

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U110118\
 Data File : VU027896.D
 Acq On : 01 Nov 2018 12:00
 Operator : MD/SY
 Sample : MDL03
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL03

Manual Integrations
 APPROVED

MMDadoda
 11/5/2018 4:10:23 PM

Quant Time: Nov 02 07:20:08 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR103018WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Fri Nov 02 08:02:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.89	114	85911	5.00	ug/L	0.00
28) Chlorobenzene-d5	9.09	117	81721	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.48	152	40261	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.40	65	23835	3.55	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	71.00%
7) Chloroethane-d5	1.68	69	19481	3.76	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	75.20%
11) 1,1-Dichloroethene-d2	2.28	63	37516	2.94	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	58.80%#
20) 2-Butanone-d5	4.20	46	86918	43.41	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	86.82%
24) Chloroform-d	4.66	84	43275	3.58	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	71.60%
26) 1,2-Dichloroethane-d4	5.31	65	31219	4.50	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	90.00%
32) Benzene-d6	5.34	84	88182	3.94	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	78.80%
36) 1,2-Dichloropropane-d6	6.34	67	31050	3.92	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	78.40%
41) Toluene-d8	7.57	98	81345	3.82	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	76.40%
43) trans-1,3-Dichloropropene-	7.85	79	11870	3.87	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	77.40%
46) 2-Hexanone-d5	8.31	63	69335	41.34	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	82.68%
57) 1,1,2,2-Tetrachloroethane-	10.43	84	22477	4.01	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	80.20%
64) 1,2-Dichlorobenzene-d4	11.86	152	30933	4.02	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	80.40%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	1.21	85	2479	0.293	ug/L	99
3) Chloromethane	1.33	50	2788	0.332	ug/L	92
5) Vinyl chloride	1.40	62	2672	0.320	ug/L #	75
6) Bromomethane	1.63	94	1003	0.366	ug/L	78
8) Chloroethane	1.70	64	1682m	0.344	ug/L	
9) Trichlorofluoromethane	1.89	101	3735	0.337	ug/L	98
10) 1,1,2-Trichloro-1,2,2-trif	2.29	101	2084	0.368	ug/L #	83
12) 1,1-Dichloroethene	2.29	96	1875	0.351	ug/L #	1
13) Acetone	2.33	43	4682	3.709	ug/L	84
14) Carbon disulfide	2.48	76	5647	0.296	ug/L	96
15) Methyl Acetate	2.62	43	1164m	0.345	ug/L	
16) Methylene chloride	2.71	84	2256	0.376	ug/L	88
17) Methyl tert-butyl Ether	3.01	73	5495	0.345	ug/L #	84
18) trans-1,2-Dichloroethene	2.99	96	1777	0.313	ug/L	92
19) 1,1-Dichloroethane	3.45	63	4257	0.356	ug/L	98
21) 2-Butanone	4.29	43	6887	3.138	ug/L	80
22) cis-1,2-Dichloroethene	4.23	96	1895	0.289	ug/L	65

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110118\
 Data File : VU027896.D
 Acq On : 01 Nov 2018 12:00
 Operator : MD/SY
 Sample : MDL03
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL03

Manual Integrations
 APPROVED

MMDadoda
 11/5/2018 4:10:23 PM

Quant Time: Nov 02 07:20:08 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR103018WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Fri Nov 02 08:02:52 2018
 Response via : Initial Calibration

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

23) Bromochloromethane	4.55	128	811	0.299	ug/L	84
25) Chloroform	4.68	83	9290	0.765	ug/L	95
27) 1,2-Dichloroethane	5.41	62	3027	0.357	ug/L #	78
29) 1,1,1-Trichloroethane	4.91	97	3724m	0.353	ug/L	
30) Cyclohexane	4.99	56	3851	0.311	ug/L #	84
31) Carbon tetrachloride	5.14	117	3151	0.348	ug/L	90
33) Benzene	5.39	78	8384	0.324	ug/L	100
34) Trichloroethene	6.19	95	2071	0.307	ug/L	94
35) Methylcyclohexane	6.42	83	3690	0.322	ug/L	90
37) 1,2-Dichloropropane	6.44	63	2551	0.346	ug/L #	90
38) Bromodichloromethane	6.76	83	2802	0.325	ug/L #	93
39) cis-1,3-Dichloropropene	7.28	75	3321	0.314	ug/L	97
40) 4-Methyl-2-pentanone	7.47	43	16668	3.200	ug/L #	89
42) Toluene	7.64	91	8853	0.327	ug/L	96
44) trans-1,3-Dichloropropene	7.89	75	2967m	0.327	ug/L	
45) 1,1,2-Trichloroethane	8.07	97	1371	0.302	ug/L	89
47) Tetrachloroethene	8.23	164	1876	0.363	ug/L	87
48) 2-Hexanone	8.37	43	12835	3.418	ug/L	98
49) Dibromochloromethane	8.48	129	1750	0.320	ug/L	97
50) 1,2-Dibromoethane	8.59	107	1483	0.342	ug/L #	91
51) Chlorobenzene	9.12	112	5352	0.319	ug/L	97
52) Ethylbenzene	9.25	91	9711	0.315	ug/L	97
53) m,p-xylene	9.38	106	3502	0.313	ug/L	95
54) o-xylene	9.78	106	3428	0.312	ug/L	94
55) Styrene	9.79	104	5309	0.288	ug/L	96
56) Isopropylbenzene	10.17	105	9122	0.305	ug/L	99
58) 1,1,2,2-Tetrachloroethane	10.46	83	1950	0.330	ug/L #	94
59) 1,2,3-Trichloropropane	10.50	75	1193	0.267	ug/L	93
61) Bromoform	9.96	173	873	0.295	ug/L #	77
62) 1,3-Dichlorobenzene	11.42	146	4345	0.336	ug/L	94
63) 1,4-Dichlorobenzene	11.51	146	4412	0.341	ug/L	96
65) 1,2-Dichlorobenzene	11.88	146	3754	0.307	ug/L #	91
66) 1,2-Dibromo-3-chloropropan	12.66	75	203m	0.212	ug/L	
67) 1,3,5-Trichlorobenzene	12.89	180	3087	0.296	ug/L	92
68) 1,2,4-trichlorobenzene	13.51	180	2445	0.276	ug/L	89
69) Naphthalene	13.75	128	3682	0.267	ug/L	95
70) 1,2,3-Trichlorobenzene	13.99	180	2522	0.312	ug/L	96

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

