Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV111821\

Data File : VV023630.D

: 19 Nov 2021 01:38 Acq On

Operator : SY/MD : VSTDCCC005EC Sample

Misc : 25.0mL/MSVOA V/WATER ALS Vial : 39 Sample Multiplier: 1

Quant Time: Nov 19 04:12:08 2021

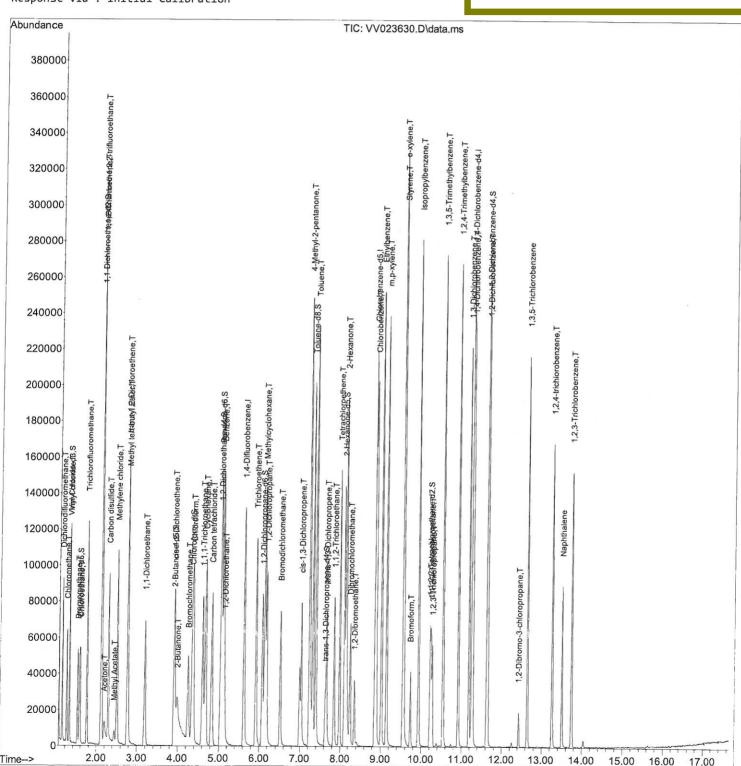
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Nov 19 03:51:44 2021 Response via: Initial Calibration

Instrument: MSVOA_V **LabSampleld**:

Manual IntegrationsAPPROVED

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



17.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111821\

Data File: VV023630.D

Acq On : 19 Nov 2021 01:38

Operator : SY/MD Sample : VSTDCCC005EC

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 39 Sample Multiplier: 1

Quant Time: Nov 19 04:12:08 2021

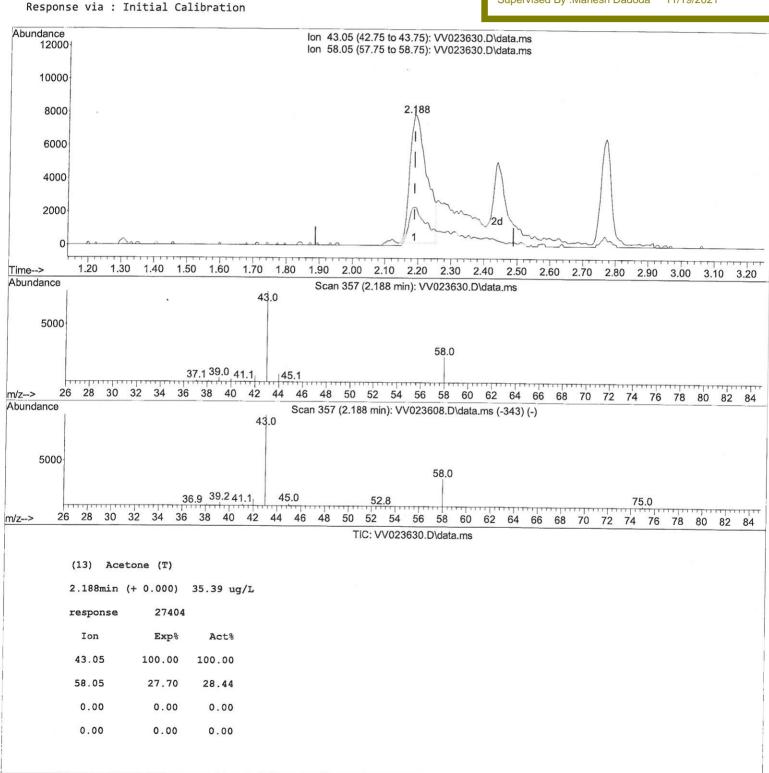
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Quant Title : TRACE VOA SFAM1.0

QLast Update : Fri Nov 19 03:51:44 2021

Instrument: MSVOA_V LabSampleId: VSTDCCC005EC

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111821\

Data File: VV023630.D

: 19 Nov 2021 01:38 Acq On

: SY/MD Operator Sample : VSTDCCC005EC

: 25.0mL/MSVOA V/WATER Misc Sample Multiplier: 1 ALS Vial : 39

Quant Time: Nov 19 04:12:08 2021

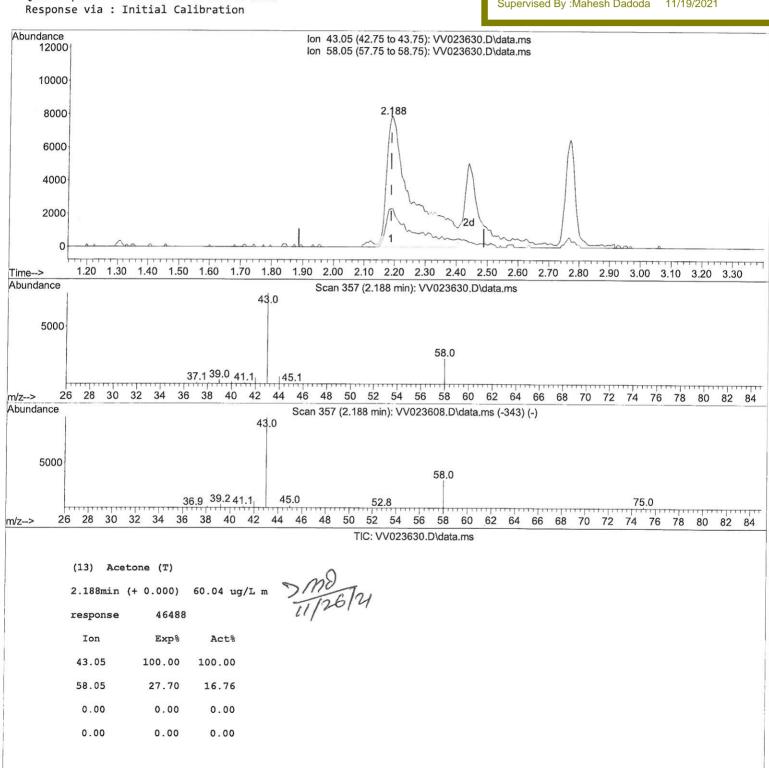
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Quant Title : TRACE VOA SFAM1.0

QLast Update : Fri Nov 19 03:51:44 2021

Instrument: MSVOA_V **LabSampleld**: VSTDCCC005E0

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111821\

Data File: VV023630.D

Acq On : 19 Nov 2021 01:38

Operator : SY/MD Sample : VSTDCCC005EC

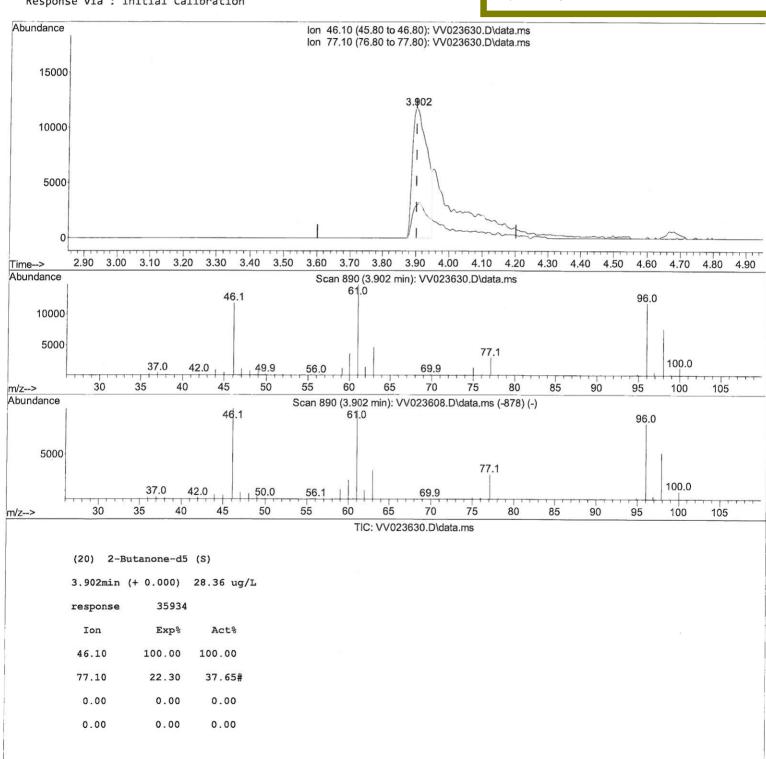
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 39 Sample Multiplier: 1

Quant Time: Nov 19 04:12:08 2021

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

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Manual IntegrationsAPPROVED



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Operator : SY/MD Sample : VSTDCCC005EC

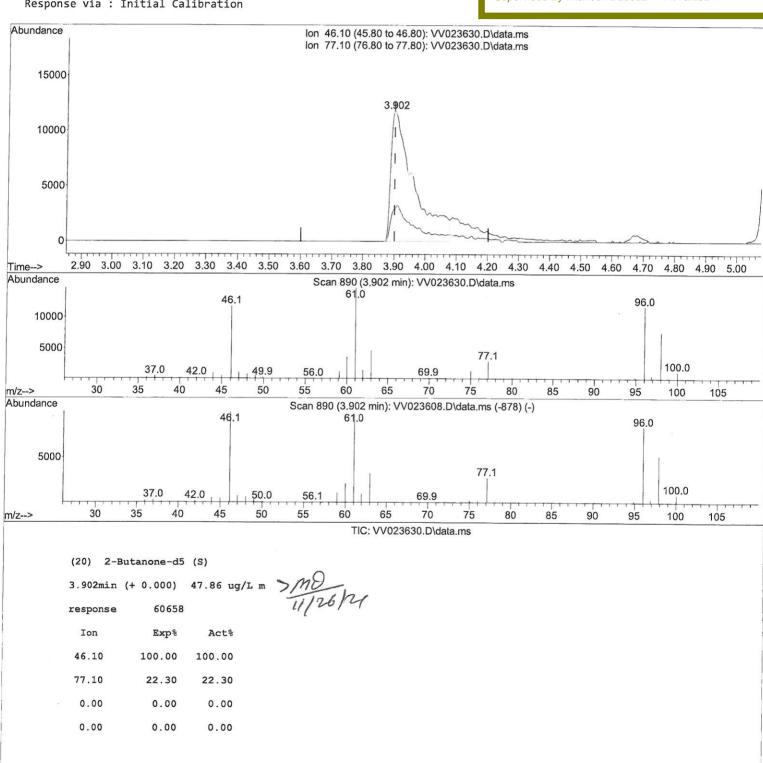
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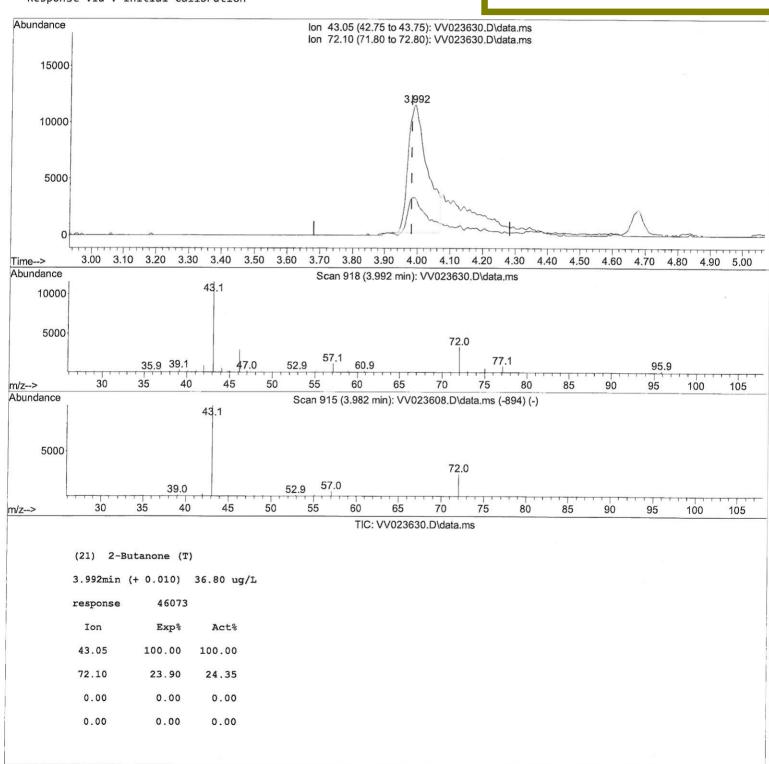
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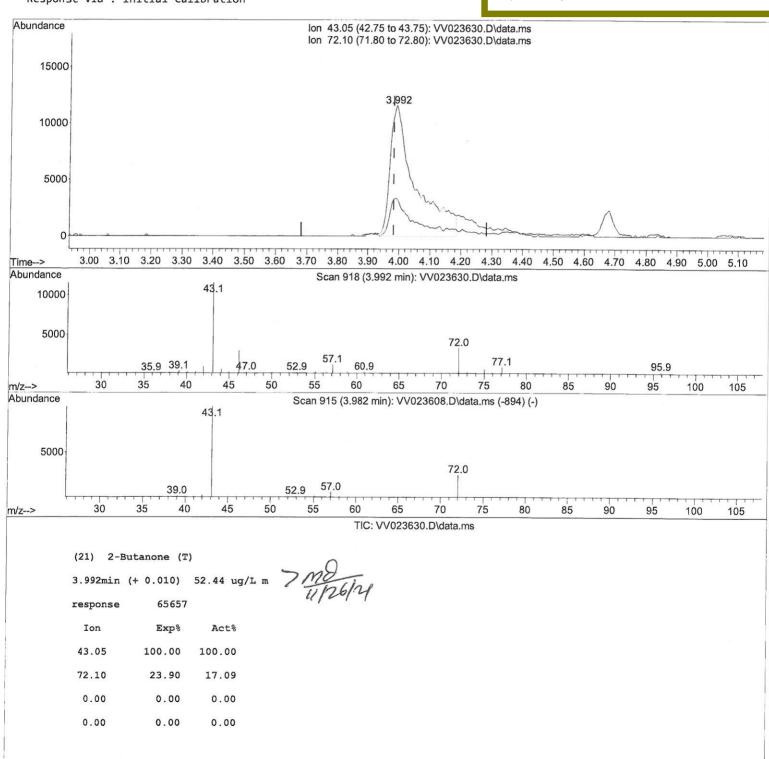
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Instrument: MSVOA_V **LabSampleld**: VSTDCCC005EC

Manual IntegrationsAPPROVED

Compound	R.T. QIon	Response Conc Units Dev(Min)
Internal Standards		
1) 1,4-Difluorobenzene	5.619 114	117419 5.000 ug/L 0.00
28) Chlorobenzene-d5	8.854 117	
58) 1,4-Dichlorobenzene-d4		6.4.T.T.
, -,	11.2-15 152	64/73 5.000 ug/L 0.00
System Monitoring Compounds		
Vinyl Chloride-d3	1.304 65	35572 4.836 ug/L 0.00
Spiked Amount 5.000	Range 40 - 130	
7) Chloroethane-d5	1.568 69	29425 4.908 ug/L 0.00
Spiked Amount 5.000	Range 65 - 130	
11) 1,1-Dichloroethene-d2	2.108 63	66469 4.827 ug/L 0.00
Spiked Amount 5.000	Range 60 - 125	Recovery = 96.600%
20) 2-Butanone-d5	3.902 46	Recovery = 96.600% 60658m 47.865 ug/L 0.00 7 mB Recovery = 95.720% 68631 4.378 ug/L 0.00
Spiked Amount 50.000	Range 40 - 130	Recovery = 95.720% $11/26/9$
24) Chloroform-d	4.349 84	68631 4.378 ug/L 0.00 /
Spiked Amount 5.000	Range 70 - 125	Recovery = 87.600%
26) 1,2-Dichloroethane-d4 Spiked Amount 5.000	5.037 65	34733 4.927 ug/L 0.00
32) Benzene-d6	Range 70 - 130	Recovery = 98.600%
Spiked Amount 5.000	5.050 84	134503 4.421 ug/L 0.00
36) 1,2-Dichloropropane-d6	Range 70 - 125 6.069 67	Recovery = 88.400% 39631 4.425 ug/L 0.00
Spiked Amount 5.000	Range 60 - 140	39631
41) Toluene-d8	7.317 98	424440
Spiked Amount 5.000	Range 70 - 130	131612 4.616 ug/L 0.00 Recovery = 92.400%
43) trans-1,3-Dichloroprop.		14670 4.320 ug/L 0.00
Spiked Amount 5.000	Range 55 - 130	Recovery = 86.400%
46) 2-Hexanone-d5	8.092 63	57508 46.025 ug/L 0.00
Spiked Amount 50.000	Range 45 - 130	Recovery = 92.040%
56) 1,1,2,2-Tetrachloroeth.		
Spiked Amount 5.000	Range 65 - 120	Recovery = 93.000%
66) 1,2-Dichlorobenzene-d4	11.625 152	49266 4.568 ug/L 0.00
Spiked Amount 5.000	Range 80 - 120	Recovery = 91.400%
Target Compounds		A 1
2) Dichlorodifluoromethane	1.127 85	Qvalue
3) Chloromethane	1.240 50	47032 4.108 ug/L 98 40234 4.133 ug/L 98
5) Vinyl chloride	1.311 62	8,
6) Bromomethane	1.520 94	42208 4.341 ug/L 97 20406 3.284 ug/L 97
8) Chloroethane	1.584 64	26562 4.734 ug/L 98
Trichlorofluoromethane		71563 4.899 ug/L 98
10) 1,1,2-Trichloro-1,2,2		35419 4.816 ug/L 95
12) 1,1-Dichloroethene	2.118 96	32771 4.680 ug/L 95
13) Acetone	2.188 43	46488m 60.035 ug/L > MO
14) Carbon disulfide	2.294 76	102778 3.890 ug/L 100 11126/1
15) Methyl Acetate	2.439 43	11019 5.028 ug/L # 88
16) Methylene chloride	2.507 84	45064 4.410 ug/L 98
17) Methyl tert-butyl Ether	2.770 73	73250 4.752 ug/L 97
18) trans-1,2-Dichloroethene	2.761 96	37707 4.381 ug/L 98
19) 1,1-Dichloroethane	3.188 63	67617 4.653 ug/L 98
21) 2-Butanone	3.992 43	65657m 52.445 ug/L 38546 4.653 ug/L # 91 11 26 2
22) cis-1,2-Dichloroethene	3.912 96	
23) Bromochloromethane	4.253 128	18508 4.845 ug/L # 79

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ALS Vial : 39 Sample Multiplier: 1

Quant Time: Nov 19 04:12:08 2021

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Quant Title : TRACE VOA SFAM1.0 QLast Update : Fri Nov 19 03:51:44 2021 Response via : Initial Calibration Instrument : MSVOA_V LabSampleId : VSTDCCC005EC

Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
25) Chloroform	4.375	83	77636	5.012 ug/L	97
27) 1,2-Dichloroethane	5.134	62	39909	4.843 ug/L	99
29) 1,1,1-Trichloroethane	4.606	97	67097	4.659 ug/L	99
30) Cyclohexane	4.677	56	51797	4.014 ug/L	96
31) Carbon tetrachloride	4.828	117	61141	4.727 ug/L	98
33) Benzene	5.101	78	152303	4.595 ug/L	100
34) Trichloroethene	5.915	95	39498	4.481 ug/L	98
35) Methylcyclohexane	6.133	83	57538	4.136 ug/L	96
37) 1,2-Dichloropropane	6.175	63	36993	4.781 ug/L	98
38) Bromodichloromethane	6.510	83	49575	4.781 ug/L	95
39) cis-1,3-Dichloropropene	7.031	75	48656	4.372 ug/L	98
40) 4-Methyl-2-pentanone	7.227	43	187001	52.111 ug/L	98
42) Toluene	7.387	91	169887	4.792 ug/L	98
44) trans-1,3-Dichloropropene	7.654	75	42748	4.629 ug/L	99
45) 1,1,2-Trichloroethane	7.841	97	26417	4.752 ug/L	97
47) Tetrachloroethene	7.976	164	34286	4.489 ug/L	96
48) 2-Hexanone	8.143	43	136323	54.214 ug/L	96
49) Dibromochloromethane	8.249	129	34914	4.956 ug/L	98
50) 1,2-Dibromoethane	8.352	107	25099	4.872 ug/L #	97
51) Chlorobenzene	8.883	112	109277	4.638 ug/L	99
52) Ethylbenzene	9.014	91	173167	4.632 ug/L	98
53) m,p-xylene	9.140	106	69385	4.729 ug/L	94
54) o-xylene	9.545	106	67010	4.868 ug/L	97
55) Styrene	9.561	104	115145	4.883 ug/L	96
57) 1,1,2,2-Tetrachloroethane	10.243	83	29181	4.791 ug/L	99
59) Bromoform	9.731	173	19509	5.043 ug/L	99
60) Isopropylbenzene	9.931	105	177164	4.766 ug/L	99
61) 1,2,3-Trichloropropane	10.275	75	20769	4.827 ug/L	96
62) 1,3,5-Trimethylbenzene	10.538	105	143888	4.669 ug/L	99
63) 1,2,4-Trimethylbenzene	10.915	105	145541	4.745 ug/L	100
64) 1,3-Dichlorobenzene	11.182	146	91064	4.795 ug/L	99
65) 1,4-Dichlorobenzene	11.275	146	89785	4.629 ug/L	99
67) 1,2-Dichlorobenzene	11.645	146	82560	4.858 ug/L	98
68) 1,2-Dibromo-3-chloropr	12.429	75	4664	5.088 ug/L	97
69) 1,3,5-Trichlorobenzene	12.645	180	68030	4.575 ug/L	98
70) 1,2,4-trichlorobenzene	13.262	180	49288	4.139 ug/L	99
71) Naphthalene	13.503	128	70420	4.011 ug/L	99
72) 1,2,3-Trichlorobenzene	13.744	180	46715	4.484 ug/L	99

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