

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL081820\
 Data File : VL035452.D
 Acq On : 18 Aug 2020 21:48
 Operator : SY/AP
 Sample : L3725-04
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SS-04

Manual Integrations
 APPROVED

MMDadoda
 8/20/2020 11:07:24 AM

Quant Time: Aug 19 11:15:27 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL081720AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Mon Aug 18 2020
 QLast Update : Mon Aug 17 19:28:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.69	49	1152614	10.00	ppbv	0.01
33) 1,4-Difluorobenzene	7.20	114	5381964	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.11	117	5488377	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.45	95	4049917	10.54	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	105.40%

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	3.05	85	55392	0.259	ppbv	95
3) Chlorodifluoromethane	2.99	51	38231	0.265	ppbv	100
4) Chloromethane	3.14	50	19305	0.256	ppbv	92
5) Vinyl Chloride	3.26	62	4838m	0.054	ppbv	
9) Propene	3.01	41	1743709	30.018	ppbv	99
10) Heptane	8.14	43	292389	1.920	ppbv #	82
11) Trichlorofluoromethane	3.96	101	108785	0.325	ppbv	99
13) Ethanol	3.62	45	2715163	127.578	ppbv	99
15) Acetone	3.85	43	10511405	79.232	ppbv #	38
17) tert-Butyl alcohol	4.31	59	1203234	7.017	ppbv	99
20) Methylene Chloride	4.36	84	618796	5.781	ppbv #	63
26) Hexane	5.70	57	676458	4.643	ppbv #	33
27) Carbon Disulfide	4.56	76	1016206	4.826	ppbv #	85
28) Methyl tert-Butyl Ether	5.05	73	30897	0.116	ppbv #	1
29) Chloroform	5.78	83	247290	0.845	ppbv	100
30) Cyclohexane	7.15	84	208848	1.292	ppbv #	62
34) 2-Butanone	5.25	43	6890187	38.318	ppbv #	54
35) Carbon Tetrachloride	7.03	117	11717m	0.041	ppbv	
36) Benzene	6.91	78	1025924	2.597	ppbv #	88
41) Tetrahydrofuran	6.06	42	2588820	26.100	ppbv #	53
43) Methyl Methacrylate	8.03	69	298491	1.932	ppbv #	64
44) 2,2,4-Trimethylpentane	7.89	57	353717	0.686	ppbv #	87
50) 4-Methyl-2-Pentanone	8.76	43	4800023	18.609	ppbv #	76
51) 2-Hexanone	10.16	43	778923	3.583	ppbv #	77
52) Tetrachloroethene	11.25	164	50639	0.210	ppbv #	85
53) Toluene	9.83	91	4600899	9.043	ppbv	97
58) Ethyl Benzene	12.74	91	1421662	2.146	ppbv	93
59) m/p-Xylene	13.00	91	3989795	7.124	ppbv #	82
60) o-Xylene	13.72	91	1580522	2.904	ppbv	88
61) Styrene	13.56	104	304133	0.691	ppbv #	87
62) Isopropylbenzene	14.70	105	227216	0.255	ppbv	96
64) n-propylbenzene	15.59	120	218874m	0.796	ppbv	
65) tert-Butylbenzene	16.80	119	494281	0.570	ppbv #	89
67) sec-Butylbenzene	17.33	105	182084	0.160	ppbv	95
69) p-Isopropyltoluene	17.69	119	313796	0.310	ppbv	93
70) n-Butylbenzene	18.61	91	549131	0.574	ppbv #	31
72) 4-Ethyltoluene	15.87	105	1179764	1.652	ppbv #	91
73) 1,3,5-Trimethylbenzene	16.03	105	1151521	1.732	ppbv	90
74) 1,2,4-Trimethylbenzene	16.80	105	3887158	5.382	ppbv #	70
79) Naphthalene	22.26	128	1257597	1.237	ppbv	98
80) Naphthalene,2-methyl-	25.01	142	426796	1.015	ppbv	96

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

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