

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U100419\
 Data File : VU034947.D
 Acq On : 04 Oct 2019 20:28
 Operator : JC/SP
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD00527

Quant Time: Oct 05 00:36:21 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR093019WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Oct 05 00:29:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.86	114	194779	5.00	ug/L	0.00
28) Chlorobenzene-d5	9.07	117	195862	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.46	152	97167	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.39	65	67722	4.91	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	98.20%
7) Chloroethane-d5	1.67	69	56914	5.06	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	101.20%
11) 1,1-Dichloroethene-d2	2.26	63	111095	4.70	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	94.00%
20) 2-Butanone-d5	4.16	46	174640	44.49	ug/L	-0.02
Spiked Amount	50.000	Range	40 - 130	Recovery	=	88.98%
24) Chloroform-d	4.62	84	116318	4.99	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	99.80%
26) 1,2-Dichloroethane-d4	5.29	65	64759	4.46	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	89.20%
32) Benzene-d6	5.32	84	235570	4.88	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	97.60%
36) 1,2-Dichloropropane-d6	6.31	67	77904	4.83	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	96.60%
41) Toluene-d8	7.54	98	228230	4.92	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	98.40%
43) trans-1,3-Dichloropropene-	7.83	79	26803	4.51	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	90.20%
46) 2-Hexanone-d5	8.29	63	133588	45.62	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	91.24%
57) 1,1,2,2-Tetrachloroethane-	10.41	84	64985	5.05	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	101.00%
64) 1,2-Dichlorobenzene-d4	11.84	152	82801	5.14	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	102.80%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
2) Dichlorodifluoromethane	1.20	85	52418	4.602	ug/L	98
3) Chloromethane	1.32	50	40570	4.313	ug/L	97
5) Vinyl chloride	1.40	62	49557	4.869	ug/L	100
6) Bromomethane	1.62	94	21781	4.388	ug/L	97
8) Chloroethane	1.69	64	30595	5.277	ug/L	100
9) Trichlorofluoromethane	1.88	101	69977	5.031	ug/L	98
10) 1,1,2-Trichloro-1,2,2-trif	2.28	101	43942	5.083	ug/L	94
12) 1,1-Dichloroethene	2.27	96	33876	5.181	ug/L	77
13) Acetone	2.31	43	119561	41.611	ug/L	94
14) Carbon disulfide	2.46	76	52391	4.818	ug/L	100
15) Methyl Acetate	2.60	43	26366	4.372	ug/L	97
16) Methylene chloride	2.68	84	48693	4.720	ug/L	89
17) Methyl tert-butyl Ether	2.98	73	141890	4.916	ug/L	95
18) trans-1,2-Dichloroethene	2.96	96	34328	5.109	ug/L	91
19) 1,1-Dichloroethane	3.42	63	97466	4.687	ug/L	97
21) 2-Butanone	4.25	43	173313	40.059	ug/L	86
22) cis-1,2-Dichloroethene	4.20	96	50677	4.921	ug/L	93

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U100419\
 Data File : VU034947.D
 Acq On : 04 Oct 2019 20:28
 Operator : JC/SP
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD00527

Quant Time: Oct 05 00:36:21 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR093019WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Oct 05 00:29:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
23) Bromochloromethane	4.52	128	22105	5.260	ug/L	92
25) Chloroform	4.65	83	111623	4.461	ug/L	96
27) 1,2-Dichloroethane	5.39	62	63752	4.325	ug/L	100
29) 1,1,1-Trichloroethane	4.89	97	82650	4.590	ug/L	96
30) Cyclohexane	4.97	56	56549	4.218	ug/L	95
31) Carbon tetrachloride	5.11	117	64097	4.592	ug/L	99
33) Benzene	5.37	78	187400	4.764	ug/L	100
34) Trichloroethene	6.16	95	48093	4.621	ug/L	91
35) Methylcyclohexane	6.39	83	58478	4.500	ug/L	94
37) 1,2-Dichloropropane	6.41	63	61894	4.606	ug/L	100
38) Bromodichloromethane	6.74	83	75827	4.520	ug/L	99
39) cis-1,3-Dichloropropene	7.25	75	76746	4.435	ug/L	99
40) 4-Methyl-2-pentanone	7.44	43	461152	39.598	ug/L	94
42) Toluene	7.61	91	206164	4.797	ug/L	100
44) trans-1,3-Dichloropropene	7.86	75	62460	4.298	ug/L	100
45) 1,1,2-Trichloroethane	8.05	97	45896	4.850	ug/L	95
47) Tetrachloroethene	8.21	164	35372	5.132	ug/L	98
48) 2-Hexanone	8.34	43	319343	41.325	ug/L	94
49) Dibromochloromethane	8.45	129	53220	4.963	ug/L	97
50) 1,2-Dibromoethane	8.57	107	37065	4.855	ug/L	98
51) Chlorobenzene	9.10	112	145929	5.014	ug/L	96
52) Ethylbenzene	9.23	91	236494	4.624	ug/L	96
53) m,p-Xylene	9.35	106	83303	4.672	ug/L	97
54) o-Xylene	9.76	106	89026	4.717	ug/L	91
55) Styrene	9.77	104	155642	4.851	ug/L	92
56) Isopropylbenzene	10.14	105	251591	4.652	ug/L	98
58) 1,1,1,2,2-Tetrachloroethane	10.43	83	68243	4.914	ug/L	97
59) 1,2,3-Trichloropropane	10.47	75	44910	4.673	ug/L	99
61) Bromoform	9.94	173	29383	4.870	ug/L	97
62) 1,3-Dichlorobenzene	11.40	146	122462	4.883	ug/L	99
63) 1,4-Dichlorobenzene	11.49	146	121066	4.697	ug/L	98
65) 1,2-Dichlorobenzene	11.85	146	122831	4.774	ug/L	98
66) 1,2-Dibromo-3-chloropropan	12.64	75	8479	4.385	ug/L	95
67) 1,3,5-Trichlorobenzene	12.86	180	95807	5.207	ug/L	98
68) 1,2,4-trichlorobenzene	13.49	180	65838	5.010	ug/L	97
69) Naphthalene	13.73	128	92995	4.215	ug/L	98
70) 1,2,3-Trichlorobenzene	13.97	180	66986	5.078	ug/L	98

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

