

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU020720\
 Data File : VU036696.D
 Acq On : 07 Feb 2020 12:27
 Operator : JC/MD
 Sample : MDL03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL03

Manual Integrations
 APPROVED

MMDadoda
 2/10/2020 11:54:10 AM

Quant Time: Feb 07 23:16:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SFAMULM020620W.M
 Quant Title : VOC Analysis
 QLast Update : Fri Feb 07 22:36:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	6.29	114	261788	50.00	ug/L	0.00
28) Chlorobenzene-d5	9.45	117	245713	50.00	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.83	152	104680	50.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.61	65	77312	44.28	ug/L	0.00
Spiked Amount	50.000	Range	60 - 135	Recovery	=	88.56%
7) Chloroethane-d5	1.93	69	74395	45.96	ug/L	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	91.92%
11) 1,1-Dichloroethene-d2	2.59	63	112777	34.21	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	68.42%
21) 2-Butanone-d5	4.68	46	113363	76.11	ug/L	0.00
Spiked Amount	100.000	Range	40 - 130	Recovery	=	76.11%
24) Chloroform-d	5.11	84	136041	43.05	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	86.10%
26) 1,2-Dichloroethane-d4	5.75	65	102558	47.45	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	94.90%
32) Benzene-d6	5.77	84	312278	48.34	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	96.68%
36) 1,2-Dichloropropane-d6	6.73	67	104926	48.93	ug/L	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	97.86%
41) Toluene-d8	7.93	98	282359	47.73	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	95.46%
43) trans-1,3-Dichloropropene-	8.21	79	47258	46.02	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	92.04%
47) 2-Hexanone-d5	8.67	63	124851	102.01	ug/L	0.00
Spiked Amount	100.000	Range	45 - 130	Recovery	=	102.01%
56) 1,1,2,2-Tetrachloroethane-	10.78	84	152438	48.24	ug/L	0.00
Spiked Amount	50.000	Range	65 - 120	Recovery	=	96.48%
66) 1,2-Dichlorobenzene-d4	12.21	152	94195	49.53	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	99.06%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.39	85	4664	2.579 ug/L	100
3) Chloromethane	1.54	50	5768	2.691 ug/L	99
5) Vinyl chloride	1.62	62	6450	2.795 ug/L #	63
6) Bromomethane	1.87	94	3645	2.519 ug/L	94
8) Chloroethane	1.95	64	4711	3.216 ug/L	86
9) Trichlorofluoromethane	2.16	101	7078	2.581 ug/L	97
10) 1,1,2-Trichloro-1,2,2-trif	2.60	101	5064	3.083 ug/L #	86
12) 1,1-Dichloroethene	2.61	96	5172	3.234 ug/L #	1
13) Acetone	2.66	43	11816	7.468 ug/L	93
14) Carbon disulfide	2.82	76	11926	2.496 ug/L	99
15) Methyl Acetate	2.99	43	5861	2.562 ug/L	98
16) Methylene chloride	3.08	84	5523	2.798 ug/L	94
17) trans-1,2-Dichloroethene	3.39	96	4486	2.637 ug/L	95
18) Methyl tert-butyl Ether	3.41	73	14915	2.526 ug/L	98
19) 1,1-Dichloroethane	3.92	63	8706	2.568 ug/L	96
20) cis-1,2-Dichloroethene	4.72	96	4899	2.527 ug/L	97
22) 2-Butanone	4.76	43	18075	9.625 ug/L	82

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU020720\
 Data File : VU036696.D
 Acq On : 07 Feb 2020 12:27
 Operator : JC/MD
 Sample : MDL03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL03

Manual Integrations
 APPROVED

MMDadoda
 2/10/2020 11:54:10 AM

Quant Time: Feb 07 23:16:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SFAMULM020620W.M
 Quant Title : VOC Analysis
 QLast Update : Fri Feb 07 22:36:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	5.02	128	2399	2.641	ug/L	91
25) Chloroform	5.14	83	21436	6.501	ug/L	98
27) 1,2-Dichloroethane	5.84	62	7579	2.733	ug/L	97
29) Cyclohexane	5.43	56	8265	2.619	ug/L	98
30) 1,1,1-Trichloroethane	5.36	97	6406	2.454	ug/L #	74
31) Carbon tetrachloride	5.57	117	5297	2.506	ug/L	98
33) Benzene	5.82	78	19855	2.672	ug/L	100
34) Trichloroethene	6.58	95	4829	2.628	ug/L	97
35) Methylcyclohexane	6.80	83	7983	2.467	ug/L	97
37) 1,2-Dichloropropane	6.83	63	5398	2.694	ug/L #	88
38) Bromodichloromethane	7.14	83	6231	2.559	ug/L	99
39) cis-1,3-Dichloropropene	7.64	75	8185	2.556	ug/L	100
40) 4-Methyl-2-pentanone	7.83	43	16493	5.209	ug/L	99
42) Toluene	8.00	91	20923	2.605	ug/L	96
44) trans-1,3-Dichloropropene	8.24	75	7219m	2.581	ug/L	
45) 1,1,2-Trichloroethane	8.43	97	4508	2.466	ug/L	97
46) Tetrachloroethene	8.58	164	3840	2.978	ug/L	92
48) 2-Hexanone	8.72	43	14461	5.599	ug/L	97
49) Dibromochloromethane	8.84	129	4113	2.300	ug/L	86
50) 1,2-Dibromoethane	8.95	107	4495	2.340	ug/L #	98
51) Chlorobenzene	9.47	112	12765	2.640	ug/L	99
52) Ethylbenzene	9.60	91	21684	2.457	ug/L	96
53) m,p-Xylene	9.72	106	8100	2.488	ug/L	93
54) o-xylene	10.13	106	7714	2.392	ug/L	95
55) Styrene	10.14	104	12091	2.250	ug/L	97
57) 1,1,2,2-Tetrachloroethane	10.81	83	8879	2.695	ug/L	99
59) Bromoform	10.32	173	2832	2.449	ug/L #	94
60) 1,2,3-Trichloropropane	10.85	75	6895	2.800	ug/L	99
61) Isopropylbenzene	10.51	105	20328	2.608	ug/L	99
62) 1,3,5-Trimethylbenzene	11.11	105	16896	2.528	ug/L	95
63) 1,2,4-Trimethylbenzene	11.49	105	16440	2.464	ug/L	99
64) 1,3-Dichlorobenzene	11.77	146	8474	2.585	ug/L	98
65) 1,4-Dichlorobenzene	11.86	146	8694	2.634	ug/L	97
67) 1,2-Dichlorobenzene	12.24	146	8965	2.701	ug/L	98
68) 1,2-Dibromo-3-chloropropan	13.02	75	1669	2.372	ug/L	91
69) 1,3,5-Trichlorobenzene	13.24	180	5054	2.223	ug/L	99
70) 1,2,4-trichlorobenzene	13.85	180	3480	1.836	ug/L	95
71) Naphthalene	14.10	128	10769	1.612	ug/L	99
72) 1,2,3-Trichlorobenzene	14.34	180	3452	1.828	ug/L	94

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

