

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_L\DATA\VL120718\
 Data File : VL032921.D
 Acq On : 7 Dec 2018 4:19
 Operator : SY/AP
 Sample : VSTDICC002
 Misc : 400ml/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC002

Manual Integrations
 APPROVED

MMDadoda
 12/11/2018 10:07:59 AM

Quant Time: Dec 07 06:01:23 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_L\METHODS\VL120718AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Dec 07 2018
 QLast Update : Fri Dec 07 05:41:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	5.77	49	1188175	10.00	ppbv	0.00
33) 1,4-Difluorobenzene	7.30	114	3221788	10.00	ppbv	0.00
55) Chlorobenzene-d5	12.25	117	3117268	10.00	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	14.59	95	2282888	10.03	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	100.30%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	3.08	85	359006	2.204	ppbv	99
3) Chlorodifluoromethane	3.02	51	197081	2.126	ppbv	97
4) Chloromethane	3.17	50	111524	2.041	ppbv	93
5) Vinyl Chloride	3.29	62	106403	1.988	ppbv	92
6) Bromomethane	3.52	94	71500	2.061	ppbv	92
7) Chloroethane	3.61	64	39894	1.924	ppbv #	87
8) Dichlorotetrafluoroethane	3.22	85	284726	2.147	ppbv	99
9) Propene	3.04	41	81396	2.194	ppbv	97
10) Heptane	8.25	43	296768	2.103	ppbv	99
11) Trichlorofluoromethane	4.01	101	319759	1.982	ppbv	98
12) 1,1,2-Trichlorotrifluoroet	4.56	101	276424	2.254	ppbv	98
13) Ethanol	3.66	45	30598	2.436	ppbv	98
14) Bromoethene	3.79	108	87587	1.951	ppbv	99
15) Acetone	3.91	43	232901	2.196	ppbv	97
16) 1,3-Butadiene	3.36	39	93858	2.217	ppbv	94
17) tert-Butyl alcohol	4.38	59	48059	1.632	ppbv #	82
18) 1,1-Dichloroethene	4.36	96	96309	1.808	ppbv	94
19) Isopropyl Alcohol	4.04	45	134499	1.753	ppbv #	97
20) Methylene Chloride	4.41	84	92363	1.550	ppbv	97
21) Allyl Chloride	4.48	41	172478	2.205	ppbv	98
22) trans-1,2-Dichloroethene	4.97	96	142081	2.208	ppbv	94
23) Vinyl Acetate	5.16	43	384999	1.926	ppbv	97
24) 1,1-Dichloroethane	5.09	63	319475	2.077	ppbv	99
25) Ethyl Acetate	5.78	43	498187	1.991	ppbv	99
26) Hexane	5.78	57	229484	1.953	ppbv	98
27) Carbon Disulfide	4.62	76	284979	1.918	ppbv	98
28) Methyl tert-Butyl Ether	5.12	73	378898	2.013	ppbv	98
29) Chloroform	5.85	83	414586	2.125	ppbv	99
30) Cyclohexane	7.25	84	199368	2.078	ppbv #	82
31) cis-1,2-Dichloroethene	5.64	61	224268	2.098	ppbv	98
32) 1,1,1-Trichloroethane	6.62	97	439733	2.013	ppbv	99
34) 2-Butanone	5.34	43	299639	1.879	ppbv	100
35) Carbon Tetrachloride	7.14	117	480966	2.007	ppbv	94
36) Benzene	7.01	78	492085	1.950	ppbv	97
37) 1,2-Dichloroethane	6.42	62	327670	2.019	ppbv	100
38) Trichloroethene	7.96	130	195446	1.944	ppbv	98
39) 1,2-Dichloropropane	7.74	63	182253	1.970	ppbv	92
40) 1,4-Dioxane	7.98	88	70608m	2.021	ppbv	
41) Tetrahydrofuran	6.16	42	184288	2.029	ppbv	98
42) Bromodichloromethane	7.92	83	455700	2.019	ppbv	99
43) Methyl Methacrylate	8.15	69	180080	1.930	ppbv	97

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_L\DATA\VL120718\
 Data File : VL032921.D
 Acq On : 7 Dec 2018 4:19
 Operator : SY/AP
 Sample : VSTDIC002
 Misc : 400ml/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC002

Manual Integrations
 APPROVED

MMDadoda
 12/11/2018 10:07:59 AM

Quant Time: Dec 07 06:01:23 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_L\METHODS\VL120718AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Dec 07 06:01:23 2018
 QLast Update : Fri Dec 07 05:41:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,2,4-Trimethylpentane	7.99	57	855744	2.078	ppbv	100
45) t-1,3-Dichloropropene	9.43	75	206884	1.688	ppbv	92
46) cis-1,3-Dichloropropene	8.84	75	280910	2.022	ppbv	96
47) 1,1,2-Trichloroethane	9.62	97	203668	1.863	ppbv	97
48) Dibromochloromethane	10.45	129	376978	1.866	ppbv	99
49) Bromoform	13.20	173	308760	1.821	ppbv	98
50) 4-Methyl-2-Pentanone	8.88	43	495047	2.122	ppbv	99
51) 2-Hexanone	10.29	43	319506	1.852	ppbv #	98
52) Tetrachloroethene	11.38	164	171978	1.821	ppbv	94
53) Toluene	9.95	91	528208	1.847	ppbv	100
54) 1,2-Dibromoethane	10.76	107	322367	1.977	ppbv	99
56) 1,1,1,2-Tetrachloroethane	12.28	131	282517m	2.006	ppbv	
57) Chlorobenzene	12.31	112	448771	1.928	ppbv	96
58) Ethyl Benzene	12.87	91	726521	2.045	ppbv	99
59) m/p-Xylene	13.13	91	1301459m	4.058	ppbv	
60) o-Xylene	13.86	91	691111	2.206	ppbv	99
61) Styrene	13.69	104	439815	2.102	ppbv	98
62) Isopropylbenzene	14.84	105	909612	2.046	ppbv	99
63) 1,1,2,2-Tetrachloroethane	13.84	83	494683	2.094	ppbv	100
64) n-propylbenzene	15.72	120	234352	2.091	ppbv	95
65) tert-Butylbenzene	16.92	119	819167	2.148	ppbv	97
66) Benzyl Chloride	17.17	91	451544m	2.233	ppbv	
67) sec-Butylbenzene	17.47	105	1261628	2.333	ppbv	100
69) p-Isopropyltoluene	17.82	119	1052632	2.449	ppbv	98
70) n-Butylbenzene	18.72	91	956081	2.261	ppbv	98
71) 2-Chlorotoluene	15.62	91	686891	2.139	ppbv	99
72) 4-Ethyltoluene	16.01	105	767946	2.323	ppbv	98
73) 1,3,5-Trimethylbenzene	16.16	105	752943	2.364	ppbv	97
74) 1,2,4-Trimethylbenzene	16.93	105	778681	2.248	ppbv	97
75) 1,3-Dichlorobenzene	17.17	146	478606	2.359	ppbv	100
76) 1,4-Dichlorobenzene	17.32	146	434587	2.287	ppbv	98
77) 1,2-Dichlorobenzene	18.00	146	445708	2.330	ppbv	99
78) Hexachloro-1,3-Butadiene	23.55	225	239839	1.985	ppbv	99
79) Naphthalene	22.42	128	587547	2.869	ppbv	99
80) Naphthalene,2-methyl-	25.13	142	102391	2.237	ppbv	98
81) 1,2,4-Trichlorobenzene	22.15	180	285551	2.644	ppbv	98

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

