

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL070219\
 Data File : VL033992.D
 Acq On : 2 Jul 2019 15:36
 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
 7/3/2019 11:03:31 PM

Quant Time: Jul 02 19:33:31 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL070219AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Tue Jul 02 2019
 QLast Update : Tue Jul 02 16:27:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.73	49	1094816	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.26	114	2894061	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.20	117	2742008	10.00	ppbv	0.00

System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	14.54	95	2227084	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	3.06	85	146969	0.827	ppbv	91
3) Chlorodifluoromethane	3.00	51	78121	0.606	ppbv	100
4) Chloromethane	3.16	50	44966	0.612	ppbv #	88
5) Vinyl Chloride	3.27	62	42845	0.628	ppbv	93
6) Bromomethane	3.50	94	26335	0.581	ppbv	82
7) Chloroethane	3.58	64	16437	0.609	ppbv	94
8) Dichlorotetrafluoroethane	3.20	85	122206	0.644	ppbv	100
9) Propene	3.02	41	26658m	0.553	ppbv	
10) Heptane	8.22	43	53657m	0.440	ppbv	
11) Trichlorofluoromethane	3.98	101	132975	0.580	ppbv	89
12) 1,1,2-Trichlorotrifluoroet	4.53	101	90152	0.582	ppbv	97
13) Ethanol	3.64	45	8474m	0.540	ppbv	
14) Bromoethene	3.77	108	32443	0.562	ppbv #	91
15) Acetone	3.89	43	101547	0.636	ppbv	95
16) 1,3-Butadiene	3.35	39	32185	0.530	ppbv	95
17) tert-Butyl alcohol	4.35	59	66363	0.532	ppbv #	92
18) 1,1-Dichloroethene	4.33	96	39959	0.611	ppbv	85
19) Isopropyl Alcohol	4.03	45	65160m	0.714	ppbv	
20) Methylene Chloride	4.38	84	44896m	0.721	ppbv	
21) Allyl Chloride	4.45	41	52446	0.578	ppbv	95
22) trans-1,2-Dichloroethene	4.94	96	39622	0.574	ppbv	88
23) Vinyl Acetate	5.13	43	72098	0.388	ppbv #	91
24) 1,1-Dichloroethane	5.06	63	85396	0.592	ppbv #	96
25) Ethyl Acetate	5.75	43	119831	0.493	ppbv	98
26) Hexane	5.75	57	52808	0.477	ppbv	90
27) Carbon Disulfide	4.60	76	86736	0.564	ppbv #	83
28) Methyl tert-Butyl Ether	5.10	73	88813	0.530	ppbv	100
29) Chloroform	5.82	83	116159	0.613	ppbv	88
30) Cyclohexane	7.22	84	38920m	0.498	ppbv	
31) cis-1,2-Dichloroethene	5.61	61	50313m	0.518	ppbv	
32) 1,1,1-Trichloroethane	6.59	97	101699	0.556	ppbv	98
34) 2-Butanone	5.31	43	66803	0.409	ppbv	96
35) Carbon Tetrachloride	7.10	117	109865m	0.491	ppbv	
36) Benzene	6.97	78	99893	0.452	ppbv #	92
37) 1,2-Dichloroethane	6.38	62	83500	0.511	ppbv #	97
38) Trichloroethene	7.93	130	45952m	0.534	ppbv	
39) 1,2-Dichloropropane	7.71	63	37757	0.443	ppbv #	88
40) 1,4-Dioxane	7.97	88	20103m	0.578	ppbv	
41) Tetrahydrofuran	6.14	42	29489m	0.356	ppbv	
42) Bromodichloromethane	7.89	83	107154	0.490	ppbv	90
43) Methyl Methacrylate	8.11	69	34033m	0.412	ppbv	

Data Path : Z:\VOASRV\HPCHEM1\MSVOA L\DATA\VL070219\
 Data File : VL033992.D
 Acq On : 2 Jul 2019 15:36
 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
 7/3/2019 11:03:31 PM

Quant Time: Jul 02 19:33:31 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA L\METHODS\VL070219AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Tue Jul 0
 QLast Update : Tue Jul 02 16:27:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

44) 2,2,4-Trimethylpentane	7.96	57	169305m	0.431	ppbv	
45) t-1,3-Dichloropropene	9.38	75	39573m	0.365	ppbv	
46) cis-1,3-Dichloropropene	8.80	75	48215m	0.388	ppbv	
47) 1,1,2-Trichloroethane	9.58	97	51106m	0.502	ppbv	
48) Dibromochloromethane	10.42	129	117419	0.642	ppbv	97
49) Bromoform	13.16	173	73870m	0.458	ppbv	
50) 4-Methyl-2-Pentanone	8.86	43	89690m	0.368	ppbv	
51) 2-Hexanone	10.26	43	85628m	0.450	ppbv	
52) Tetrachloroethene	11.34	164	44860m	0.577	ppbv	
53) Toluene	9.91	91	91753	0.375	ppbv	94
54) 1,2-Dibromoethane	10.72	107	66286	0.432	ppbv	86
56) 1,1,1,2-Tetrachloroethane	12.25	131	62389	0.521	ppbv #	1
57) Chlorobenzene	12.26	112	114132m	0.556	ppbv	
58) Ethyl Benzene	12.83	91	124140	0.402	ppbv	90
59) m/p-Xylene	13.11	91	256564m	0.913	ppbv	
60) o-Xylene	13.81	91	127079	0.438	ppbv	97
61) Styrene	13.65	104	71032m	0.378	ppbv	
62) Isopropylbenzene	14.79	105	183018m	0.468	ppbv	
63) 1,1,2,2-Tetrachloroethane	13.79	83	116208m	0.532	ppbv	
64) n-propylbenzene	15.69	120	45865m	0.458	ppbv	
65) tert-Butylbenzene	16.88	119	158635	0.438	ppbv	95
66) Benzyl Chloride	17.11	91	80166m	0.437	ppbv	
67) sec-Butylbenzene	17.42	105	226159	0.445	ppbv	99
69) p-Isopropyltoluene	17.78	119	181934	0.431	ppbv	98
70) n-Butylbenzene	18.68	91	195721m	0.423	ppbv	
71) 2-Chlorotoluene	15.58	91	119999m	0.416	ppbv	
72) 4-Ethyltoluene	15.96	105	134045	0.406	ppbv #	95
73) 1,3,5-Trimethylbenzene	16.13	105	137920	0.424	ppbv	87
74) 1,2,4-Trimethylbenzene	16.89	105	171736	0.458	ppbv #	82
75) 1,3-Dichlorobenzene	17.13	146	125568m	0.535	ppbv	
76) 1,4-Dichlorobenzene	17.27	146	112562m	0.522	ppbv	
77) 1,2-Dichlorobenzene	17.95	146	108510	0.501	ppbv	99
78) Hexachloro-1,3-Butadiene	23.50	225	106035m	0.626	ppbv	
79) Naphthalene	22.38	128	119870m	0.400	ppbv	
80) Naphthalene, 2-methyl-	25.08	142	40685	0.340	ppbv	95
81) 1,2,4-Trichlorobenzene	22.10	180	90654m	0.479	ppbv	

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

