

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV041620\
 Data File : VV015526.D
 Acq On : 16 Apr 2020 10:21
 Operator : SY/MD
 Sample : VSTDCCC005
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00549

Quant Time: Apr 17 00:43:53 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR040920WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Apr 16 14:00:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.64	114	110399	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.87	117	103728	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.27	152	52497	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.31	65	31126	5.43	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	108.60%
7) Chloroethane-d5	1.58	69	27597	5.23	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	104.60%
11) 1,1-Dichloroethene-d2	2.12	63	62875	5.21	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	104.20%
20) 2-Butanone-d5	3.93	46	82540	51.32	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	102.64%
24) Chloroform-d	4.37	84	65091	5.13	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	102.60%
26) 1,2-Dichloroethane-d4	5.06	65	34165	5.00	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	100.00%
32) Benzene-d6	5.07	84	123501	5.37	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	107.40%
36) 1,2-Dichloropropane-d6	6.09	67	36571	5.10	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	102.00%
41) Toluene-d8	7.33	98	113552	5.17	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	103.40%
43) trans-1,3-Dichloropropene-	7.64	79	16312	5.23	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	104.60%
46) 2-Hexanone-d5	8.11	63	65866	43.81	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	87.62%
57) 1,1,2,2-Tetrachloroethane-	10.24	84	26340	4.63	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	92.60%
64) 1,2-Dichlorobenzene-d4	11.65	152	41067	4.87	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	97.40%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.14	85	53335	4.881 ug/L	99
3) Chloromethane	1.25	50	50208	4.825 ug/L	98
5) Vinyl chloride	1.32	62	46756	4.666 ug/L	97
6) Bromomethane	1.53	94	25699	4.641 ug/L	100
8) Chloroethane	1.59	64	29122	5.011 ug/L	99
9) Trichlorofluoromethane	1.76	101	65040	5.073 ug/L	100
10) 1,1,2-Trichloro-1,2,2-trif	2.13	101	37617	5.195 ug/L	99
12) 1,1-Dichloroethene	2.13	96	33971	4.902 ug/L	93
13) Acetone	2.21	43	67616	59.504 ug/L #	53
14) Carbon disulfide	2.31	76	106837	4.422 ug/L	99
15) Methyl Acetate	2.46	43	14827	4.410 ug/L	94
16) Methylene chloride	2.52	84	37128	4.573 ug/L	98
17) Methyl tert-butyl Ether	2.78	73	92185	4.977 ug/L	99
18) trans-1,2-Dichloroethene	2.78	96	38465	4.925 ug/L	96
19) 1,1-Dichloroethane	3.21	63	74266	5.045 ug/L	98
21) 2-Butanone	4.01	43	106337	57.908 ug/L #	76
22) cis-1,2-Dichloroethene	3.94	96	40343	4.976 ug/L	99

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV041620\
 Data File : VV015526.D
 Acq On : 16 Apr 2020 10:21
 Operator : SY/MD
 Sample : VSTDCCC005
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00549

Quant Time: Apr 17 00:43:53 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR040920WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Apr 16 14:00:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.27	128	16856	4.996	ug/L	97
25) Chloroform	4.40	83	77039	5.114	ug/L	98
27) 1,2-Dichloroethane	5.15	62	49083	5.039	ug/L	98
29) 1,1,1-Trichloroethane	4.63	97	69762	5.183	ug/L	100
30) Cyclohexane	4.70	56	67751	4.742	ug/L	100
31) Carbon tetrachloride	4.85	117	60784	5.189	ug/L	100
33) Benzene	5.12	78	159338	5.132	ug/L	100
34) Trichloroethene	5.94	95	43929	5.257	ug/L	99
35) Methylcyclohexane	6.15	83	70766	5.026	ug/L	99
37) 1,2-Dichloropropane	6.20	63	40640	5.278	ug/L	99
38) Bromodichloromethane	6.53	83	52773	5.139	ug/L	98
39) cis-1,3-Dichloropropene	7.05	75	58770	4.974	ug/L	98
40) 4-Methyl-2-pentanone	7.25	43	248840	49.215	ug/L	99
42) Toluene	7.41	91	169789	5.037	ug/L	98
44) trans-1,3-Dichloropropene	7.67	75	51295	5.134	ug/L	98
45) 1,1,2-Trichloroethane	7.86	97	26205	5.063	ug/L	97
47) Tetrachloroethene	8.00	164	32527	5.254	ug/L	98
48) 2-Hexanone	8.16	43	175715	49.739	ug/L	100
49) Dibromochloromethane	8.27	129	31884	5.142	ug/L	97
50) 1,2-Dibromoethane	8.37	107	24856	5.011	ug/L	100
51) Chlorobenzene	8.90	112	108991	5.073	ug/L	99
52) Ethylbenzene	9.03	91	196518	5.065	ug/L	99
53) m,p-xylene	9.16	106	73468	5.014	ug/L	99
54) o-xylene	9.56	106	69823	4.923	ug/L	96
55) Styrene	9.58	104	118304	4.963	ug/L	100
56) Isopropylbenzene	9.95	105	195827	5.036	ug/L	99
58) 1,1,1,2,2-Tetrachloroethane	10.26	83	29749	4.722	ug/L	97
59) 1,2,3-Trichloropropane	10.30	75	23902	5.020	ug/L	99
61) Bromoform	9.75	173	15459	5.021	ug/L	99
62) 1,3-Dichlorobenzene	11.20	146	87248	5.155	ug/L	96
63) 1,4-Dichlorobenzene	11.30	146	87211	5.064	ug/L	98
65) 1,2-Dichlorobenzene	11.67	146	78699	5.100	ug/L	97
66) 1,2-Dibromo-3-chloropropan	12.45	75	5769	4.886	ug/L	89
67) 1,3,5-Trichlorobenzene	12.67	180	66353	5.425	ug/L	99
68) 1,2,4-trichlorobenzene	13.29	180	58141	5.277	ug/L	99
69) Naphthalene	13.53	128	94986	4.576	ug/L	99
70) 1,2,3-Trichlorobenzene	13.77	180	50766	5.217	ug/L	99

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

