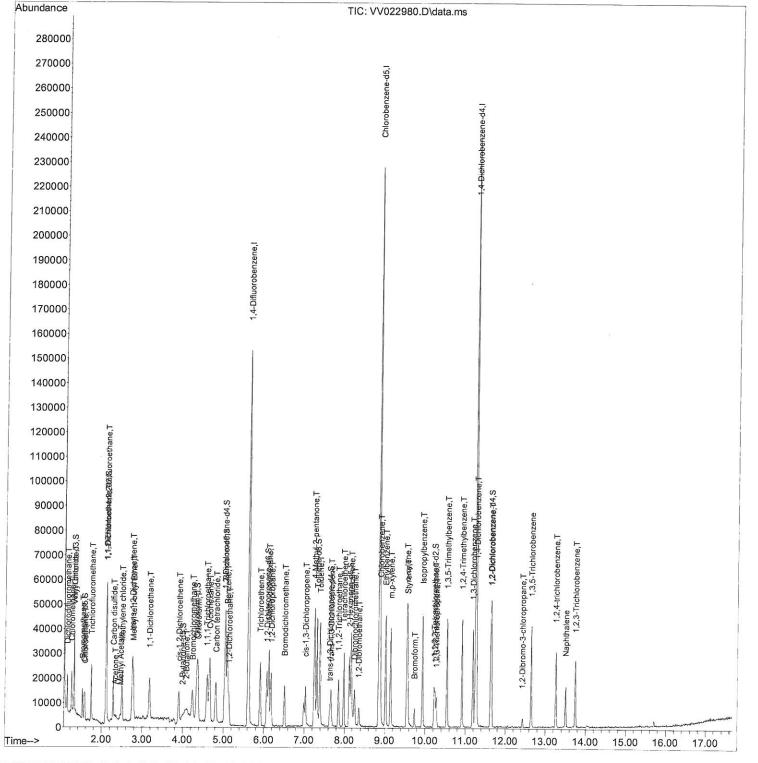
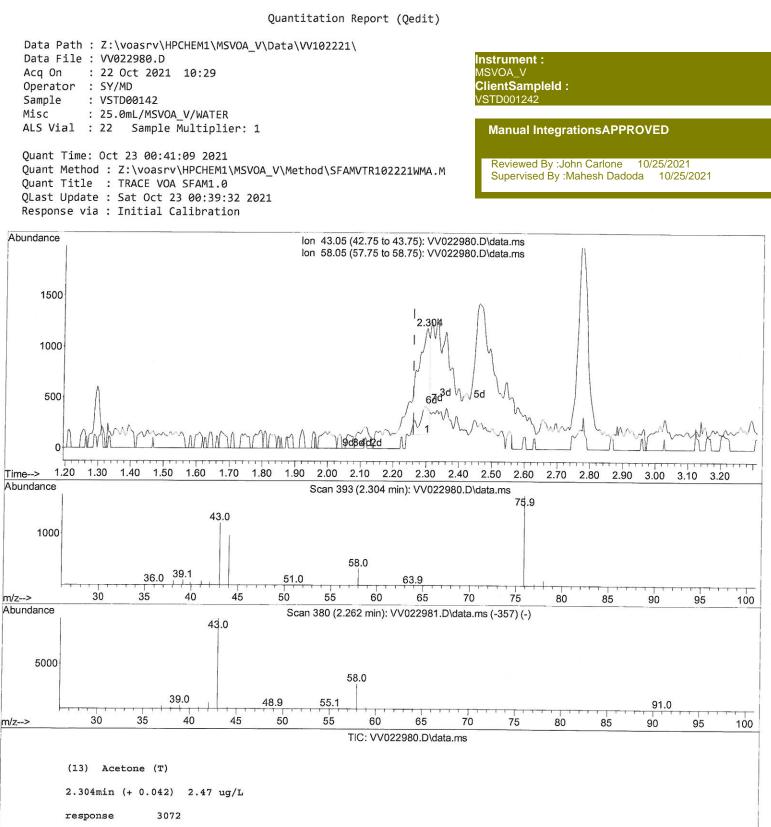
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102221\						
Data File : VV022980.D	Instrument :					
Acq On : 22 Oct 2021 10:29	MSVOA_V					
Operator : SY/MD	ClientSampleId :					
Sample : VSTD00142	VSTD001242					
Misc : 25.0mL/MSVOA_V/WATER						
ALS Vial : 22 Sample Multiplier: 1	Manual IntegrationsAPPROVED					
Quant Time: Oct 23 00:41:09 2021 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102221WMA.M Quant Title : TRACE VOA SFAM1.0 QLast Update : Sat Oct 23 00:39:32 2021	Reviewed By :John Carlone 10/25/2021 Supervised By :Mahesh Dadoda 10/25/2021					
Response via : Initial Calibration						



SFAMVTR102221WMA.M Sat Oct 23 01:05:09 2021



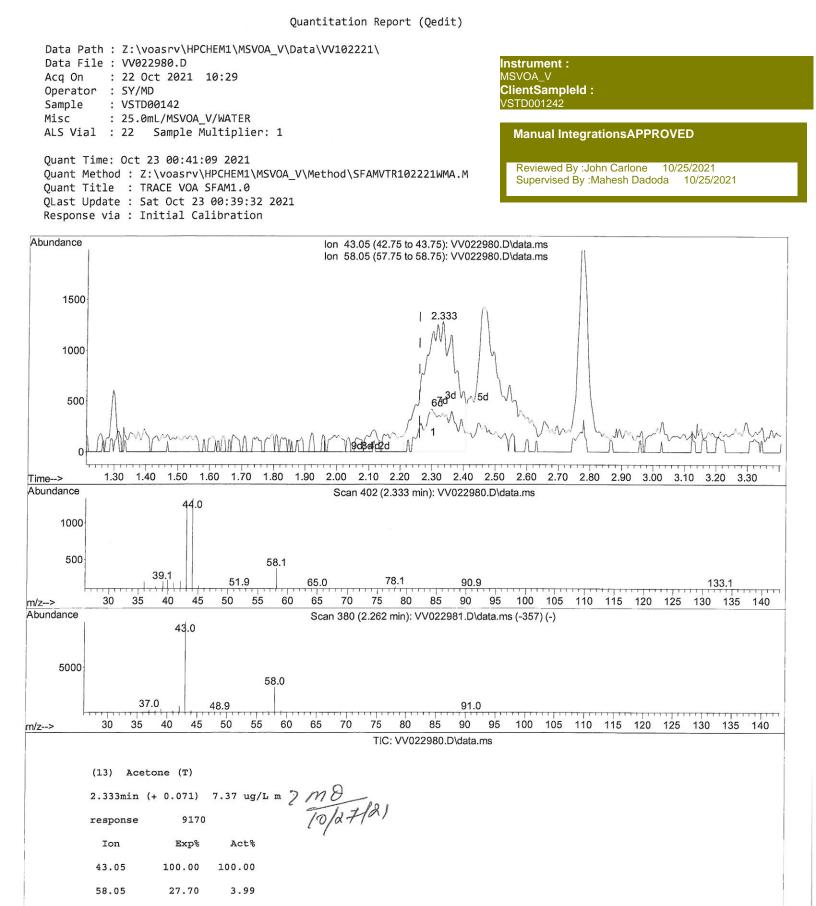
 Ion
 Exp%
 Act%

 43.05
 100.00
 100.00

 58.05
 27.70
 11.91

 0.00
 0.00
 0.00

 0.00
 0.00
 0.00



0.00

0.00

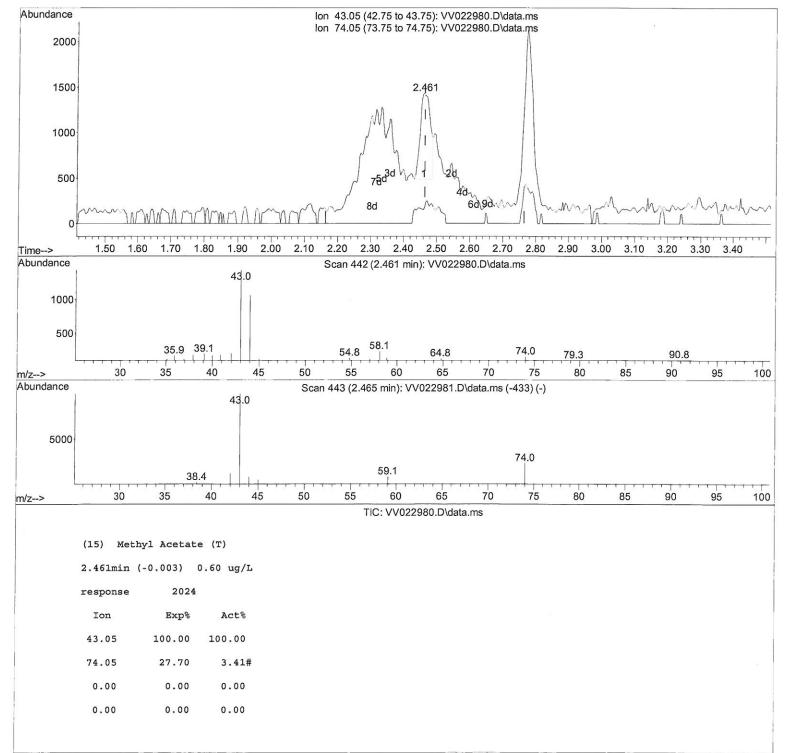
0.00

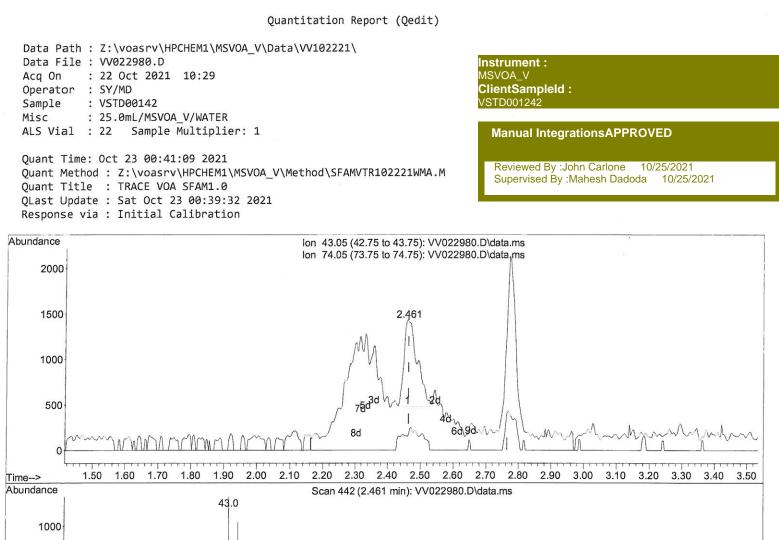
0.00

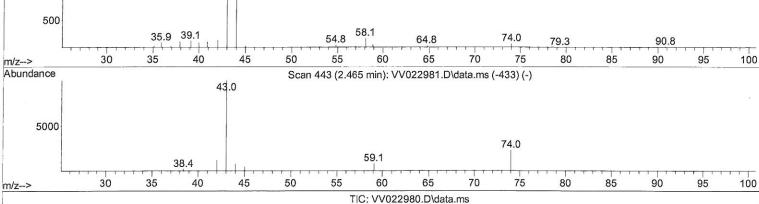
0.00

0.00









(15) Methyl Acetate (T)

2670

Exp%

100.00

27.70

0.00

0.00

0.79 ug/L m

Act%

2.58#

0.00

0.00

100.00

2.461min (-0.003)

response

Ion

43.05

74.05

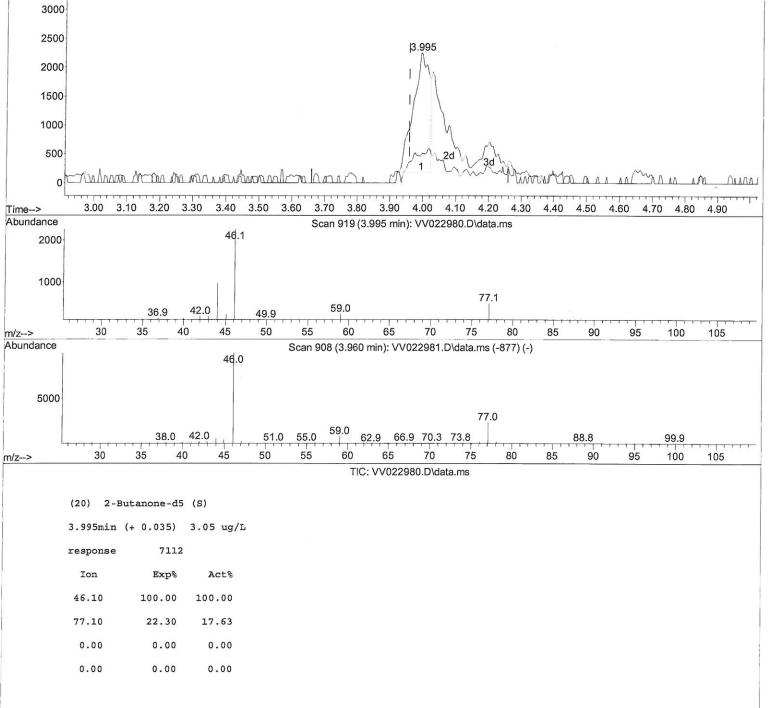
0.00

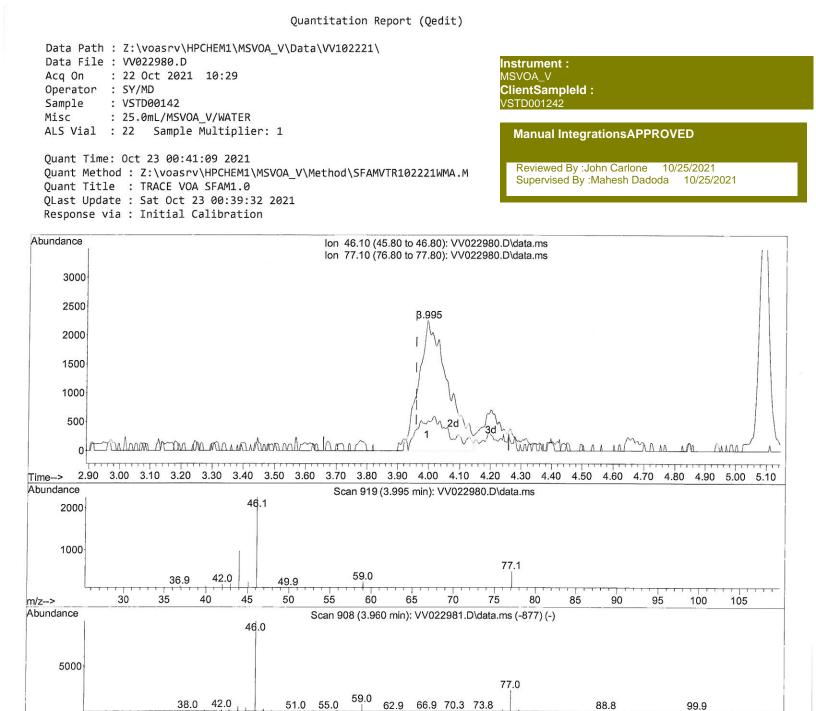
0.00

87/21

SFAMVTR102221WMA.M Sat Oct 23 01:03:32 2021







70

TIC: VV022980.D\data.ms

75

80

85

90

95

100

105

SFAMVTR102221WMA.M Sat Oct 23 01:03:57 2021

m/z-->

30

(20)

response

Ion

46.10

77.10

0.00

0.00

35

3.995min (+ 0.035)

2-Butanone-d5 (S)

14946

Exp%

100.00

22.30

0.00

0.00

40

45

6.42 ug/L m

Act%

8.39#

0.00

0.00

100.00

50

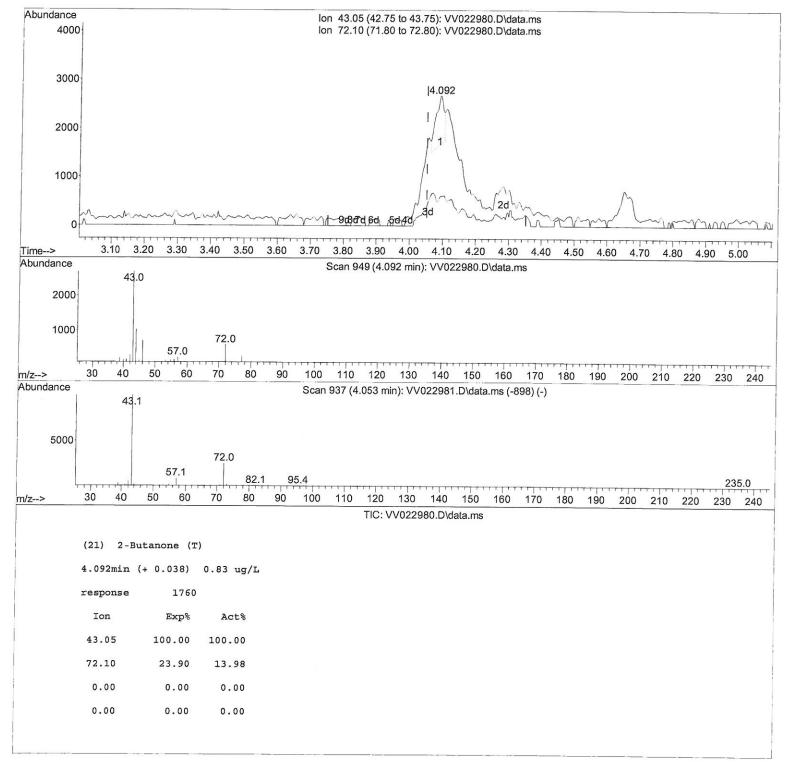
55

10/27/2

60

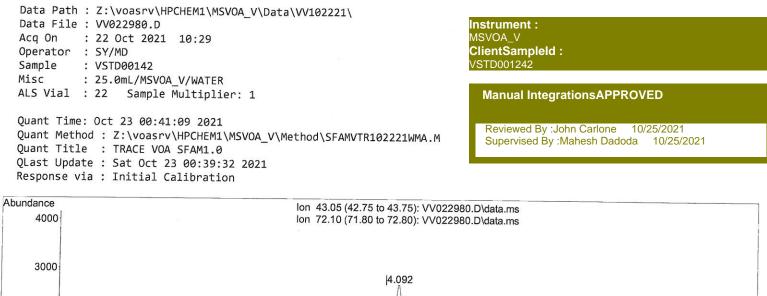
65

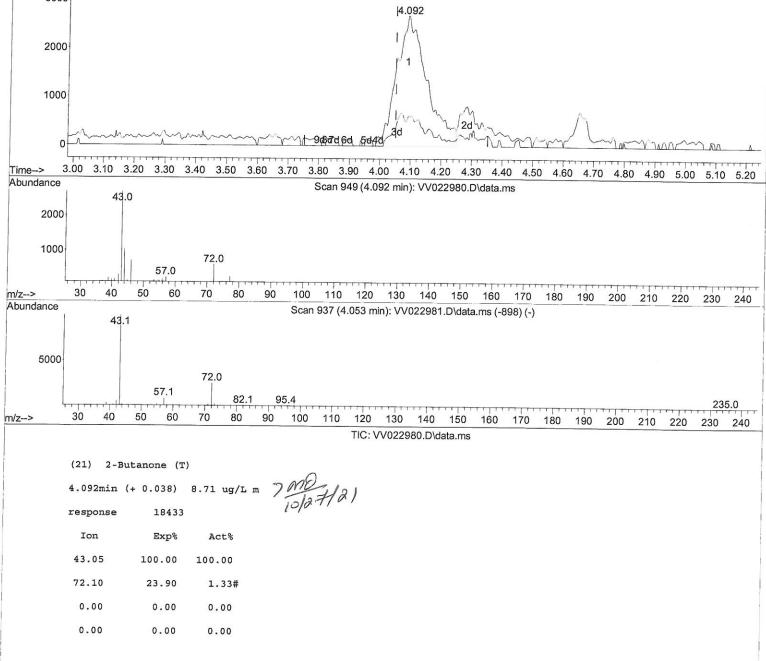




## SFAMVTR102221WMA.M Sat Oct 23 01:04:06 2021

Page: 1





Data Path : Z:\voasrv\HPCHEM1\M Data File : VV022980.D Acq On : 22 Oct 2021 10:29 Operator : SY/MD Sample : VSTD00142 Misc : 25.0mL/MSVOA_V/WATE ALS Vial : 22 Sample Multipl Quant Time: Oct 23 00:41:09 202	Instrument : MSVOA_V ClientSampleId : VSTD001242 Manual IntegrationsAPPROVED Reviewed By :John Carlone 10/25/2021				
Quant Method : Z:\voasrv\HPCHEM Quant Title : TRACE VOA SFAM1. QLast Update : Sat Oct 23 00:39 Response via : Initial Calibrat	Supervised By :Mahesh Dadoda 10/25/2021				
Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
1) 1,4-Difluorobenzene	5.612		130838	5.000 ug/L	0.00
28) Chlorobenzene-d5	8.853		126496	5.000 ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249	152	63052	5.000 ug/L	0.00
System Monitoring Compounds					
4) Vinyl Chloride-d3	1.297	65	10750	1.107 ug/L	0.00
7) Chloroethane-d5	1.558	69	6329	0.793 ug/L	-0.02
11) 1,1-Dichloroethene-d2	2.101	63	15496	0.878 ug/L	-0.02
20) 2-Butanone-d5	3.995	46	14946m	6.416 ug/L	0.02 0.04 MD 0.00 10/27/21 0.00 10/27/21
24) Chloroform-d	4.342	84	16763	0.891 ug/L	0.00 5/27/21
26) 1,2-Dichloroethane-d4	5.034	65	8854	0.975 ug/L	
32) Benzene-d6	5.040	84	35568	0.925 ug/L	0.00
36) 1,2-Dichloropropane-d6	6.069	67	10222	0.898 ug/L	-0.02
41) Toluene-d8	7.316	98	26805	0.801 ug/L	-0.01
<pre>43) trans-1,3-Dichloroprop 46) 2-Hexanone-d5</pre>	7.632 8.114	79 63	3561 10457	0.934 ug/L 7.015 ug/L	0.00 0.00
56) 1,1,2,2-Tetrachloroeth		84	7060	0.924 ug/L	0.00
66) 1,2-Dichlorobenzene-d4	11.625		10200	0.867 ug/L	0.00
, .				ġ.	
Target Compounds				•	alue
<ol><li>Dichlorodifluoromethane</li></ol>	1.127	85	8165	0.878 ug/L	99
3) Chloromethane	1.236	50	9573	1.014 ug/L	100
5) Vinyl chloride	1.301	62	9265	0.946 ug/L	98
<ul><li>6) Bromomethane</li><li>8) Chloroethane</li></ul>	1.513 1.577	94 64	4720 5055	0.805 ug/L 0.863 ug/L	99 99
<ol> <li>9) Trichlorofluoromethane</li> </ol>	1.744	101	12861	0.920 ug/L	97
10) 1,1,2-Trichloro-1,2,2	2.111	101	7067	0.861 ug/L	99
12) 1,1-Dichloroethene	2.111	96	6980	0.907 ug/L	94 0
13) Acetone	2.333	43	9170m	7.369 ug/L	
14) Carbon disulfide	2.285	76	18875	0.878 ug/L	98 CMD-
15) Methyl Acetate	2.461	43	2670m	0.786 ug/L	98 MU 91 (10/2-7/21
16) Methylene chloride	2.500	84	7696	0.752 ug/L	91 10 2-7121
17) Methyl tert-butyl Ether	2.776	73	17745	0.994 ug/L	97
18) trans-1,2-Dichloroethene	2.748	96	7934	0.980 ug/L	95
19) 1,1-Dichloroethane 21) 2-Butanone	3.182 4.092	63 43	15116 18433m	1.035 ug/L 8.713 ug/L	96
22) cis-1,2-Dichloroethene	3.905	96	7147	0.797 ug/L #	92
23) Bromochloromethane	4.239	128	3881	0.986 ug/L	86
25) Chloroform	4.374	83	17637	0.998 ug/L	91
27) 1,2-Dichloroethane	5.133	62	10317	1.160 ug/L	100
29) 1,1,1-Trichloroethane	4.593	97	14303	0.987 ug/L	98
30) Cyclohexane	4.657	56	13163	0.995 ug/L	98
31) Carbon tetrachloride	4.815	117	12082	0.964 ug/L	98
33) Benzene	5.091	78	36003	1.054 ug/L	100
34) Trichloroethene	5.908	95	8356	0.910 ug/L	98
35) Methylcyclohexane	6.117	83	11069	0.858 ug/L	99
37) 1,2-Dichloropropane	6.175	63	8591	1.046 ug/L #	91
38) Bromodichloromethane	6.509	83 75	10793	1.062 ug/L	99 100
<pre>39) cis-1,3-Dichloropropene 40) 4-Methyl-2-pentanone</pre>	7.030 7.252	75 43	9852 35576	0.917 ug/L 8.238 ug/L	98
42) Toluene	7.387	45 91	31288	0.900 ug/L	98
			22200		0.7

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV102221\ Data File : VV022980.D Acq On : 22 Oct 2021 10:29 Operator : SY/MD Sample : VSTD00142 Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 22 Sample Multiplier: 1

Compound

Instrument: MSVOA\_V ClientSampleId: VSTD001242

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 10/25/2021 Supervised By :Mahesh Dadoda 10/25/2021

Quant Time: Oct 23 00:41:09 2021 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR102221WMA.M Quant Title : TRACE VOA SFAM1.0 QLast Update : Sat Oct 23 00:39:32 2021 Response via : Initial Calibration

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

							and a second state
44)	trans-1,3-Dichloropropene	7.661	75	8224	0.931	ug/L	100
45)	1,1,2-Trichloroethane	7.847	97	6167	1.019		98
47)	Tetrachloroethene	7.972	164	6875	0.921	-	94
48)	2-Hexanone	8.162	43	25328	8.021		93
49)	Dibromochloromethane	8.249	129	7155	1.007		99
50)	1,2-Dibromoethane	8.355	107	5336	0.952		<sup>#</sup> 90
51)	Chlorobenzene	8.882	112	21938	0.968		98
52)	Ethylbenzene	9.014	91	30028	0.863 (		100
53)	m,p-xylene	9.140	106	11792	0.856 (		91
54)	o-xylene	9.545	106	11129	0.866 ι		98
55)	Styrene	9.564	104	18250	0.823 i		99
57)	1,1,2,2-Tetrachloroethane	10.246	83	7017	1.133 u		99
59)	Bromoform	9.734	173	3774	1.069 ı		97
60)	Isopropylbenzene	9.931	105	28860	0.902 L		99
61)	1,2,3-Trichloropropane	10.275	75	4666	1.073 u		97
62)	1,3,5-Trimethylbenzene	10.538	105	23377	0.913 u	Jg/L	95
63)	1,2,4-Trimethylbenzene	10.914	105	22252	0.867 u		98
64)	1,3-Dichlorobenzene	11.181	146	16223	0.967 L		89
65)	1,4-Dichlorobenzene	11.275	146	17712	1.043 u		100
67)	1,2-Dichlorobenzene	11.644	146	15487	0.987 L	lg/L	97
68)	1,2-Dibromo-3-chloropr	12.429	75	921	1.132 u	lg/L	95
69)	1,3,5-Trichlorobenzene	12.647	180	12992	1.028 u		99
70)	1,2,4-trichlorobenzene	13.262	180	9502	0.994 u	ig/L	95
71)	Naphthalene	13.506	128	12626	0.831 u	ig/L	100
72)	1,2,3-Trichlorobenzene	13.747	180	8432	0.937 u		96

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