Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File: VV023526.D

Acq On : 16 Nov 2021 12:02

Operator : SY/MD Sample : M4643-08

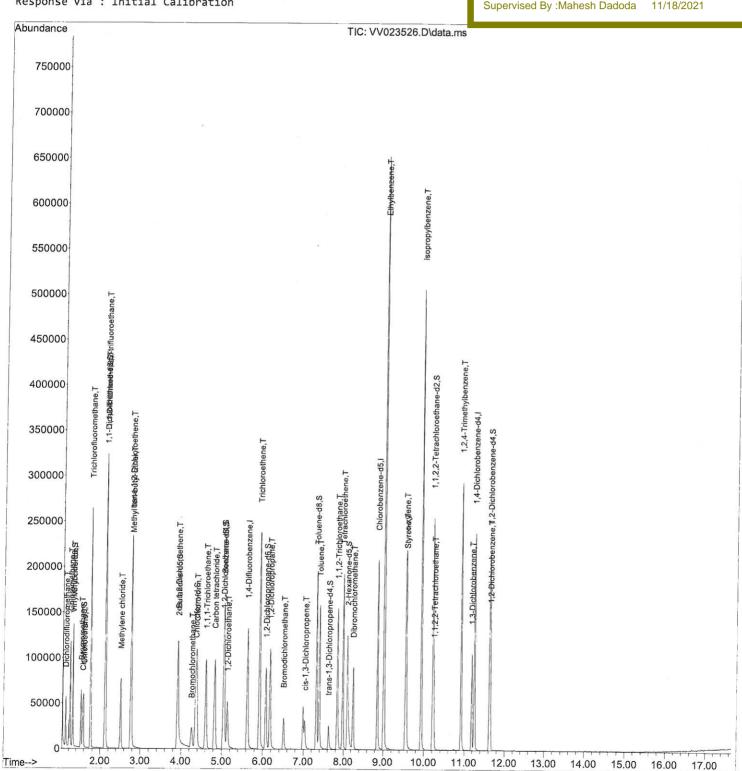
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 17 00:50:45 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId :

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File: VV023526.D

Acq On : 16 Nov 2021 12:02

Operator : SY/MD Sample : M4643-08

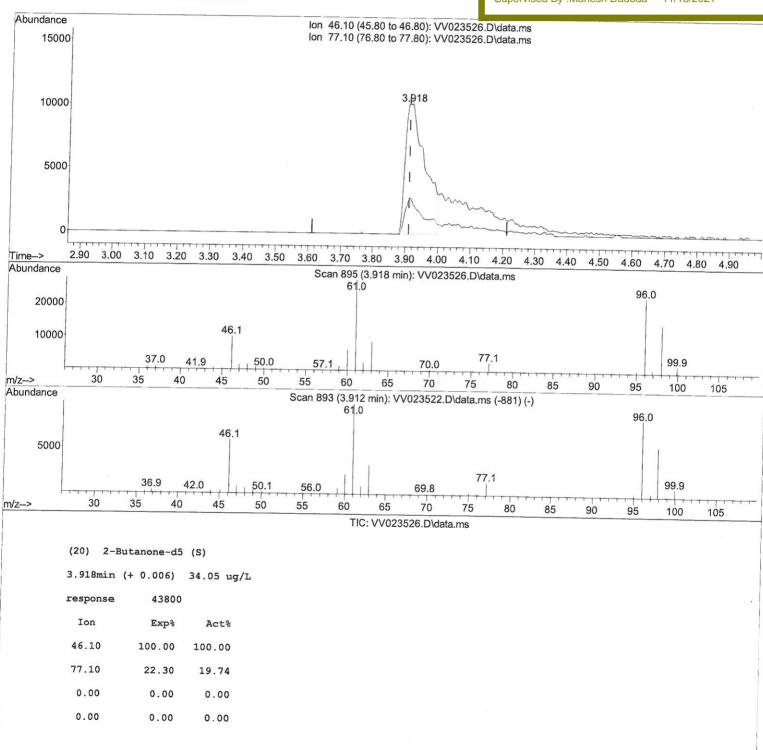
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Quant Time: Nov 17 00:50:45 2021

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Operator : SY/MD Sample : M4643-08

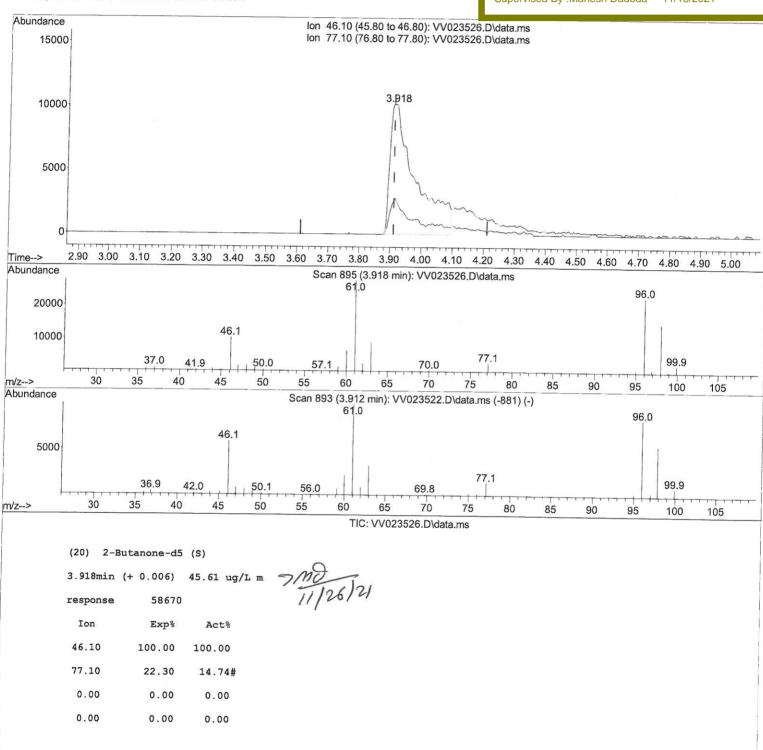
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 17 00:50:45 2021

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

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId :

Manual IntegrationsAPPROVED



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Acq On : 16 Nov 2021 12:02

Operator : SY/MD

Sample : M4643-08 Misc : 25.0mL/MSVOA_V/WATER Misc ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 17 00:50:45 2021

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

Quant Title : TRACE VOA SFAM1.0 ·QLast Update : Wed Nov 17 00:48:57 2021 Response via: Initial Calibration

Instrument : MSVOA_V ClientSampleId: GB8K3

Manual IntegrationsAPPROVED

*				
Compound	R.T. QIon	Response	Conc Units Dev(Min)
Internal Standards				
1) 1,4-Difluorobenzene	5.619 114	119182	5.000 ug/L	0.00
28) Chlorobenzene-d5	8.854 117	118164	5.000 ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249 152	63101	5.000 ug/L	0.00
, -,		03101	3.000 ag/ c	0.00
System Monitoring Compounds				
4) Vinyl Chloride-d3	1.304 65	34820	4.664 ug/L	0.00
Spiked Amount 5.000	Range 40 - 130	Recove	ry = 93.200%	
7) Chloroethane-d5	1.568 69	29312	4.817 ug/L	0.00
Spiked Amount 5.000	Range 65 - 130	Recove	ry = 96.400%	
<pre>11) 1,1-Dichloroethene-d2</pre>	2.111 63	74450	5.327 ug/L	0.00
Spiked Amount 5.000	Range 60 - 125	Recove		- MO-
20) 2-Butanone-d5	3.918 46		45.611 ug/L	0.00 7 Mg
Spiked Amount 50.000	Range 40 - 130		ry = 91.220%	- /
24) Chloroform-d	4.352 84	69950	4.396 ug/L	0.00
Spiked Amount 5.000	Range 70 - 125	Recover	(- 73)	0.00
26) 1,2-Dichloroethane-d4 Spiked Amount 5.000	5.037 65		4.999 ug/L	0.00
32) Benzene-d6	Range 70 - 130 5.050 84	Recover 145099	^y = 100.000% 4.786 ug/L	0.00
Spiked Amount 5.000	Range 70 - 125	Recover		0.00
36) 1,2-Dichloropropane-d6	6.069 67		4.965 ug/L	0.00
Spiked Amount 5.000	Range 60 - 140	Recover		0.00
41) Toluene-d8	7.317 98	127781	4.498 ug/L	0.00
Spiked Amount 5.000	Range 70 - 130	Recover		
43) trans-1,3-Dichloroprop.		15834	4.679 ug/L	0.00
Spiked Amount 5.000	Range 55 - 130	Recover		
46) 2-Hexanone-d5	8.095 63	52259	41.971 ug/L	0.00
Spiked Amount 50.000	Range 45 - 130	Recover	y = 83.940%	
56) 1,1,2,2-Tetrachloroeth.	10.217 84	30610	4.769 ug/L	0.00
Spiked Amount 5.000	Range 65 - 120	Recover	•	
66) 1,2-Dichlorobenzene-d4	11.625 152	52561	5.002 ug/L	0.00
Spiked Amount 5.000	Range 80 - 120	Recover	y = 100.000%	
Target Compounds			Qva]	Luo
2) Dichlorodifluoromethane	1.127 85	28669	2.467 ug/L	100
3) Chloromethane	1.240 50	71266	7.213 ug/L	97
5) Vinyl chloride	1.310 62	53899	5.462 ug/L	100
6) Bromomethane	1.523 94	26140	4.144 ug/L	98
8) Chloroethane	1.584 64	30000	5.268 ug/L	98
9) Trichlorofluoromethane	1.754 101	153920	10.381 ug/L	99
10) 1,1,2-Trichloro-1,2,2	. 2.117 101	37542	5.029 ug/L	95
12) 1,1-Dichloroethene	2.117 96	46546	6.549 ug/L	89
16) Methylene chloride	2.507 84	32571	3.140 ug/L	97
17) Methyl tert-butyl Ether	2.770 73	94778	6.058 ug/L	97
18) trans-1,2-Dichloroethene		61572	7.047 ug/L	99
22) cis-1,2-Dichloroethene	3.912 96	59075	7.026 ug/L #	93
23) Bromochloromethane	4.259 128	8937	2.305 ug/L #	77
25) Chloroform	4.375 83	89837	5.713 ug/L	95
27) 1,2-Dichloroethane29) 1,1,1-Trichloroethane	5.133 62 4.609 97	48398 82022	5.787 ug/L	97
31) Carbon tetrachloride	4.828 117	75455	5.715 ug/L 5.854 ug/L	99 97
34) Trichloroethene	5.915 95	80054	9.115 ug/L	94
5 ty 11 Tell for occilence	3.713 73	30034	J.IIJ UB/L	J-4

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Instrument : MSVOA_V ClientSampleId:

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Uni	ts Dev	/(Min)
37) 1,2-Dichloropropane	6.172	63	42096	5.460	ug/L	100
38) Bromodichloromethane	6.513	83	23695	2.293	A	100
39) cis-1,3-Dichloropropene	7.034	75	20592	1.857	ug/L	94
42) Toluene	7.391	91	117666	3.331	_	98
45) 1,1,2-Trichloroethane	7.837	97	50548	9.124	ug/L	95
47) Tetrachloroethene	7.976	164	44341	5.825		99
49) Dibromochloromethane	8.246	129	47825	6.813	ug/L	100
52) Ethylbenzene	9.011	91	444345	11.926	ug/L	98
54) o-xylene	9.545	106	46148	3.364	_	95
55) Styrene	9.561	104	73118	3.112		96
57) 1,1,2,2-Tetrachloroethane	10.243	83	31655	5.215	_	99
60) Isopropylbenzene	9.931	105	311800	8.611	•	99
63) 1,2,4-Trimethylbenzene	10.915	105	158004	5.287	ug/L	99
64) 1,3-Dichlorobenzene	11.181	146	44284	2.394 (•	98
67) 1,2-Dichlorobenzene	11.644	146	39555	2.389 ı		96
	·					

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