

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV091020\
 Data File : VV018189.D
 Acq On : 10 Sep 2020 19:19
 Operator : SY/MD
 Sample : MDL07
 Misc : 5.0mL/MSVOA V/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 MDL07

Manual Integrations
 APPROVED

MMDadoda
 9/11/2020 10:29:17 AM

Quant Time: Sep 11 03:26:52 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM090920WMA.M
 Quant Title : VOC Analysis
 QLast Update : Fri Sep 11 01:30:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.64	114	286518	50.00	ug/L	0.00
28) Chlorobenzene-d5	8.87	117	285863	50.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.27	152	144311	50.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	96292	41.24	ug/L	0.00
Spiked Amount	50.000	Range	60 - 135	Recovery	=	82.48%
7) Chloroethane-d5	1.58	69	84510	44.23	ug/L	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	88.46%
11) 1,1-Dichloroethene-d2	2.12	63	147331	34.40	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	68.80%
21) 2-Butanone-d5	3.91	46	142838	99.94	ug/L	0.00
Spiked Amount	100.000	Range	40 - 130	Recovery	=	99.94%
24) Chloroform-d	4.38	84	182710	45.10	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	90.20%
26) 1,2-Dichloroethane-d4	5.06	65	127349	49.04	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	98.08%
32) Benzene-d6	5.08	84	379919	46.42	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	92.84%
36) 1,2-Dichloropropane-d6	6.10	67	124545	47.90	ug/L	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.80%
41) Toluene-d8	7.34	98	343204	46.02	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	92.04%
43) trans-1,3-Dichloropropene-	7.64	79	59648	45.14	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	90.28%
47) 2-Hexanone-d5	8.11	63	89468	90.84	ug/L	0.00
Spiked Amount	100.000	Range	45 - 130	Recovery	=	90.84%
57) 1,1,2,2-Tetrachloroethane-	10.24	84	151178	46.32	ug/L	0.00
Spiked Amount	50.000	Range	65 - 120	Recovery	=	92.64%
64) 1,2-Dichlorobenzene-d4	11.65	152	147308	49.70	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	99.40%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	1.14	85	4973	2.380	ug/L	95
3) Chloromethane	1.25	50	5705	2.501	ug/L	91
5) Vinyl chloride	1.32	62	5944	2.541	ug/L #	64
6) Bromomethane	1.53	94	3480	2.434	ug/L	99
8) Chloroethane	1.60	64	3597	2.454	ug/L	99
9) Trichlorofluoromethane	1.77	101	7983	2.480	ug/L	98
10) 1,1,2-Trichloro-1,2,2-trif	2.13	101	6238	3.226	ug/L #	82
12) 1,1-Dichloroethene	2.13	96	5890	3.131	ug/L #	1
13) Acetone	2.20	43	5685	4.967	ug/L	99
14) Carbon disulfide	2.31	76	14057	2.442	ug/L	97
15) Methyl Acetate	2.46	43	4952	2.408	ug/L	94
16) Methylene chloride	2.53	84	9277	4.397	ug/L	97
17) trans-1,2-Dichloroethene	2.78	96	4491	2.316	ug/L	99
18) Methyl tert-butyl Ether	2.79	73	13773	2.296	ug/L	99
19) 1,1-Dichloroethane	3.22	63	9058	2.471	ug/L	97
20) cis-1,2-Dichloroethene	3.95	96	5166	2.493	ug/L	96
22) 2-Butanone	4.02	43	6059m	4.471	ug/L	

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

23) Bromochloromethane	4.29	128	2628m	2.424	uq/L	
25) Chloroform	4.41	83	17319	4.648	uq/L	99
27) 1,2-Dichloroethane	5.17	62	7381	2.559	uq/L	100
29) Cyclohexane	4.70	56	7208	2.184	uq/L	95
30) 1,1,1-Trichloroethane	4.64	97	7467	2.251	uq/L	96
31) Carbon tetrachloride	4.86	117	6710	2.300	uq/L	96
33) Benzene	5.13	78	19646	2.404	uq/L	100
34) Trichloroethene	5.95	95	5649	2.537	uq/L	98
35) Methylcyclohexane	6.16	83	7694	2.295	uq/L	98
37) 1,2-Dichloropropane	6.21	63	4782	2.208	uq/L #	91
38) Bromodichloromethane	6.54	83	6686	2.372	uq/L	89
39) cis-1,3-Dichloropropene	7.05	75	6902	2.068	uq/L	96
40) 4-Methyl-2-pentanone	7.25	43	10969	4.175	uq/L	99
42) Toluene	7.41	91	20514	2.372	uq/L	94
44) trans-1,3-Dichloropropene	7.68	75	7798	2.400	uq/L	96
45) 1,1,2-Trichloroethane	7.86	97	4931	2.456	uq/L	93
46) Tetrachloroethene	8.00	164	4103	2.350	uq/L	95
48) 2-Hexanone	8.16	43	16095	7.770	uq/L	95
49) Dibromochloromethane	8.27	129	5037	2.235	uq/L	99
50) 1,2-Dibromoethane	8.38	107	4772	2.259	uq/L	97
51) Chlorobenzene	8.90	112	13676	2.435	uq/L	96
52) Ethylbenzene	9.04	91	20701	2.192	uq/L	99
53) m,p-Xylene	9.16	106	7697	2.184	uq/L	95
54) o-xylene	9.57	106	7188	2.117	uq/L	97
55) Styrene	9.59	104	11227	1.902	uq/L	99
56) Isopropylbenzene	9.95	105	18339	2.004	uq/L	98
58) 1,1,2,2-Tetrachloroethane	10.26	83	7591	2.556	uq/L #	97
59) 1,2,3-Trichloropropane	10.30	75	6205	2.436	uq/L	98
61) Bromoform	9.76	173	3209	2.070	uq/L	96
62) 1,3-Dichlorobenzene	11.21	146	10525	2.453	uq/L	99
63) 1,4-Dichlorobenzene	11.30	146	11446	2.595	uq/L	99
65) 1,2-Dichlorobenzene	11.67	146	11375	2.623	uq/L #	90
66) 1,2-Dibromo-3-chloropropan	12.45	75	1244	2.032	uq/L	92
67) 1,3,5-Trichlorobenzene	12.67	180	7977	2.358	uq/L	99
68) 1,2,4-trichlorobenzene	13.29	180	6735	2.236	uq/L	99
69) Naphthalene	13.53	128	13658	1.792	uq/L	100
70) 1,2,3-Trichlorobenzene	13.77	180	6792	2.289	uq/L	94

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

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