

Data Path : W:\HPCHEM1\MSVOA L\DATA\VL012218\
 Data File : VL031427.D
 Acq On : 22 Jan 2018 22:27
 Operator : SY/AP
 Sample : J1089-09
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 MDL-AIR-2

Manual Integrations
 APPROVED

MMDadoda
 1/24/2018 5:10:23 PM

Quant Time: Jan 23 10:26:56 2018
 Quant Method : W:\HPCHEM1\MSVOA L\METHODS\VL012218AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Mon Jan 22 2018
 QLast Update : Mon Jan 22 17:52:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.90	49	1205423	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.44	114	2730302	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.41	117	2639952	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.75	95	2203725	10.42	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	104.20%

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	3.17	85	34840	0.20	ppbv	91
3) Chlorodifluoromethane	3.11	51	21035	0.13	ppbv	99
4) Chloromethane	3.27	50	10126	0.14	ppbv	94
5) Vinyl Chloride	3.40	62	9570	0.14	ppbv	90
6) Bromomethane	3.62	94	7079m	0.14	ppbv	
7) Chloroethane	3.72	64	4167	0.14	ppbv #	78
8) Dichlorotetrafluoroethane	3.32	85	27684	0.15	ppbv	97
9) Propene	3.14	41	7465m	0.13	ppbv	
10) Heptane	8.40	43	15464	0.09	ppbv	98
11) Trichlorofluoromethane	4.13	101	32017	0.14	ppbv	94
12) 1,1,2-Trichlorotrifluoroet	4.68	101	21004	0.15	ppbv	99
13) Ethanol	3.79	45	4070m	0.24	ppbv	
14) Bromoethene	3.90	108	7728m	0.13	ppbv	
15) Acetone	4.04	43	32716	0.20	ppbv	93
16) 1,3-Butadiene	3.47	39	8168	0.12	ppbv	97
17) tert-Butyl alcohol	4.52	59	11346m	0.15	ppbv	
18) 1,1-Dichloroethene	4.48	96	8646	0.14	ppbv	85
19) Isopropyl Alcohol	4.18	45	13096	0.13	ppbv #	96
20) Methylene Chloride	4.54	84	16663	0.27	ppbv	94
21) Allyl Chloride	4.61	41	12055	0.13	ppbv #	84
22) trans-1,2-Dichloroethene	5.10	96	9777m	0.15	ppbv	
23) Vinyl Acetate	5.30	43	19323	0.10	ppbv #	93
24) 1,1-Dichloroethane	5.22	63	21661	0.13	ppbv	98
25) Ethyl Acetate	5.92	43	34292m	0.12	ppbv	
26) Hexane	5.92	57	15049m	0.12	ppbv	
27) Carbon Disulfide	4.76	76	19510	0.12	ppbv #	77
28) Methyl tert-Butyl Ether	5.28	73	17103	0.11	ppbv	100
29) Chloroform	5.98	83	28607m	0.13	ppbv	
30) Cyclohexane	7.41	84	15936m	0.16	ppbv	
31) cis-1,2-Dichloroethene	5.78	61	13662	0.11	ppbv #	77
32) 1,1,1-Trichloroethane	6.77	97	25324	0.12	ppbv	90
34) 2-Butanone	5.49	43	17458	0.10	ppbv	95
35) Carbon Tetrachloride	7.29	117	27959	0.13	ppbv	95
36) Benzene	7.16	78	27683	0.11	ppbv	98
37) 1,2-Dichloroethane	6.56	62	23374	0.12	ppbv #	97
38) Trichloroethene	8.12	130	13941m	0.15	ppbv	
39) 1,2-Dichloropropane	7.89	63	10287	0.11	ppbv #	81
40) 1,4-Dioxane	8.20	88	3665m	0.10	ppbv	
41) Tetrahydrofuran	6.34	42	9790m	0.10	ppbv	
42) Bromodichloromethane	8.07	83	28393m	0.13	ppbv	
43) Methyl Methacrylate	8.30	69	7761m	0.09	ppbv	

Data Path : W:\HPCHEM1\MSVOA L\DATA\VL012218\
 Data File : VL031427.D
 Acq On : 22 Jan 2018 22:27
 Operator : SY/AP
 Sample : J1089-09
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 MDL-AIR-2

Manual Integrations
 APPROVED

MMDadoda
 1/24/2018 5:10:23 PM

Quant Time: Jan 23 10:26:56 2018
 Quant Method : W:\HPCHEM1\MSVOA L\METHODS\VL012218AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Mon Jan 22 2018
 QLast Update : Mon Jan 22 17:52:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,2,4-Trimethylpentane	8.15	57	37939	0.09	ppbv	# 82
45) t-1,3-Dichloropropene	9.58	75	11038m	0.09	ppbv	
46) cis-1,3-Dichloropropene	9.00	75	13807m	0.09	ppbv	
47) 1,1,2-Trichloroethane	9.78	97	13511m	0.13	ppbv	
48) Dibromochloromethane	10.61	129	20952m	0.11	ppbv	
49) Bromoform	13.37	173	15654m	0.11	ppbv	
50) 4-Methyl-2-Pentanone	9.07	43	22133m	0.10	ppbv	
51) 2-Hexanone	10.47	43	16650m	0.08	ppbv	
52) Tetrachloroethene	11.54	164	11520	0.13	ppbv	# 75
53) Toluene	10.11	91	24452m	0.09	ppbv	
54) 1,2-Dibromoethane	10.93	107	18753m	0.11	ppbv	
56) 1,1,1,2-Tetrachloroethane	12.45	131	17248m	0.14	ppbv	
57) Chlorobenzene	12.47	112	31136	0.14	ppbv	91
58) Ethyl Benzene	13.03	91	35246	0.09	ppbv	90
59) m/p-Xylene	13.32	91	64134m	0.19	ppbv	
60) o-Xylene	14.02	91	28175m	0.09	ppbv	
61) Styrene	13.85	104	5969m	0.08	ppbv	
62) Isopropylbenzene	14.99	105	41431	0.10	ppbv	97
63) 1,1,2,2-Tetrachloroethane	14.00	83	27998m	0.12	ppbv	
64) n-propylbenzene	15.89	120	9526m	0.09	ppbv	
65) tert-Butylbenzene	17.08	119	30391m	0.08	ppbv	
66) Benzyl Chloride	17.32	91	3786m	0.10	ppbv	
67) sec-Butylbenzene	17.63	105	46588	0.09	ppbv	96
69) p-Isopropyltoluene	17.98	119	35279m	0.09	ppbv	
70) n-Butylbenzene	18.89	91	38238	0.09	ppbv	97
71) 2-Chlorotoluene	15.79	91	29474m	0.09	ppbv	
72) 4-Ethyltoluene	16.18	105	28876m	0.09	ppbv	
73) 1,3,5-Trimethylbenzene	16.32	105	25340m	0.08	ppbv	
74) 1,2,4-Trimethylbenzene	17.09	105	30507m	0.09	ppbv	
75) 1,3-Dichlorobenzene	17.34	146	30352m	0.14	ppbv	
76) 1,4-Dichlorobenzene	17.48	146	25919m	0.13	ppbv	
77) 1,2-Dichlorobenzene	18.16	146	23871	0.12	ppbv	96
78) Hexachloro-1,3-Butadiene	23.70	225	13698m	0.19	ppbv	
79) Naphthalene	22.62	128	17449m	0.08	ppbv	
80) Naphthalene, 2-methyl-	25.26	142	3925	0.07	ppbv	94
81) 1,2,4-Trichlorobenzene	22.34	180	13148m	0.12	ppbv	

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

