

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL021021\
 Data File : VL036317.D
 Acq On : 10 Feb 2021 13:24
 Operator : SY/AP
 Sample : VSTDIC0.5
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC0.5

Manual Integrations
 APPROVED

MMDadoda
 2/12/2021 1:41:57 PM

Quant Time: Feb 10 16:02:14 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL021021AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LWed Feb 10
 QLast Update : Wed Feb 10 13:51:24 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.65	49	1529841	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.15	114	3444037	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.06	117	3121658	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.39	95	2589043	10.02	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	100.20%

Target Compounds

						Ovalue
2) Dichlorodifluoromethane	3.03	85	146039	0.712	ppbv	94
3) Chlorodifluoromethane	2.97	51	107403	0.572	ppbv	100
4) Chloromethane	3.12	50	50118m	0.540	ppbv	
5) Vinyl Chloride	3.24	62	46241m	0.526	ppbv	
6) Bromomethane	3.46	94	28636m	0.583	ppbv	
7) Chloroethane	3.54	64	14536m	0.480	ppbv	
8) Dichlorotetrafluoroethane	3.17	85	110880	0.558	ppbv	93
9) Propene	3.28	41	15353m	0.172	ppbv	
10) Heptane	8.10	43	102050	0.460	ppbv	97
11) Trichlorofluoromethane	3.93	101	122350	0.587	ppbv	84
12) 1,1,2-Trichlorotrifluoroet	4.46	101	78347	0.620	ppbv	99
13) Ethanol	3.60	45	7187m	0.512	ppbv	
14) Bromoethene	3.72	108	31385	0.545	ppbv #	96
15) Acetone	3.84	43	104995	0.560	ppbv #	86
16) 1,3-Butadiene	3.31	39	44639m	0.489	ppbv	
17) tert-Butyl alcohol	4.29	59	85642m	0.554	ppbv	
18) 1,1-Dichloroethene	4.27	96	33950m	0.579	ppbv	
19) Isopropyl Alcohol	3.96	45	65978m	0.584	ppbv	
20) Methylene Chloride	4.32	84	36499m	0.610	ppbv	
21) Allyl Chloride	4.39	41	54597m	0.535	ppbv	
22) trans-1,2-Dichloroethene	4.86	96	32258m	0.596	ppbv	
23) Vinyl Acetate	5.05	43	115620m	0.509	ppbv	
24) 1,1-Dichloroethane	4.99	63	92806m	0.564	ppbv	
25) Ethyl Acetate	5.66	43	196306m	0.522	ppbv	
26) Hexane	5.66	57	80795m	0.516	ppbv	
27) Carbon Disulfide	4.53	76	62734m	0.501	ppbv	
28) Methyl tert-Butyl Ether	5.02	73	101006m	0.477	ppbv	
29) Chloroform	5.73	83	131131m	0.585	ppbv	
30) Cyclohexane	7.11	84	65265m	0.545	ppbv	
31) cis-1,2-Dichloroethene	5.53	61	77809m	0.520	ppbv	
32) 1,1,1-Trichloroethane	6.49	97	121238m	0.559	ppbv	
34) 2-Butanone	5.23	43	123644m	0.488	ppbv	
35) Carbon Tetrachloride	6.99	117	114341m	0.525	ppbv	
36) Benzene	6.87	78	147327	0.467	ppbv	96
37) 1,2-Dichloroethane	6.28	62	95456	0.503	ppbv	99
38) Trichloroethene	7.81	130	54839m	0.469	ppbv	
39) 1,2-Dichloropropane	7.59	63	59913	0.468	ppbv	94
40) 1,4-Dioxane	7.84	88	25182m	0.543	ppbv	
41) Tetrahydrofuran	6.05	42	60220m	0.411	ppbv	
42) Bromodichloromethane	7.77	83	125498m	0.526	ppbv	
43) Methyl Methacrylate	8.00	69	58768m	0.438	ppbv	

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL021021\
 Data File : VL036317.D
 Acq On : 10 Feb 2021 13:24
 Operator : SY/AP
 Sample : VSTDIC0.5
 Misc : 400ml/MSVOA L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC0.5

Manual Integrations
 APPROVED

MMDadoda
 2/12/2021 1:41:57 PM

Quant Time: Feb 10 16:02:14 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL021021AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LWed Feb 10
 QLast Update : Wed Feb 10 13:51:24 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,2,4-Trimethylpentane	7.84	57	284371	0.491	ppbv	98
45) t-1,3-Dichloropropene	9.25	75	46861	0.368	ppbv	97
46) cis-1,3-Dichloropropene	8.68	75	68393m	0.418	ppbv	
47) 1,1,2-Trichloroethane	9.45	97	66334m	0.541	ppbv	
48) Dibromochloromethane	10.28	129	86181m	0.481	ppbv	
49) Bromoform	13.01	173	65736m	0.439	ppbv	
50) 4-Methyl-2-Pentanone	8.74	43	168576m	0.446	ppbv	
51) 2-Hexanone	10.13	43	112806m	0.378	ppbv	
52) Tetrachloroethene	11.20	164	52479m	0.451	ppbv	
53) Toluene	9.79	91	152164m	0.431	ppbv	
54) 1,2-Dibromoethane	10.59	107	91949m	0.506	ppbv	
56) 1,1,1,2-Tetrachloroethane	12.10	131	78252	0.562	ppbv #	1
57) Chlorobenzene	12.12	112	138205	0.564	ppbv	99
58) Ethyl Benzene	12.69	91	209557	0.452	ppbv	93
59) m/p-Xylene	12.97	91	390724m	0.982	ppbv	
60) o-Xylene	13.67	91	190059	0.495	ppbv	97
61) Styrene	13.51	104	101266m	0.418	ppbv	
62) Isopropylbenzene	14.64	105	289629	0.508	ppbv	100
63) 1,1,2,2-Tetrachloroethane	13.66	83	142036	0.497	ppbv	93
64) n-propylbenzene	15.54	120	67727	0.483	ppbv	76
65) tert-Butylbenzene	16.73	119	237829	0.492	ppbv	92
66) Benzyl Chloride	16.97	91	49161m	0.398	ppbv	
67) sec-Butylbenzene	17.27	105	357681m	0.498	ppbv	
69) p-Isopropyltoluene	17.64	119	255646	0.449	ppbv	97
70) n-Butylbenzene	18.54	91	274232m	0.452	ppbv	
71) 2-Chlorotoluene	15.43	91	213879	0.477	ppbv	97
72) 4-Ethyltoluene	15.82	105	194787	0.450	ppbv	98
73) 1,3,5-Trimethylbenzene	15.98	105	199566m	0.484	ppbv	
74) 1,2,4-Trimethylbenzene	16.74	105	202878	0.479	ppbv #	90
75) 1,3-Dichlorobenzene	16.98	146	120874m	0.532	ppbv	
76) 1,4-Dichlorobenzene	17.12	146	122605m	0.547	ppbv	
77) 1,2-Dichlorobenzene	17.80	146	111722m	0.504	ppbv	
78) Hexachloro-1,3-Butadiene	23.36	225	104069m	0.587	ppbv	
79) Naphthalene	22.19	128	91820m	0.371	ppbv	
80) Naphthalene, 2-methyl-	24.99	142	20545m	0.389	ppbv	
81) 1,2,4-Trichlorobenzene	21.91	180	70201m	0.486	ppbv	

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

