

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL061621\
 Data File : VL037165.D
 Acq On : 16 Jun 2021 10:04
 Operator : SY/AP
 Sample : VSTDIC002
 Misc : 400mL/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC002

Manual Integrations
 APPROVED

MMDadoda
 6/17/2021 5:37:15 PM

Quant Time: Jun 16 12:20:31 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL061621AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LWed Jun 16 12:20:31 2021
 QLast Update : Wed Jun 16 11:33:05 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.67	49	1572013	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.18	114	4642565	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.10	117	4448230	10.00	ppbv	0.00

System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	14.43	95	3239582	9.89	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	98.90%

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	3.05	85	514193	2.054	ppbv	97
3) Chlorodifluoromethane	2.99	51	641221	2.174	ppbv	99
4) Chloromethane	3.14	50	212698	2.117	ppbv	98
5) Vinyl Chloride	3.26	62	234536	2.413	ppbv	98
6) Bromomethane	3.47	94	131990	2.135	ppbv	98
7) Chloroethane	3.56	64	85338m	2.250	ppbv	
8) Dichlorotetrafluoroethane	3.19	85	572876	2.142	ppbv	98
9) Propene	3.01	41	209125	2.150	ppbv	100
10) Heptane	8.13	43	516080	2.226	ppbv	98
11) Trichlorofluoromethane	3.95	101	605599	2.132	ppbv	99
12) 1,1,2-Trichlorotrifluoroet	4.49	101	429660	2.205	ppbv	98
13) Ethanol	3.62	45	33486m	2.492	ppbv	
14) Bromoethene	3.74	108	177010	2.099	ppbv	94
15) Acetone	3.86	43	432819	2.301	ppbv	# 84
16) 1,3-Butadiene	3.33	39	227092	2.236	ppbv	98
17) tert-Butyl alcohol	4.30	59	366702	2.253	ppbv	98
18) 1,1-Dichloroethene	4.29	96	185693	2.112	ppbv	99
19) Isopropyl Alcohol	3.98	45	234549m	2.272	ppbv	
20) Methylene Chloride	4.35	84	265031m	3.233	ppbv	
21) Allyl Chloride	4.42	41	288152m	2.293	ppbv	
22) trans-1,2-Dichloroethene	4.89	96	192686	2.250	ppbv	98
23) Vinyl Acetate	5.08	43	561873	2.275	ppbv	100
24) 1,1-Dichloroethane	5.02	63	438064	2.210	ppbv	99
25) Ethyl Acetate	5.69	43	838234	2.274	ppbv	98
26) Hexane	5.69	57	420968	2.297	ppbv	95
27) Carbon Disulfide	4.56	76	404107	2.174	ppbv	96
28) Methyl tert-Butyl Ether	5.04	73	489928	2.134	ppbv	100
29) Chloroform	5.76	83	629477	2.127	ppbv	98
30) Cyclohexane	7.14	84	342026	2.140	ppbv	88
31) cis-1,2-Dichloroethene	5.55	61	367755	2.082	ppbv	100
32) 1,1,1-Trichloroethane	6.52	97	649989	2.392	ppbv	100
34) 2-Butanone	5.25	43	412895m	2.368	ppbv	
35) Carbon Tetrachloride	7.03	117	673975	2.362	ppbv	94
36) Benzene	6.90	78	797796	2.166	ppbv	98
37) 1,2-Dichloroethane	6.31	62	476048m	2.225	ppbv	
38) Trichloroethene	7.84	130	356258m	2.238	ppbv	
39) 1,2-Dichloropropane	7.62	63	297822	2.117	ppbv	96
40) 1,4-Dioxane	7.86	88	78705	1.998	ppbv	65
41) Tetrahydrofuran	6.07	42	176113	2.075	ppbv	96
42) Bromodichloromethane	7.80	83	646144	2.192	ppbv	99
43) Methyl Methacrylate	8.02	69	276657	2.189	ppbv	97

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL061621\
 Data File : VL037165.D
 Acq On : 16 Jun 2021 10:04
 Operator : SY/AP
 Sample : VSTDIC002
 Misc : 400mL/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC002

Manual Integrations
 APPROVED

MMDadoda
 6/17/2021 5:37:15 PM

Quant Time: Jun 16 12:20:31 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL061621AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LWed Jun 16 2021
 QLast Update : Wed Jun 16 11:33:05 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,2,4-Trimethylpentane	7.87	57	1392244	2.258	ppbv	99
45) t-1,3-Dichloropropene	9.28	75	241512m	2.121	ppbv	
46) cis-1,3-Dichloropropene	8.71	75	333300	2.193	ppbv	95
47) 1,1,2-Trichloroethane	9.48	97	330404	2.071	ppbv	98
48) Dibromochloromethane	10.32	129	545693	2.120	ppbv	96
49) Bromoform	13.05	173	417351	2.076	ppbv	100
50) 4-Methyl-2-Pentanone	8.76	43	540451	2.107	ppbv	98
51) 2-Hexanone	10.16	43	271527	2.121	ppbv #	99
52) Tetrachloroethene	11.23	164	341343	2.240	ppbv	95
53) Toluene	9.82	91	979146	2.236	ppbv	98
54) 1,2-Dibromoethane	10.62	107	504195	2.158	ppbv	100
56) 1,1,1,2-Tetrachloroethane	12.14	131	406214m	2.094	ppbv	
57) Chlorobenzene	12.15	112	749452	2.103	ppbv	100
58) Ethyl Benzene	12.72	91	1319861	2.191	ppbv	99
59) m/p-Xylene	13.00	91	2212857m	4.322	ppbv	
60) o-Xylene	13.70	91	1065122	2.167	ppbv	99
61) Styrene	13.54	104	539220	2.080	ppbv	98
62) Isopropylbenzene	14.67	105	1502881	2.145	ppbv	98
63) 1,1,2,2-Tetrachloroethane	13.69	83	677532	2.120	ppbv	99
64) n-propylbenzene	15.57	120	375050	2.099	ppbv	82
65) tert-Butylbenzene	16.75	119	1395250	2.197	ppbv	100
66) Benzyl Chloride	17.00	91	57652m	2.080	ppbv	
67) sec-Butylbenzene	17.30	105	1850998	2.194	ppbv	98
69) p-Isopropyltoluene	17.66	119	1584845	2.247	ppbv	98
70) n-Butylbenzene	18.56	91	1458294	2.260	ppbv	98
71) 2-Chlorotoluene	15.46	91	1052843	2.106	ppbv	100
72) 4-Ethyltoluene	15.85	105	1205907	2.112	ppbv	99
73) 1,3,5-Trimethylbenzene	16.00	105	1135065	2.095	ppbv	94
74) 1,2,4-Trimethylbenzene	16.77	105	1208673	2.193	ppbv	95
75) 1,3-Dichlorobenzene	17.00	146	717761	2.115	ppbv	99
76) 1,4-Dichlorobenzene	17.15	146	742835m	2.204	ppbv	
77) 1,2-Dichlorobenzene	17.82	146	698267	2.087	ppbv	99
78) Hexachloro-1,3-Butadiene	23.38	225	614123	2.589	ppbv	98
79) Naphthalene	22.21	128	1065846	3.213	ppbv	99
80) Naphthalene, 2-methyl-	24.99	142	238299	2.969	ppbv #	83
81) 1,2,4-Trichlorobenzene	21.94	180	598560	2.836	ppbv	100

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

