

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071819\
 Data File : VV011931.D
 Acq On : 18 Jul 2019 11:15
 Operator : SY/MD
 Sample : VSTDCCC005
 Misc : 25ML/MSVOA V/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00541

Quant Time: Jul 19 01:06:22 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071719WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Jul 18 02:17:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.66	114	188863	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.89	117	178988	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.29	152	86105	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	60637	4.37	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	87.40%
7) Chloroethane-d5	1.58	69	47484	4.13	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	82.60%
11) 1,1-Dichloroethene-d2	2.13	63	102471	4.52	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	90.40%
20) 2-Butanone-d5	3.95	46	176993	50.52	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	101.04%
24) Chloroform-d	4.40	84	132750	4.77	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	95.40%
26) 1,2-Dichloroethane-d4	5.08	65	69334	4.86	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	97.20%
32) Benzene-d6	5.09	84	253577	4.87	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	97.40%
36) 1,2-Dichloropropane-d6	6.12	67	77958	4.89	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	97.80%
41) Toluene-d8	7.36	98	231034	4.86	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	97.20%
43) trans-1,3-Dichloropropene-	7.66	79	28706	4.86	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	97.20%
46) 2-Hexanone-d5	8.13	63	126720	53.32	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	106.64%
57) 1,1,2,2-Tetrachloroethane-	10.26	84	51843	5.17	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	103.40%
64) 1,2-Dichlorobenzene-d4	11.67	152	82430	4.79	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	95.80%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.14	85	96312	4.980 ug/L	97
3) Chloromethane	1.25	50	89917	4.797 ug/L	100
5) Vinyl chloride	1.32	62	88476	4.907 ug/L	99
6) Bromomethane	1.54	94	41832	3.999 ug/L	100
8) Chloroethane	1.60	64	46851	4.569 ug/L	99
9) Trichlorofluoromethane	1.77	101	111432	4.915 ug/L	100
10) 1,1,2-Trichloro-1,2,2-trif	2.14	101	55824	5.097 ug/L	98
12) 1,1-Dichloroethene	2.14	96	50952	5.061 ug/L	97
13) Acetone	2.21	43	88411	49.726 ug/L	98
14) Carbon disulfide	2.32	76	138545	4.855 ug/L	98
15) Methyl Acetate	2.47	43	20651	5.463 ug/L	100
16) Methylene chloride	2.53	84	55167	5.228 ug/L	96
17) Methyl tert-butyl Ether	2.81	73	152954	5.023 ug/L	99
18) trans-1,2-Dichloroethene	2.79	96	69120	4.884 ug/L	98
19) 1,1-Dichloroethane	3.23	63	132751	4.975 ug/L	99
21) 2-Butanone	4.04	43	175886	50.455 ug/L	97
22) cis-1,2-Dichloroethene	3.96	96	77627	5.328 ug/L	99

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071819\
 Data File : VV011931.D
 Acq On : 18 Jul 2019 11:15
 Operator : SY/MD
 Sample : VSTDCCC005
 Misc : 25ML/MSVOA V/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00541

Quant Time: Jul 19 01:06:22 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071719WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Jul 18 02:17:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.30	128	30009	5.034	ug/L	97
25) Chloroform	4.42	83	135702	4.715	ug/L	95
27) 1,2-Dichloroethane	5.18	62	84037	4.957	ug/L	99
29) 1,1,1-Trichloroethane	4.66	97	112957	5.135	ug/L	98
30) Cyclohexane	4.72	56	120258	5.134	ug/L	99
31) Carbon tetrachloride	4.87	117	97678	5.085	ug/L	99
33) Benzene	5.14	78	287611	5.135	ug/L	100
34) Trichloroethene	5.95	95	78396	5.037	ug/L	99
35) Methylcyclohexane	6.17	83	121118	5.130	ug/L	99
37) 1,2-Dichloropropane	6.22	63	71789	5.252	ug/L	100
38) Bromodichloromethane	6.55	83	85032	5.078	ug/L	100
39) cis-1,3-Dichloropropene	7.07	75	99750	5.330	ug/L	99
40) 4-Methyl-2-pentanone	7.27	43	441911	53.793	ug/L	100
42) Toluene	7.43	91	311641	5.303	ug/L	98
44) trans-1,3-Dichloropropene	7.69	75	76518	5.144	ug/L	99
45) 1,1,2-Trichloroethane	7.88	97	46858	5.070	ug/L	97
47) Tetrachloroethene	8.02	164	59311	5.177	ug/L	98
48) 2-Hexanone	8.18	43	316086	53.751	ug/L	100
49) Dibromochloromethane	8.29	129	51391	5.128	ug/L	98
50) 1,2-Dibromoethane	8.40	107	42852	5.038	ug/L	98
51) Chlorobenzene	8.92	112	189785	5.116	ug/L	98
52) Ethylbenzene	9.05	91	336941	5.216	ug/L	100
53) m,p-xylene	9.18	106	125128	5.201	ug/L	94
54) o-xylene	9.59	106	119534	5.333	ug/L	97
55) Styrene	9.60	104	202475	5.317	ug/L	98
56) Isopropylbenzene	9.97	105	327722	5.324	ug/L	100
58) 1,1,1,2,2-Tetrachloroethane	10.29	83	47724	5.414	ug/L	99
59) 1,2,3-Trichloropropane	10.32	75	41175	5.078	ug/L	99
61) Bromoform	9.77	173	24611	5.108	ug/L	98
62) 1,3-Dichlorobenzene	11.22	146	142192	5.130	ug/L	100
63) 1,4-Dichlorobenzene	11.31	146	144068	5.051	ug/L	98
65) 1,2-Dichlorobenzene	11.69	146	135369	5.214	ug/L	100
66) 1,2-Dibromo-3-chloropropan	12.47	75	7921	5.373	ug/L	96
67) 1,3,5-Trichlorobenzene	12.69	180	111810	5.140	ug/L	99
68) 1,2,4-trichlorobenzene	13.31	180	89358	5.205	ug/L	99
69) Naphthalene	13.55	128	137250	5.205	ug/L	100
70) 1,2,3-Trichlorobenzene	13.79	180	81927	5.293	ug/L	98

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

