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

Data File : VV022972.D

Acq On : 21 Oct 2021 22:25

Operator : SY/MD Sample : M4265-11MSD

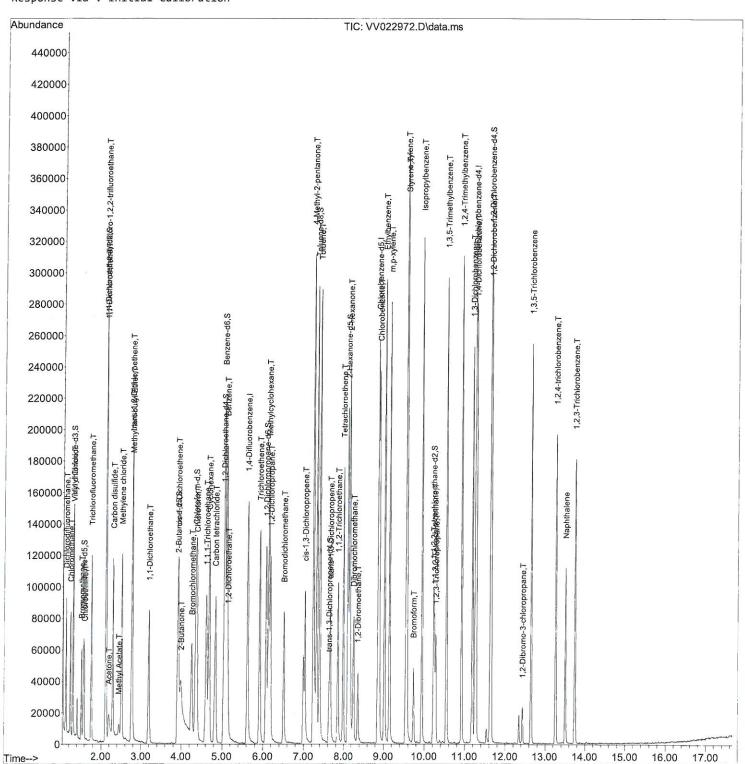
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 22 04:58:48 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:
GB7J3MSD

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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

Data File : VV022972.D

Acq On : 21 Oct 2021 22:25

Operator : SY/MD Sample : M4265-11MSD

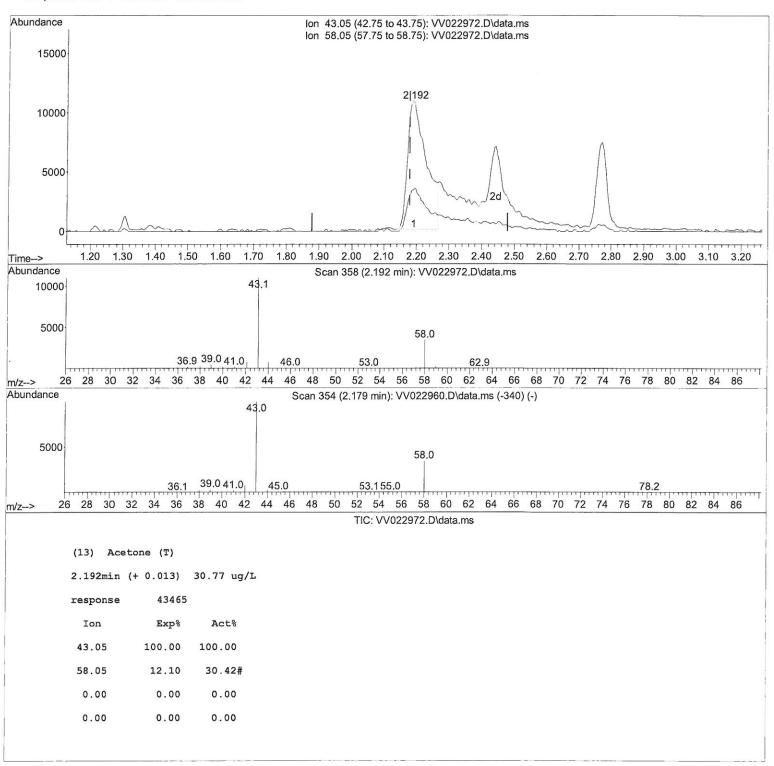
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Manual IntegrationsAPPROVED



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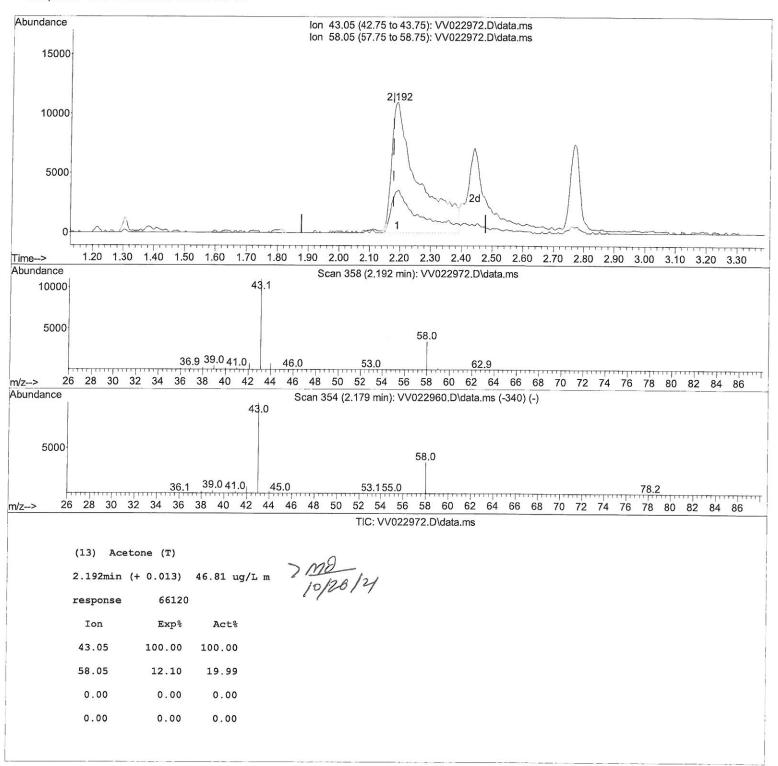
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Manual IntegrationsAPPROVED

Compound	R.T. QIon	Response Conc Units Dev((Min)
Internal Standards			.===
1) 1,4-Difluorobenzene	5.619 114	140414 5.000 ug/L	0.00
28) Chlorobenzene-d5	8.854 117	138946 5.000 ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.249 152	73706 5.000 ug/L	0.00
System Monitoring Compounds		*	
4) Vinyl Chloride-d3	1.304 65		0.00
Spiked Amount 5.000	Range 40 - 13		5
7) Chloroethane-d5	1.568 69	Ç.	0.00
Spiked Amount 5.000	Range 65 - 130		
11) 1,1-Dichloroethene-d2	2.108 63	0.	0.00
Spiked Amount 5.000	Range 60 - 12!		
20) 2-Butanone-d5	3.902 46	O.	0.01
Spiked Amount 50.000	Range 40 - 130		
24) Chloroform-d	4.349 84		0.00
Spiked Amount 5.000	Range 70 - 125	4.50	
26) 1,2-Dichloroethane-d4 Spiked Amount 5.000	5.034 65 Range 70 - 130		0.00
32) Benzene-d6	5.050 84		0.00
Spiked Amount 5.000	Range 70 - 125		
36) 1,2-Dichloropropane-d6	6.069 67		0.00
Spiked Amount 5.000	Range 60 - 140		
41) Toluene-d8	7.317 98		0.00
Spiked Amount 5.000	Range 70 - 136		
43) trans-1,3-Dichloroprop			0.00
Spiked Amount 5.000	Range 55 - 136		
46) 2-Hexanone-d5	8.092 63		0.00
Spiked Amount 50.000	Range 45 - 136		
56) 1,1,2,2-Tetrachloroeth	. 10.217 84	44895 5.410 ug/L	0.00
Spiked Amount 5.000	Range 65 - 126		
66) 1,2-Dichlorobenzene-d4	11.625 152	69647 5.096 ug/L	0.00
Spiked Amount 5.000	Range 80 - 120	Recovery = 102.000%	
Target Compounds		Qva	lue
Dichlorodifluoromethane	1.127 85	44462 4.345 ug/L	99
Chloromethane	1.240 50	49025 4.766 ug/L	98
Vinyl chloride	1.311 62	51271 4.825 ug/L	99
6) Bromomethane	1.520 94	23137 3.492 ug/L	100
8) Chloroethane	1.584 64	27696 4.212 ug/L	100
	1.751 101	69382 4.457 ug/L	100
10) 1,1,2-Trichloro-1,2,2		35810 3.955 ug/L	98
12) 1,1-Dichloroethene	2.118 96	32037 3.777 ug/L	84 md
13) Acetone	2.192 43	66120m 46.811 ug/L	20 146/21
14) Carbon disulfide	2.291 76	124319 5.215 ug/L	99 10/26/21 96 10/26/21
15) Methyl Acetate16) Methylene chloride	2.442 43 2.507 84	13374 3.309 ug/L 49129 4.208 ug/L	94
17) Methyl tert-butyl Ether	2.770 73	87880 4.282 ug/L	95
18) trans-1,2-Dichloroethene		47471 5.147 ug/L	97
19) 1,1-Dichloroethane	3.192 63	85581 5.151 ug/L	100
21) 2-Butanone	3.986 43	77028 33.306 ug/L	90
22) cis-1,2-Dichloroethene	3.912 96	49067 5.066 ug/L	94
23) Bromochloromethane	4.253 128	22520 5.234 ug/L	93
			#1000 M

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Operator : SY/MD : M4265-11MSD Sample

Misc : 25.0mL/MSVOA V/WATER ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 22 04:58:48 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 : GB7J3MSD

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
25) Chloroform	4.375	83	89115	4.627 ug/L	100
27) 1,2-Dichloroethane	5.134	62	47080	4.898 ug/L	98
29) 1,1,1-Trichloroethane	4.609	97	77200	4.866 ug/L	98
30) Cyclohexane	4.677	56	65565	4.478 ug/L	99
31) Carbon tetrachloride	4.828	117	67378	4.892 ug/L	99
33) Benzene	5.101	78	192701	5.150 ug/L	100
34) Trichloroethene	5.915	95	47578	4.675 ug/L	99
35) Methylcyclohexane	6.130	83	70206	4.924 ug/L	94
37) 1,2-Dichloropropane	6.175	63	46395	5.155 ug/L	100
38) Bromodichloromethane	6.510	83	58786	5.301 ug/L	99
39) cis-1,3-Dichloropropene	7.031	75	60492	5.077 ug/L	99
40) 4-Methyl-2-pentanone	7.230	43	243286	50.308 ug/L	98
42) Toluene	7.387	91	211035	5.555 ug/L	99
44) trans-1,3-Dichloropropene	7.654	75	50559	5.205 ug/L	96
45) 1,1,2-Trichloroethane	7.841	97	33832	5.066 ug/L	98
47) Tetrachloroethene	7.976	164	40620	4.936 ug/L	100
48) 2-Hexanone	8.143	43	170928	48.341 ug/L	99
49) Dibromochloromethane	8.249	129	40549	5.187 ug/L	88
50) 1,2-Dibromoethane	8.355	107	30930	4.960 ug/L	98
51) Chlorobenzene	8.883	112	131334	5.261 ug/L	98
52) Ethylbenzene	9.014	91	200305	5.193 ug/L	98
53) m,p-xylene	9.140	106	79404	5.246 ug/L	98
54) o-xylene	9.545	106	76194	5.351 ug/L	99
55) Styrene	9.561	104	130593	5.322 ug/L	97
57) 1,1,2,2-Tetrachloroethane	10.243	83	38338	5.674 ug/L	100
59) Bromoform	9.731	173	21899	5.304 ug/L	100
60) Isopropylbenzene	9.934	105	200726	5.362 ug/L	99
61) 1,2,3-Trichloropropane	10.275	75	26183	5.113 ug/L	98
62) 1,3,5-Trimethylbenzene	10.542	105	158279	5.233 ug/L	99
63) 1,2,4-Trimethylbenzene	10.915	105	162956	5.383 ug/L	98
64) 1,3-Dichlorobenzene	11.182	146	104570	5.322 ug/L	99
65) 1,4-Dichlorobenzene	11.275	146	104122	5.219 ug/L	98
67) 1,2-Dichlorobenzene	11.645	146	96631	5.243 ug/L	100
68) 1,2-Dibromo-3-chloropr	12.432	75	5516	5.918 ug/L	89
69) 1,3,5-Trichlorobenzene	12.644	180	82439	5.570 ug/L	95
70) 1,2,4-trichlorobenzene	13.262	180	59215	5.237 ug/L	99
71) Naphthalene	13.503	128	88989	4.952 ug/L	98
72) 1,2,3-Trichlorobenzene	13.744	180	57180	5.363 ug/L	97

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