

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV102418\
 Data File : VV008391.D
 Acq On : 24 Oct 2018 14:40
 Operator : SY/MD
 Sample : MDL02
 Misc : 5.0 mL/MSVOA_V/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 MDL02

Manual Integrations
 APPROVED

apatel
 10/25/2018 4:40:40 PM

Quant Time: Oct 25 10:53:26 2018

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM102318WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Oct 25 08:48:14 2018

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.67	114	192427	50.00	ug/L	0.00
28) Chlorobenzene-d5	8.90	117	173434	50.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.30	152	74152	50.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	47702	44.77	ug/L	0.00
Spiked Amount	50.000	Range	60 - 135	Recovery	=	89.54%
7) Chloroethane-d5	1.57	69	37547	48.81	ug/L	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	97.62%
11) 1,1-Dichloroethene-d2	2.13	63	67728	35.20	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	70.40%
21) 2-Butanone-d5	3.93	46	115063	99.37	ug/L	0.00
Spiked Amount	100.000	Range	40 - 130	Recovery	=	99.37%
24) Chloroform-d	4.41	84	121157	46.11	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	92.22%
26) 1,2-Dichloroethane-d4	5.09	65	83671	47.62	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	95.24%
32) Benzene-d6	5.10	84	244688	49.51	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	99.02%
36) 1,2-Dichloropropane-d6	6.12	67	84802	49.43	ug/L	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.86%
41) Toluene-d8	7.36	98	219615	48.94	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	97.88%
43) trans-1,3-Dichloropropene-	7.67	79	41158	49.38	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	98.76%
47) 2-Hexanone-d5	8.14	63	88765	99.09	ug/L	0.00
Spiked Amount	100.000	Range	45 - 130	Recovery	=	99.09%
57) 1,1,2,2-Tetrachloroethane-	10.26	84	114723	49.38	ug/L	0.00
Spiked Amount	50.000	Range	65 - 120	Recovery	=	98.76%
64) 1,2-Dichlorobenzene-d4	11.68	152	74921	49.39	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	98.78%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	1.14	85	2416	1.758 ug/L	96
3) Chloromethane	1.25	50	2748	1.867 ug/L	95
5) Vinyl chloride	1.33	62	2579	1.891 ug/L #	52
6) Bromomethane	1.52	94	853	2.279 ug/L	94
8) Chloroethane	1.59	64	1591	2.220 ug/L	91
9) Trichlorofluoromethane	1.77	101	3086	1.758 ug/L	95
10) 1,1,2-Trichloro-1,2,2-trif	2.14	101	2172	2.369 ug/L #	82
12) 1,1-Dichloroethene	2.14	96	2275	2.635 ug/L #	1
13) Acetone	2.19	43	3595	4.000 ug/L	96
14) Carbon disulfide	2.32	76	7161	1.842 ug/L	100
15) Methyl Acetate	2.46	43	3203	1.787 ug/L	98
16) Methylene chloride	2.54	84	2891	1.933 ug/L	91
17) trans-1,2-Dichloroethene	2.80	96	2381	1.809 ug/L	95
18) Methyl tert-butyl Ether	2.81	73	8390	1.781 ug/L #	90
19) 1,1-Dichloroethane	3.24	63	4814	1.774 ug/L	96
20) cis-1,2-Dichloroethene	3.97	96	2879	1.920 ug/L	96
22) 2-Butanone	4.03	43	4267	3.287 ug/L #	67

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV102418\
 Data File : VV008391.D
 Acq On : 24 Oct 2018 14:40
 Operator : SY/MD
 Sample : MDL02
 Misc : 5.0 mL/MSVOA_V/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 MDL02

Manual Integrations
 APPROVED

apatel
 10/25/2018 4:40:40 PM

Quant Time: Oct 25 10:53:26 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM102318WMA.M
 Quant Title : VOC Analysis
 QLast Update : Thu Oct 25 08:48:14 2018
 Response via : Initial Calibration

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

23) Bromochloromethane	4.31	128	1245m	1.761	ug/L	
25) Chloroform	4.43	83	5398m	2.037	ug/L	
27) 1,2-Dichloroethane	5.19	62	4298	1.942	ug/L	99
29) Cyclohexane	4.73	56	4738	1.876	ug/L #	66
30) 1,1,1-Trichloroethane	4.66	97	4181	1.859	ug/L	94
31) Carbon tetrachloride	4.88	117	3316	1.737	ug/L	95
33) Benzene	5.15	78	10538	1.859	ug/L	100
34) Trichloroethene	5.96	95	2783	1.917	ug/L	91
35) Methylcyclohexane	6.18	83	4443	1.801	ug/L	94
37) 1,2-Dichloropropane	6.23	63	2922	1.866	ug/L #	89
38) Bromodichloromethane	6.56	83	3484	1.746	ug/L	99
39) cis-1,3-Dichloropropene	7.07	75	4711	1.914	ug/L	97
40) 4-Methyl-2-pentanone	7.28	43	8775	3.785	ug/L	94
42) Toluene	7.43	91	10889	1.849	ug/L	100
44) trans-1,3-Dichloropropene	7.70	75	4381	1.902	ug/L	99
45) 1,1,2-Trichloroethane	7.89	97	2423	1.750	ug/L	98
46) Tetrachloroethene	8.02	164	1612	1.704	ug/L	88
48) 2-Hexanone	8.19	43	6972	3.897	ug/L #	92
49) Dibromochloromethane	8.30	129	2774	1.865	ug/L	97
50) 1,2-Dibromoethane	8.40	107	2615	1.801	ug/L #	100
51) Chlorobenzene	8.93	112	6566	1.811	ug/L	97
52) Ethylbenzene	9.06	91	11620	1.762	ug/L	97
53) m,p-Xylene	9.19	106	4316	1.791	ug/L	98
54) o-xylene	9.59	106	4426	1.839	ug/L	92
55) Styrene	9.61	104	7174	1.769	ug/L	95
56) Isopropylbenzene	9.98	105	10895	1.730	ug/L	99
58) 1,1,2,2-Tetrachloroethane	10.29	83	5022	2.071	ug/L #	97
59) 1,2,3-Trichloropropane	10.32	75	3519	1.793	ug/L	92
61) Bromoform	9.78	173	1689	1.805	ug/L #	94
62) 1,3-Dichlorobenzene	11.23	146	4523	1.845	ug/L	96
63) 1,4-Dichlorobenzene	11.32	146	4827	1.964	ug/L	92
65) 1,2-Dichlorobenzene	11.69	146	4815	1.926	ug/L	93
66) 1,2-Dibromo-3-chloropropan	12.48	75	898m	1.747	ug/L	
67) 1,3,5-Trichlorobenzene	12.70	180	2462	1.538	ug/L #	96
68) 1,2,4-trichlorobenzene	13.32	180	1602	1.237	ug/L	99
69) Naphthalene	13.56	128	4678	1.117	ug/L #	93
70) 1,2,3-Trichlorobenzene	13.80	180	1716	1.326	ug/L	92

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

