

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\
 Data File : VV017608.D
 Acq On : 23 Jul 2020 17:26
 Operator : SY/MD
 Sample : MDL03
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 MDL03

Manual Integrations
 APPROVED

sam
 7/24/2020 7:04:48 PM

Quant Time: Jul 24 11:50:16 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR072320WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Fri Jul 24 02:08:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.64	114	188128	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.87	117	181071	5.00	ug/L	0.00
61) 1,4-Dichlorobenzene-d4	11.27	152	89469	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.31	65	46492	5.18	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	103.60%
7) Chloroethane-d5	1.58	69	36156	5.26	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	105.20%
11) 1,1-Dichloroethene-d2	2.12	63	76377	3.79	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	75.80%
20) 2-Butanone-d5	3.95	46	106826	49.71	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	99.42%
24) Chloroform-d	4.38	84	105974	4.87	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	97.40%
26) 1,2-Dichloroethane-d4	5.06	65	58644	5.23	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	104.60%
32) Benzene-d6	5.07	84	198917	4.98	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	99.60%
36) 1,2-Dichloropropane-d6	6.09	67	56757	5.11	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	102.20%
41) Toluene-d8	7.34	98	185615	4.85	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	97.00%
45) trans-1,3-Dichloropropene-	7.65	79	24204	4.81	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	96.20%
48) 2-Hexanone-d5	8.11	63	73080	46.90	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	93.80%
59) 1,1,2,2-Tetrachloroethane-	10.24	84	38211	4.95	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	99.00%
65) 1,2-Dichlorobenzene-d4	11.65	152	71441	5.28	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	105.60%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.14	85	5552	0.269 ug/L	95
3) Chloromethane	1.25	50	4135	0.308 ug/L	94
5) Vinyl chloride	1.32	62	4134	0.308 ug/L #	77
6) Bromomethane	1.53	94	2649	0.307 ug/L	96
8) Chloroethane	1.60	64	2508	0.323 ug/L	95
9) Trichlorofluoromethane	1.76	101	6740	0.280 ug/L	94
10) 1,1,2-Trichloro-1,2,2-trif	2.13	101	3986	0.355 ug/L	90
12) 1,1-Dichloroethene	2.14	96	3505	0.346 ug/L #	1
13) Acetone	2.23	43	6516m	3.873 ug/L	
14) Carbon disulfide	2.31	76	8746	0.284 ug/L	99
15) Methyl Acetate	2.46	43	1082m	0.331 ug/L	
16) Methylene chloride	2.52	84	5115	0.454 ug/L	80
17) Methyl tert-butyl Ether	2.79	73	7619	0.293 ug/L #	85
18) trans-1,2-Dichloroethene	2.78	96	3315	0.300 ug/L	95
19) 1,1-Dichloroethane	3.21	63	6463	0.278 ug/L	97
21) 2-Butanone	4.05	43	7462m	2.627 ug/L	
22) cis-1,2-Dichloroethene	3.93	96	3790	0.273 ug/L	92

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

23) Bromochloromethane	4.28	128	1841	0.287	ug/L #	95
25) Chloroform	4.40	83	8222	0.316	ug/L	92
27) 1,2-Dichloroethane	5.16	62	5111	0.303	ug/L	97
29) 1,1,1-Trichloroethane	4.64	97	6943	0.274	ug/L	94
30) Cyclohexane	4.69	56	4838	0.241	ug/L #	83
31) Carbon tetrachloride	4.85	117	5987	0.266	ug/L	94
33) Benzene	5.13	78	13710	0.267	ug/L	100
34) Trichloroethene	5.94	95	4151	0.290	ug/L	93
35) Methylcyclohexane	6.15	83	4978	0.230	ug/L #	86
37) 1,2-Dichloropropane	6.20	63	3159	0.262	ug/L #	92
38) Bromodichloromethane	6.54	83	4748	0.268	ug/L #	94
39) cis-1,3-Dichloropropene	7.05	75	5056	0.278	ug/L	93
40) 4-Methyl-2-pentanone	7.26	43	16408	2.484	ug/L #	87
42) Toluene	7.41	91	14486	0.259	ug/L	92
43) 1,3,5-Trimethylbenzene	10.56	105	11031	0.211	ug/L	98
44) 1,2,4-Trimethylbenzene	10.94	105	10613	0.199	ug/L	97
46) trans-1,3-Dichloropropene	7.68	75	4361	0.267	ug/L	93
47) 1,1,2-Trichloroethane	7.87	97	2554	0.280	ug/L	93
49) Tetrachloroethene	8.00	164	3449	0.273	ug/L	87
50) 2-Hexanone	8.17	43	16417	3.513	ug/L	98
51) Dibromochloromethane	8.27	129	3363	0.289	ug/L	99
52) 1,2-Dibromoethane	8.38	107	2509	0.291	ug/L #	97
53) Chlorobenzene	8.90	112	10307	0.277	ug/L	90
54) Ethylbenzene	9.04	91	15133	0.242	ug/L	96
55) m,p-xylene	9.16	106	6153	0.261	ug/L	96
56) o-xylene	9.57	106	5543	0.242	ug/L	87
57) Styrene	9.59	104	9253	0.242	ug/L	81
58) Isopropylbenzene	9.95	105	14935	0.237	ug/L	98
60) 1,1,2,2-Tetrachloroethane	10.26	83	2719	0.273	ug/L #	97
62) Bromoform	9.76	173	1655	0.289	ug/L #	95
63) 1,3-Dichlorobenzene	11.21	146	7795	0.271	ug/L	95
64) 1,4-Dichlorobenzene	11.30	146	8809	0.305	ug/L	96
66) 1,2-Dichlorobenzene	11.67	146	7986	0.304	ug/L	94
67) 1,2-Dibromo-3-chloropropan	12.45	75	467	0.285	ug/L #	80
68) 1,3,5-Trichlorobenzene	12.67	180	6393	0.273	ug/L	97
69) 1,2,4-trichlorobenzene	13.29	180	5532	0.283	ug/L	94
70) Naphthalene	13.53	128	6347	0.233	ug/L	97
71) 1,2,3-Trichlorobenzene	13.77	180	4452	0.257	ug/L	98

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

