

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV081318\
 Data File : VV007013.D
 Acq On : 13 Aug 2018 18:41
 Operator : SY/MD
 Sample : VSTDCCC005EC
 Misc : 25 mL/MSVOA V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00586

Quant Time: Aug 14 05:37:51 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR080918WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Tue Aug 14 05:35:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.67	114	243715	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.90	117	240290	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.30	152	111339	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	59890	3.81	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	76.20%
7) Chloroethane-d5	1.58	69	55875	4.19	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	83.80%
11) 1,1-Dichloroethene-d2	2.13	63	125492	4.25	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	85.00%
20) 2-Butanone-d5	3.97	46	244914	54.13	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	108.26%
24) Chloroform-d	4.40	84	157521	4.94	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	98.80%
26) 1,2-Dichloroethane-d4	5.09	65	80254	4.98	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	99.60%
32) Benzene-d6	5.10	84	281672	4.56	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	91.20%
36) 1,2-Dichloropropane-d6	6.12	67	99637	4.72	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	94.40%
41) Toluene-d8	7.36	98	246047	4.68	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	93.60%
43) trans-1,3-Dichloropropene-	7.67	79	33784	4.60	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	92.00%
46) 2-Hexanone-d5	8.15	63	147235	49.04	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	98.08%
57) 1,1,2,2-Tetrachloroethane-	10.27	84	81948	4.98	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	99.60%
64) 1,2-Dichlorobenzene-d4	11.68	152	89847	4.64	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	92.80%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.14	85	121513	5.12 ug/L	99
3) Chloromethane	1.25	50	106996	4.68 ug/L	99
5) Vinyl chloride	1.32	62	118201	5.01 ug/L	98
6) Bromomethane	1.54	94	25466	4.97 ug/L	100
8) Chloroethane	1.60	64	70875	4.95 ug/L	98
9) Trichlorofluoromethane	1.77	101	154520	5.14 ug/L	99
10) 1,1,2-Trichloro-1,2,2-trif	2.14	101	93293	5.18 ug/L	98
12) 1,1-Dichloroethene	2.14	96	89201	5.10 ug/L	87
13) Acetone	2.21	43	107557	46.17 ug/L	97
14) Carbon disulfide	2.32	76	277451	4.80 ug/L	100
15) Methyl Acetate	2.47	43	31319	4.78 ug/L	97
16) Methylene chloride	2.54	84	101701	4.82 ug/L	94
17) Methyl tert-butyl Ether	2.82	73	211651	5.04 ug/L	99
18) trans-1,2-Dichloroethene	2.79	96	97954	5.23 ug/L	98
19) 1,1-Dichloroethane	3.23	63	169069	5.00 ug/L	97
21) 2-Butanone	4.06	43	272535	53.88 ug/L	97
22) cis-1,2-Dichloroethene	3.97	96	107077	5.04 ug/L	96

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV081318\
 Data File : VV007013.D
 Acq On : 13 Aug 2018 18:41
 Operator : SY/MD
 Sample : VSTDCCC005EC
 Misc : 25 mL/MSVOA V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00586

Quant Time: Aug 14 05:37:51 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR080918WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Tue Aug 14 05:35:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.31	128	45999	5.31	ug/L	94
25) Chloroform	4.43	83	193517	5.23	ug/L	99
27) 1,2-Dichloroethane	5.19	62	116463	5.05	ug/L	98
29) 1,1,1-Trichloroethane	4.66	97	161449	5.04	ug/L	99
30) Cyclohexane	4.73	56	157350	4.72	ug/L	100
31) Carbon tetrachloride	4.88	117	137926	5.13	ug/L	100
33) Benzene	5.15	78	429495	5.16	ug/L	100
34) Trichloroethene	5.96	95	101889	5.10	ug/L	96
35) Methylcyclohexane	6.18	83	154813	4.83	ug/L	98
37) 1,2-Dichloropropane	6.23	63	116065	5.09	ug/L	100
38) Bromodichloromethane	6.56	83	133026	5.14	ug/L	99
39) cis-1,3-Dichloropropene	7.08	75	133981	4.87	ug/L	100
40) 4-Methyl-2-pentanone	7.29	43	614809	49.95	ug/L	99
42) Toluene	7.44	91	435911	5.45	ug/L	99
44) trans-1,3-Dichloropropene	7.70	75	114542	5.07	ug/L	100
45) 1,1,2-Trichloroethane	7.89	97	74802	5.05	ug/L	94
47) Tetrachloroethene	8.02	164	84812	5.24	ug/L	99
48) 2-Hexanone	8.20	43	452367	52.11	ug/L	97
49) Dibromochloromethane	8.30	129	82976	5.34	ug/L	97
50) 1,2-Dibromoethane	8.40	107	65831	5.26	ug/L #	96
51) Chlorobenzene	8.93	112	263691	5.17	ug/L	98
52) Ethylbenzene	9.06	91	410415	5.14	ug/L	99
53) m,p-xylene	9.19	106	156586	5.26	ug/L	99
54) o-xylene	9.59	106	145488	5.16	ug/L	98
55) Styrene	9.61	104	261767	5.50	ug/L	99
56) Isopropylbenzene	9.98	105	383614	5.18	ug/L	99
58) 1,1,1,2,2-Tetrachloroethane	10.29	83	93078	5.17	ug/L	98
59) 1,2,3-Trichloropropane	10.32	75	66820	5.24	ug/L	100
61) Bromoform	9.78	173	43550	5.16	ug/L	100
62) 1,3-Dichlorobenzene	11.23	146	183501	5.09	ug/L	100
63) 1,4-Dichlorobenzene	11.32	146	195386	5.29	ug/L	95
65) 1,2-Dichlorobenzene	11.70	146	186777	5.27	ug/L	99
66) 1,2-Dibromo-3-chloropropan	12.48	75	11755	4.67	ug/L	93
67) 1,3,5-Trichlorobenzene	12.70	180	142407	5.22	ug/L	100
68) 1,2,4-trichlorobenzene	13.32	180	98278	4.91	ug/L	98
69) Naphthalene	13.56	128	131656	4.58	ug/L	99
70) 1,2,3-Trichlorobenzene	13.80	180	98096	5.09	ug/L	99

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

