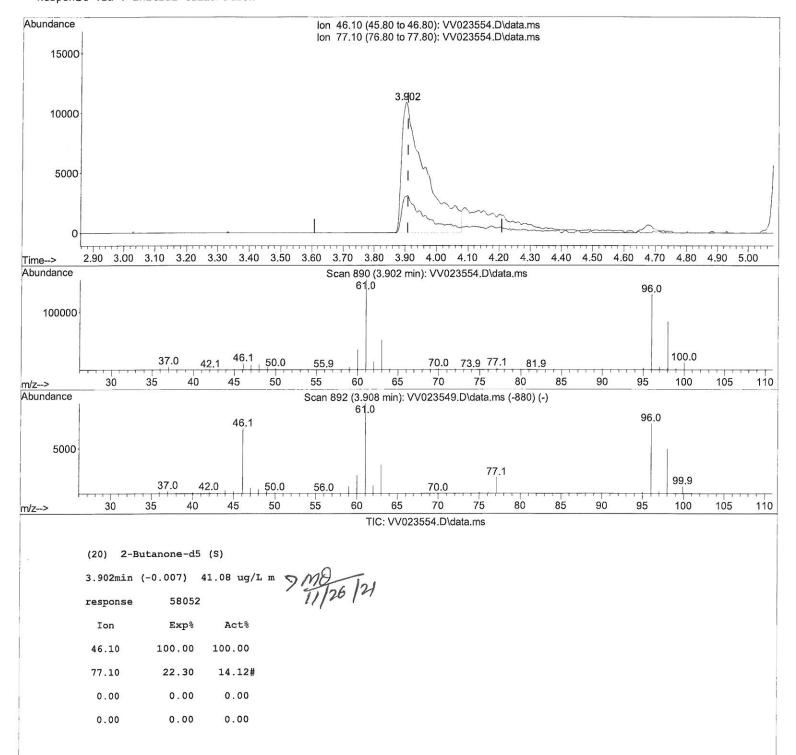
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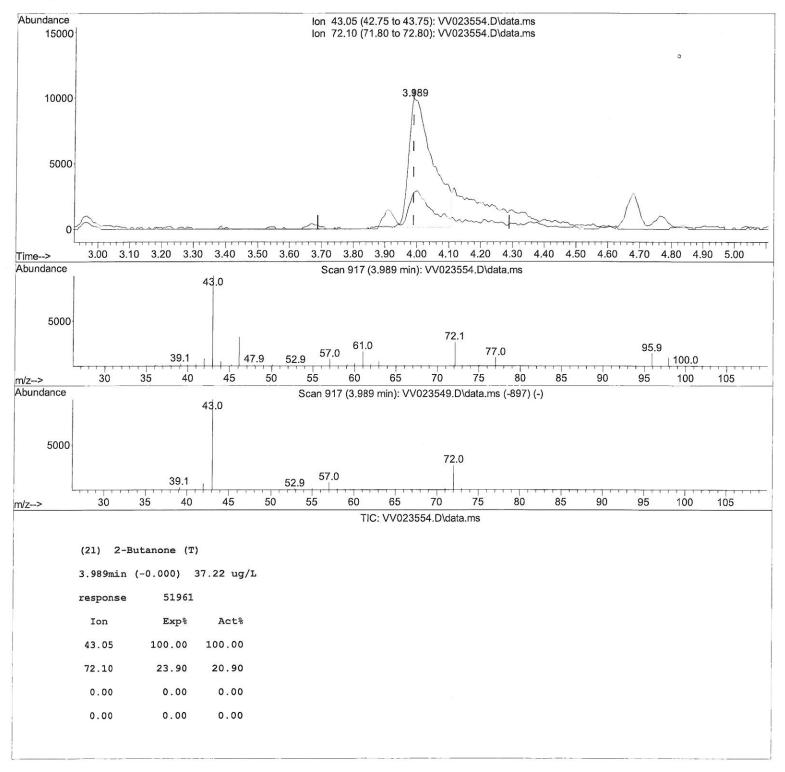
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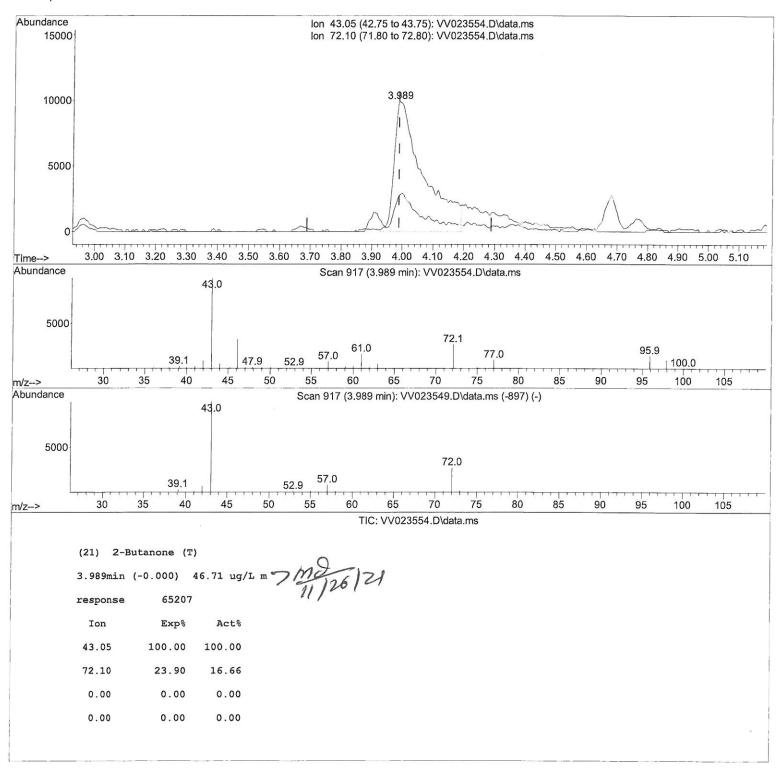
Misc : 25.0mL/MSVOA\_V/WATER
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Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 02:49:39 2021 Response via: Initial Calibration

Instrument : MSVOA\_V ClientSampleId : H4637MSD

# **Manual IntegrationsAPPROVED**

Compound	R.T. QI	on Response	Conc Uni	its Dev(	Min)
Internal Standards					
1) 1,4-Difluorobenzene	5.619 1	.14 130940	5.000	110/1	0.00
28) Chlorobenzene-d5	8.853 1		5.000		0.00
58) 1,4-Dichlorobenzene-d4		.52 69390	5.000	7000	0.00
50) 134 DICHIO ODCHZENE U4	11.2-10 1	.52 05550	31000	48/ -	5.55
System Monitoring Compounds					
4) Vinyl Chloride-d3	1.307	65 33641	4.101	ug/L	0.00
Spiked Amount 5.000	Range 40 -			82.000%	
<ol> <li>7) Chloroethane-d5</li> </ol>	1.568	69 25841	3.865	ug/L	0.00
Spiked Amount 5.000	Range 65 -	130 Recovery	y =	77.400%	
11) 1,1-Dichloroethene-d2	2.111	63 68638	4.470	ug/L	0.00
Spiked Amount 5.000	Range 60 -	125 Recovery	y =	89.400%	- not -
20) 2-Butanone-d5	3.902	46 58052m	41.078	ug/L	0.00 11/26/21
Spiked Amount 50.000	Range 40 -	(5		82.160%	-11/201
24) Chloroform-d		84 68836	3.938	ug/L	0.00
Spiked Amount 5.000	Range 70 -			78.800%	
26) 1,2-Dichloroethane-d4	5.034	65 32255			0.00
Spiked Amount 5.000	Range 70 -		63	82.000%	WWW. Marriado
32) Benzene-d6		84 128769		100	0.00
Spiked Amount 5.000	Range 70 -		A STATE SALES AND A STATE OF	77.800%	
36) 1,2-Dichloropropane-d6		67 37953	3.900		0.00
Spiked Amount 5.000	Range 60 -			78.000%	0.00
41) Toluene-d8		98 122465			0.00
Spiked Amount 5.000	Range 70 - 1			79.000%	0.00
43) trans-1,3-Dichloroprop.		79 15034	4.074		0.00
Spiked Amount 5.000	Range 55 - 1		2	81.400%	0.00
46) 2-Hexanone-d5		63 60197			0.00
Spiked Amount 50.000	Range 45 - 1	130 Recovery 84 30705	/ = 4.387	88.660%	0.00
56) 1,1,2,2-Tetrachloroeth Spiked Amount 5.000	Range 65 - 3			87.800%	0.00
66) 1,2-Dichlorobenzene-d4			4.241		0.00
Spiked Amount 5.000	Range 80 - 3			84.800%	0.00
Spiked Amount 3:000	nunge oo .	120 ((0000)		011000%	
Target Compounds				Qval	.ue
2) Dichlorodifluoromethane	1.130	85 52981	4.150	88	98
3) Chloromethane	1.240	50 48608	4.478	ug/L	96
<li>5) Vinyl chloride</li>	1.310	62 682869	62.985	0.73	99
6) Bromomethane	1.523	94 24706	3.565	ug/L	97
8) Chloroethane	1.587	64 30721	4.910	ug/L	97
<ol><li>Trichlorofluoromethane</li></ol>	1.754 16	01 75400	4.629	ug/L	98
10) 1,1,2-Trichloro-1,2,2	. 2.117 10	01 38210	4.659	ug/L	97
12) 1,1-Dichloroethene	2.121	96 51306		ug/L #	78
13) Acetone		43 45221m	52.369	177	1
14) Carbon disulfide		76 119348	4.050	-	98
15) Methyl Acetate		43 15361m	6.285		) My 104
16) Methylene chloride		84 41268	3.622		96 / 11/26/9
17) Methyl tert-butyl Ether		73 83864	4.879		92
18) trans-1,2-Dichloroethene		96 173325	18.057		99
19) 1,1-Dichloroethane		63 83694 65307	5.164		98
21) 2-Butanone			46.707		80 J
22) cis-1,2-Dichloroethene			40.766	_	89
23) Bromochloromethane	4.249 12	28 19740	4.634	ug/L #	75

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV111621\

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Operator : SY/MD Sample : M4627-04MSD

Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Nov 17 03:41:01 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 02:49:39 2021 Response via : Initial Calibration Instrument: MSVOA\_V ClientSampleId: H4637MSD

# **Manual IntegrationsAPPROVED**

Compound	R.T.	QIon	Response	Conc Units Dev(	Min)				
25) Chloroform	4.378	83	80569	4.664 ug/L	97				
27) 1,2-Dichloroethane	5.133	62	42099	4.582 ug/L	97				
29) 1,1,1-Trichloroethane	4.609	97	73007	4.665 ug/L	99				
30) Cyclohexane	4.680	56	62540	4.460 ug/L	95				
31) Carbon tetrachloride	4.828	117	65887	4.688 ug/L	96				
33) Benzene	5.101	78	185311	5.145 ug/L	100				
34) Trichloroethene	5.915	95	91907	9.596 ug/L	98				
35) Methylcyclohexane	6.130	83	66689	4.411 ug/L	97				
37) 1,2-Dichloropropane	6.175	63	38478	4.576 ug/L	99				
38) Bromodichloromethane	6.509	83	52915	4.696 ug/L	97				
39) cis-1,3-Dichloropropene	7.030	75	53867	4.454 ug/L	95				
40) 4-Methyl-2-pentanone	7.230	43	206980	53.077 ug/L	97				
42) Toluene	7.387	91	198429	5.151 ug/L	96				
44) trans-1,3-Dichloropropene	7.654	75	46992	4.683 ug/L	99				
45) 1,1,2-Trichloroethane	7.841	97	27698	4.585 ug/L	98				
47) Tetrachloroethene	7.976	164	38987	4.697 ug/L	96				
48) 2-Hexanone	8.143	43	145850	53.376 ug/L	98				
49) Dibromochloromethane	8.246	129	37236	4.864 ug/L	96				
50) 1,2-Dibromoethane	8.352	107	26476	4.729 ug/L #	99				
51) Chlorobenzene	8.882	112	120007	4.687 ug/L	99				
52) Ethylbenzene	9.014	91	200531	4.936 ug/L	99				
53) m,p-xylene	9.140	106	78971	4.953 ug/L	98				
54) o-xylene	9.545	106	72892	4.873 ug/L	99				
55) Styrene	9.561	104	127027	4.957 ug/L	94				
57) 1,1,2,2-Tetrachloroethane	10.242	83	31422	4.747 ug/L #	96				
59) Bromoform	9.734	173	19190	4.630 ug/L	96				
60) Isopropylbenzene	9.931	105	199850	5.019 ug/L	99				
61) 1,2,3-Trichloropropane	10.275	75	23506	5.099 ug/L	98				
62) 1,3,5-Trimethylbenzene	10.538	105	164140	4.971 ug/L	99				
63) 1,2,4-Trimethylbenzene	10.914	105	166498	5.067 ug/L	99				
64) 1,3-Dichlorobenzene	11.181	146	98773	4.855 ug/L	99				
65) 1,4-Dichlorobenzene	11.271	146	97394	4.687 ug/L	99				
67) 1,2-Dichlorobenzene	11.644	146	89584	4.921 ug/L	99				
68) 1,2-Dibromo-3-chloropr	12.429	75	4512	4.595 ug/L	84				
69) 1,3,5-Trichlorobenzene	12.644	180	76039	4.773 ug/L	99				
70) 1,2,4-trichlorobenzene	13.262	180	60955	4.778 ug/L	98				
71) Naphthalene	13.503	128	88097	4.683 ug/L	98				
72) 1,2,3-Trichlorobenzene	13.744	180	57207	5.125 ug/L	97				

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed