









Data Path : Z:\voasrv\HPCHEM1\ Data File : VX025091.D Acq On : 08 Nov 2021 10:36 Dperator : JC/MD Sample : VSTD10027 Misc : 5.0mL/MSVOA_X/WATEF ALS Vial : 5 Sample Multipl:	2	ata\V)	X110821\		Instrument : MSVOA_X ClientSampleId : VSTD100627 Manual IntegrationsAPPROVED
Quant Time: Nov 09 03:36:34 202 Quant Method : Z:\voasrv\HPCHEM Quant Title : VOC Analysis QLast Update : Tue Nov 09 03:33 Response via : Initial Calibrat	11\MSVOA_3	X\Meth	lod∖SFAMXL	M110821WMA.M	Reviewed By :John Carlone 11/09/2021 Supervised By :Mahesh Dadoda 11/09/2021
Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
1) 1,4-Difluorobenzene	6 762	114	107250	EQ. 000	# 0. 00
28) Chlorobenzene-d5	10.055			50.000 ug/L 50.000 ug/L	# 0.00 0.00
58) 1,4-Dichlorobenzene-d4			87880	50.000 ug/L	0.00
56) 1,4-Dichio obenzene-u4	12.024	172	87880	50.000 ug/L	0.00
System Monitoring Compounds					
4) Vinyl Chloride-d3	1.368	65	193542	151.803 ug/L	0.00
Chloroethane-d5	1.666	69	100847	115.426 ug/L	0.00
11) 1,1-Dichloroethene-d2	2.313	63	376839	127.450 ug/L	0.00
21) 2-Butanone-d5	4.459	46	267248	228.035 ug/L	0.00
24) Chloroform-d	5.062	84	401063	128.215 ug/L	0.00
26) 1,2-Dichloroethane-d4	5.964	65	228176	101.130 ug/L	0.00
32) Benzene-d6	5.977	84	532343	100.475 ug/L	0.00
36) 1,2-Dichloropropane-d6	7.312	67	173272	103.478 ug/L	0.00
41) Toluene-d8	8.653 8.952	98	499857	105.235 ug/L	0.00
<pre>43) trans-1,3-Dichloroprop 47) 2-Hexanone-d5</pre>	8.952 9.391	79 63	102668	105.798 ug/L 200.179 ug/L	0.00 9 m/ 09 2) 0.00 9 m/ 09 2)
56) 1,1,2,2-Tetrachloroeth		84	239726	107.390 ug/L	0.00
66) 1,2-Dichlorobenzene-d4	12.323		171014	100.790 ug/L	0.00
	12.525	192	1/101+	100.750 0672	0.00
Target Compounds				Qva	alue
Dichlorodifluoromethane	1.166	85	213089	120.827 ug/L	98
Chloromethane	1.288	50	147754	120.278 ug/L	90
5) Vinyl chloride	1.374	62	182746	129.531 ug/L	99
6) Bromomethane	1.611	94	97828	107.769 ug/L	95
8) Chloroethane	1.685			140.180 ug/L	99
9) Trichlorofluoromethane	1.886	101	327207	108.593 ug/L	99
10) 1,1,2-Trichloro-1,2,2 12) 1,1-Dichloroethene	2.331 2.319	101 96	171326 152797	123.698 ug/L 128.828 ug/L	92 72
13) Acetone	2.315	43	243076	216.568 ug/L	98 0
14) Carbon disulfide	2.514	76	402345	127.143 ug/L	99 000 10 1
15) Methyl Acetate	2.709	43	180331m	112.300 ug/L	2 1 9 2)
16) Methylene chloride	2.794	84	174751	129.853 ug/L	86 11011
17) trans-1,2-Dichloroethene	3.093	96	157706	131.218 ug/L	94
18) Methyl tert-butyl Ether	3.117	73	621496	126.489 ug/L #	93
19) 1,1-Dichloroethane	3.611	63	328613	119.382 ug/L	93
20) cis-1,2-Dichloroethene	4.495	96	183764	133.900 ug/L	88
22) 2-Butanone	4.562	43	308517	224.721 ug/L	85
23) Bromochloromethane	4.910	128	86466	128.345 ug/L #	74
25) Chloroform 27) 1,2-Dichloroethane	5.099 6.092	83 62	369838 259371	120.684 ug/L 97.782 ug/L	98
29) Cyclohexane	5.477	56	268938	109.920 ug/L #	98 78
30) 1,1,1-Trichloroethane	5.385	97	336244	109.491 ug/L #	93
31) Carbon tetrachloride	5.684	117	229983	88.704 ug/L	99
33) Benzene	6.044	78	539719	94.756 ug/L	100
34) Trichloroethene	7.129	95	150042	92.750 ug/L	99
35) Methylcyclohexane	7.385	83	231129	99.582 ug/L #	87
<pre>37) 1,2-Dichloropropane</pre>	7.434	63	144913	94.976 ug/L #	96
38) Bromodichloromethane	7.824	83	223140	94.448 ug/L #	94
39) cis-1,3-Dichloropropene	8.372	75	243736	99.863 ug/L	98
40) 4-Methyl-2-pentanone	8.580	43	473586	202.091 ug/L #	78
42) Toluene	8.720	91 75	594081	101.227 ug/L	99
44) trans-1,3-Dichloropropene	8.982	75	257071	100.321 ug/L	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110821\ Data File : VX025091.D Acq On : 08 Nov 2021 10:36 Dperator : JC/MD Sample : VSTD10027 Misc : 5.0mL/MSVOA_X/WATER ALS Vial : 5 Sample Multiplier: 1										
<pre>Juant Time: Nov 09 03:36:34 2021 Juant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM110821WMA.M Juant Title : VOC Analysis JLast Update : Tue Nov 09 03:33:09 2021 Response via : Initial Calibration</pre>										
Compound	R.T.	QIon	Response	Conc Un	its D	ev(I	Min)			
<pre>45) 1,1,2-Trichloroethane 46) Tetrachloroethene 48) 2-Hexanone 49) Dibromochloromethane 50) 1,2-Dibromoethane 51) Chlorobenzene 52) Ethylbenzene 53) m,p-Xylene 54) o-Xylene 55) Styrene 57) 1,1,2,2-Tetrachloroethane 59) Bromoform 60) Isopropylbenzene 61) 1,2,3-Trichloropropane 62) 1,3,5-Trimethylbenzene 63) 1,2,4-Trimethylbenzene 64) 1,3-Dichlorobenzene 65) 1,4-Dichlorobenzene 68) 1,2-Dibromo-3-chloropr 69) 1,3,5-Trichlorobenzene 70) 1,2,4-trichlorobenzene</pre>	9.153 9.275 9.433 9.525 9.610 10.086 10.195 10.305 10.646 10.659 11.213 10.805 10.963 11.244 11.457 11.756 11.969 12.043 12.335 12.945 13.116 13.591	43 129 107	136660 91345 393677 153559 150138 358546 699622 249717 238000 421299 231967 99350 686687 202321 593844 605306 266634 261133 260520 64478 179376 160679	93.360 90.134 199.585 95.392 94.088 100.396 104.330 107.184 105.338 109.690 102.429 92.403 100.061 95.601 103.654 104.563 103.258 98.201 98.454 92.420 93.802	ug/L ug/L ug/L ug/L ug/L ug/L ug/L ug/L	#	93 98 79 97 94 100 96 95 98 97 96 98 92 100 98 98 98 95 85 99			
71) Naphthalene72) 1,2,3-Trichlorobenzene	13.780 13.963		622858	96.211 105.810 93.645	ug/L		97 99 95			

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

AMXLM110821WMA.M Tue Nov 09 03:55:30 2021

Instrument : MSVOA_X ClientSampleId : /STD100627

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 11/09/2021 Supervised By :Mahesh Dadoda 11/09/2021

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