

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\U111618\
 Data File : VU028216.D
 Acq On : 17 Nov 2018 04:52
 Operator : MD/SY
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD00513

Quant Time: Nov 17 05:58:57 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR111418WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Nov 17 03:46:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.89	114	142523	5.00	ug/L	0.00
28) Chlorobenzene-d5	9.09	117	140338	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.48	152	69485	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.40	65	62839	5.31	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	106.20%
7) Chloroethane-d5	1.68	69	49853	5.31	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	106.20%
11) 1,1-Dichloroethene-d2	2.28	63	118174	5.35	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	107.00%
20) 2-Butanone-d5	4.20	46	206570	60.22	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	120.44%
24) Chloroform-d	4.65	84	104414	5.55	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	111.00%
26) 1,2-Dichloroethane-d4	5.31	65	60141	5.73	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	114.60%
32) Benzene-d6	5.34	84	211108	5.24	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	104.80%
36) 1,2-Dichloropropane-d6	6.33	67	76649	5.32	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	106.40%
41) Toluene-d8	7.57	98	188540	5.03	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	100.60%
43) trans-1,3-Dichloropropene-	7.85	79	23494	4.57	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	91.40%
46) 2-Hexanone-d5	8.31	63	153648	52.33	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	104.66%
57) 1,1,2,2-Tetrachloroethane-	10.43	84	51490	5.05	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	101.00%
64) 1,2-Dichlorobenzene-d4	11.86	152	64314	4.95	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	99.00%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	1.21	85	65144	5.182 ug/L	99
3) Chloromethane	1.33	50	78846	4.907 ug/L	99
5) Vinyl chloride	1.40	62	78200	5.123 ug/L	100
6) Bromomethane	1.63	94	47781	6.535 ug/L	99
8) Chloroethane	1.70	64	44433	4.835 ug/L	99
9) Trichlorofluoromethane	1.89	101	91034	5.467 ug/L	99
10) 1,1,2-Trichloro-1,2,2-trif	2.29	101	47280	5.058 ug/L	97
12) 1,1-Dichloroethene	2.29	96	45778	4.831 ug/L	87
13) Acetone	2.33	43	136325	59.010 ug/L	97
14) Carbon disulfide	2.48	76	143568	4.518 ug/L	97
15) Methyl Acetate	2.62	43	32375	5.240 ug/L	97
16) Methylene chloride	2.70	84	66621	5.950 ug/L	95
17) Methyl tert-butyl Ether	3.01	73	138697	5.235 ug/L	98
18) trans-1,2-Dichloroethene	2.99	96	49272	4.890 ug/L	95
19) 1,1-Dichloroethane	3.45	63	111588	5.266 ug/L	100
21) 2-Butanone	4.28	43	216247	59.313 ug/L	99
22) cis-1,2-Dichloroethene	4.23	96	59181	5.328 ug/L	95

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU111618\
 Data File : VU028216.D
 Acq On : 17 Nov 2018 04:52
 Operator : MD/SY
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD00513

Quant Time: Nov 17 05:58:57 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR111418WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Nov 17 03:46:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.55	128	22447	5.114	ug/L	89
25) Chloroform	4.68	83	109324	5.752	ug/L	96
27) 1,2-Dichloroethane	5.41	62	73554	5.888	ug/L	97
29) 1,1,1-Trichloroethane	4.92	97	81316	5.119	ug/L	98
30) Cyclohexane	5.00	56	107078	4.808	ug/L	98
31) Carbon tetrachloride	5.13	117	67812	5.144	ug/L	97
33) Benzene	5.39	78	240530	5.088	ug/L	100
34) Trichloroethene	6.19	95	68980	5.903	ug/L	97
35) Methylcyclohexane	6.42	83	91560	4.558	ug/L	93
37) 1,2-Dichloropropane	6.43	63	69411	5.309	ug/L	100
38) Bromodichloromethane	6.76	83	70223	5.048	ug/L	98
39) cis-1,3-Dichloropropene	7.27	75	83087	4.700	ug/L	96
40) 4-Methyl-2-pentanone	7.47	43	510431	56.693	ug/L	96
42) Toluene	7.64	91	246453	5.056	ug/L	98
44) trans-1,3-Dichloropropene	7.88	75	64184	4.634	ug/L	99
45) 1,1,2-Trichloroethane	8.07	97	42226	5.350	ug/L	98
47) Tetrachloroethene	8.23	164	40667	4.281	ug/L	97
48) 2-Hexanone	8.37	43	371040	57.918	ug/L	98
49) Dibromochloromethane	8.48	129	41896	4.986	ug/L	96
50) 1,2-Dibromoethane	8.59	107	38366	5.163	ug/L	95
51) Chlorobenzene	9.12	112	142729	5.007	ug/L	96
52) Ethylbenzene	9.25	91	264827	4.977	ug/L	99
53) m,p-xylene	9.38	106	95260	4.914	ug/L	96
54) o-xylene	9.78	106	93469	5.008	ug/L	96
55) Styrene	9.79	104	158928	5.036	ug/L	99
56) Isopropylbenzene	10.17	105	247593	4.905	ug/L	99
58) 1,1,2,2-Tetrachloroethane	10.46	83	45000	4.423	ug/L #	98
59) 1,2,3-Trichloropropane	10.49	75	39334	5.356	ug/L	98
61) Bromoform	9.96	173	21834	4.647	ug/L	97
62) 1,3-Dichlorobenzene	11.42	146	106272	4.837	ug/L	96
63) 1,4-Dichlorobenzene	11.51	146	110103	4.951	ug/L	100
65) 1,2-Dichlorobenzene	11.88	146	102398	4.874	ug/L	97
66) 1,2-Dibromo-3-chloropropan	12.66	75	8060	5.287	ug/L	85
67) 1,3,5-Trichlorobenzene	12.89	180	82863	4.916	ug/L	99
68) 1,2,4-trichlorobenzene	13.50	180	71255	4.850	ug/L	98
69) Naphthalene	13.74	128	130102	4.752	ug/L	98
70) 1,2,3-Trichlorobenzene	13.99	180	66819	4.937	ug/L	100

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

