

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV050820\
 Data File : VV015952.D
 Acq On : 08 May 2020 09:30
 Operator : SY/MD
 Sample : VSTDCCC005
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VSTD00560

Quant Time: May 09 00:40:28 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat May 09 00:31:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.64	114	153803	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.87	117	149667	5.00	ug/L	0.00
61) 1,4-Dichlorobenzene-d4	11.27	152	76854	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.31	65	39674	4.75	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	95.00%
7) Chloroethane-d5	1.58	69	36946	4.67	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	93.40%
11) 1,1-Dichloroethene-d2	2.12	63	88121	4.91	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	98.20%
20) 2-Butanone-d5	3.93	46	123159	51.74	ug/L	-0.01
Spiked Amount	50.000	Range	40 - 130	Recovery	=	103.48%
24) Chloroform-d	4.37	84	94319	5.17	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	103.40%
26) 1,2-Dichloroethane-d4	5.06	65	51121	4.89	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	97.80%
32) Benzene-d6	5.07	84	172472	4.84	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	96.80%
36) 1,2-Dichloropropane-d6	6.09	67	54311	5.00	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	100.00%
41) Toluene-d8	7.33	98	166039	4.93	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	98.60%
45) trans-1,3-Dichloropropene-	7.64	79	21417	5.02	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	100.40%
48) 2-Hexanone-d5	8.11	63	94392	50.36	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	100.72%
59) 1,1,2,2-Tetrachloroethane-	10.24	84	38374	5.02	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	100.40%
65) 1,2-Dichlorobenzene-d4	11.65	152	60416	4.84	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	96.80%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.14	85	62612	5.313	ug/L 98
3) Chloromethane	1.25	50	62726	5.257	ug/L 100
5) Vinyl chloride	1.32	62	60678	5.318	ug/L 99
6) Bromomethane	1.53	94	35535	5.488	ug/L 99
8) Chloroethane	1.60	64	36288	4.464	ug/L 98
9) Trichlorofluoromethane	1.76	101	85737	5.479	ug/L 100
10) 1,1,2-Trichloro-1,2,2-trif	2.13	101	48020	5.432	ug/L 98
12) 1,1-Dichloroethene	2.13	96	43529	5.143	ug/L 96
13) Acetone	2.23	43	83199	52.991	ug/L 69
14) Carbon disulfide	2.31	76	148955	5.330	ug/L 98
15) Methyl Acetate	2.46	43	20371	5.615	ug/L 95
16) Methylene chloride	2.52	84	50373	4.974	ug/L 96
17) Methyl tert-butyl Ether	2.79	73	114098	5.188	ug/L 99
18) trans-1,2-Dichloroethene	2.78	96	49483	5.290	ug/L 97
19) 1,1-Dichloroethane	3.21	63	95510	5.301	ug/L 97
21) 2-Butanone	4.02	43	131058	54.811	ug/L 99
22) cis-1,2-Dichloroethene	3.94	96	52032	5.314	ug/L 100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.27	128	21876	5.234	ug/L	96
25) Chloroform	4.40	83	98786	4.979	ug/L	99
27) 1,2-Dichloroethane	5.15	62	63210	5.185	ug/L	99
29) 1,1,1-Trichloroethane	4.63	97	88439	5.358	ug/L	99
30) Cyclohexane	4.70	56	86427	5.226	ug/L	99
31) Carbon tetrachloride	4.85	117	76547	5.374	ug/L	99
33) Benzene	5.12	78	202456	5.377	ug/L	100
34) Trichloroethene	5.94	95	56020	5.212	ug/L	98
35) Methylcyclohexane	6.15	83	87436	5.378	ug/L	99
37) 1,2-Dichloropropane	6.20	63	49788	5.204	ug/L	98
38) Bromodichloromethane	6.53	83	65175	5.365	ug/L	100
39) cis-1,3-Dichloropropene	7.05	75	70666	5.226	ug/L	100
40) 4-Methyl-2-pentanone	7.25	43	321384	53.898	ug/L	100
42) Toluene	7.41	91	219441	5.498	ug/L	98
43) 1,3,5-Trimethylbenzene	10.56	105	202815	5.535	ug/L #	100
44) 1,2,4-Trimethylbenzene	10.94	105	204320	5.580	ug/L #	100
46) trans-1,3-Dichloropropene	7.67	75	61980	5.426	ug/L	96
47) 1,1,2-Trichloroethane	7.86	97	33646	5.266	ug/L	99
49) Tetrachloroethene	8.00	164	40775	5.437	ug/L	98
50) 2-Hexanone	8.16	43	231510	54.903	ug/L	99
51) Dibromochloromethane	8.27	129	39706	5.531	ug/L	97
52) 1,2-Dibromoethane	8.37	107	32022	5.405	ug/L	92
53) Chlorobenzene	8.90	112	138280	5.387	ug/L	97
54) Ethylbenzene	9.03	91	246080	5.422	ug/L	99
55) m,p-xylene	9.16	106	92953	5.531	ug/L	98
56) o-xylene	9.56	106	87804	5.465	ug/L	96
57) Styrene	9.58	104	150541	5.602	ug/L	100
58) Isopropylbenzene	9.95	105	244311	5.535	ug/L	99
60) 1,1,1,2-Tetrachloroethane	10.26	83	37793	5.536	ug/L	98
62) Bromoform	9.75	173	19050	5.396	ug/L	97
63) 1,3-Dichlorobenzene	11.20	146	109675	5.297	ug/L	97
64) 1,4-Dichlorobenzene	11.30	146	110773	5.325	ug/L	99
66) 1,2-Dichlorobenzene	11.67	146	99103	5.288	ug/L	98
67) 1,2-Dibromo-3-chloropropan	12.45	75	6697	5.475	ug/L	92
68) 1,3,5-Trichlorobenzene	12.67	180	82082	5.422	ug/L	99
69) 1,2,4-trichlorobenzene	13.29	180	68494	5.266	ug/L	98
70) Naphthalene	13.53	128	105133	5.082	ug/L	99
71) 1,2,3-Trichlorobenzene	13.77	180	60563	5.278	ug/L	99

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

