

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL080421\
 Data File : VL037325.D
 Acq On : 5 Aug 2021 2:36
 Operator : SY/AP
 Sample : M3211-03
 Misc : 400mL/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SVP-3

Manual Integrations
 APPROVED

MMDadoda
 8/5/2021 1:40:37 PM

Quant Time: Aug 05 10:02:56 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL080421AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LWed Aug 04
 QLast Update : Wed Aug 04 22:53:01 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.66	49	1105797	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.17	114	3605754	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.07	117	3056855	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.40	95	2144032	10.27	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	102.70%

Target Compounds					Ovalue
2) Dichlorodifluoromethane	3.04	85	36814	0.164 ppbv	96
3) Chlorodifluoromethane	2.98	51	35512	0.158 ppbv	92
4) Chloromethane	3.13	50	15740	0.195 ppbv #	70
9) Propene	3.02	41	2211590m	27.204 ppbv	
10) Heptane	8.11	43	1209163	6.368 ppbv	99
11) Trichlorofluoromethane	3.94	101	60225	0.292 ppbv	97
13) Ethanol	3.61	45	1052040	85.916 ppbv	92
15) Acetone	3.84	43	10885648m	90.625 ppbv	
17) tert-Butyl alcohol	4.30	59	276296	2.290 ppbv #	72
19) Isopropyl Alcohol	3.97	45	1222856	18.395 ppbv #	93
20) Methylene Chloride	4.34	84	831617	8.362 ppbv	98
26) Hexane	5.68	57	3761084	21.519 ppbv #	47
27) Carbon Disulfide	4.55	76	238565	1.677 ppbv	97
29) Chloroform	5.75	83	101662	0.459 ppbv	90
30) Cyclohexane	7.12	84	61564	0.457 ppbv #	81
32) 1,1,1-Trichloroethane	6.50	97	10601m	0.050 ppbv	
34) 2-Butanone	5.23	43	4369794	34.924 ppbv	98
35) Carbon Tetrachloride	7.02	117	7924m	0.037 ppbv	
36) Benzene	6.89	78	236183	0.728 ppbv	97
38) Trichloroethene	7.83	130	8263m	0.058 ppbv	
40) 1,4-Dioxane	7.83	88	67373m	1.938 ppbv	
41) Tetrahydrofuran	6.03	42	7115814	102.262 ppbv	99
44) 2,2,4-Trimethylpentane	7.86	57	77749m	0.145 ppbv	
50) 4-Methyl-2-Pentanone	8.73	43	278473	1.453 ppbv	97
51) 2-Hexanone	10.10	43	613018m	7.037 ppbv	
52) Tetrachloroethene	11.21	164	669682	4.792 ppbv	99
53) Toluene	9.80	91	757446	1.977 ppbv	99
57) Chlorobenzene	12.18	112	30946m	0.120 ppbv	
58) Ethyl Benzene	12.70	91	201749	0.449 ppbv	99
59) m/p-Xylene	12.95	91	497245	1.315 ppbv	98
60) o-Xylene	13.68	91	215257	0.626 ppbv	91
61) Styrene	13.51	104	30330m	0.161 ppbv	
63) 1,1,2,2-Tetrachloroethane	13.65	83	10257m	0.041 ppbv	
64) n-propylbenzene	15.54	120	14820m	0.114 ppbv	
70) n-Butylbenzene	18.54	91	51937m	0.109 ppbv	
72) 4-Ethyltoluene	15.82	105	71742m	0.173 ppbv	
73) 1,3,5-Trimethylbenzene	15.98	105	75521m	0.202 ppbv	
74) 1,2,4-Trimethylbenzene	16.74	105	210705m	0.529 ppbv	
75) 1,3-Dichlorobenzene	16.98	146	110097m	0.450 ppbv	
79) Naphthalene	22.19	128	64061m	0.175 ppbv	
80) Naphthalene,2-methyl-	24.96	142	11849	0.200 ppbv #	70

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

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