

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030921\
 Data File : VU042533.D
 Acq On : 09 Mar 2021 10:55
 Operator : SY/MD
 Sample : VSTDIC001
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTDIC001

Manual Integrations
 APPROVED

SAM
 3/12/2021 5:19:01 PM

Quant Time: Mar 10 06:01:47 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U030921DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Mar 10 05:59:54 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Fluorobenzene	6.124	96	19592	1.00	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.642	95	8719	1.11	ug/l	0.00
Spiked Amount 1.000			Recovery	=	111.00%	
68) 1,2-Dichlorobenzene-d4	12.201	152	8557	1.07	ug/l	0.00
Spiked Amount 1.000			Recovery	=	107.00%	
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	1.388	85	8580	1.11	ug/l	100
3) Chloromethane	1.527	50	7581	1.07	ug/l	97
4) Vinyl Chloride	1.607	62	7185	1.07	ug/l	99
5) Bromomethane	1.864	94	4605	1.32	ug/l	94
6) Chloroethane	1.938	64	4490	1.17	ug/l	95
7) Trichlorofluoromethane	2.147	101	11935	1.23	ug/l	96
8) 1,1,2-Trichloro-1,2,2-...	2.591	101	6195	1.21	ug/l	98
9) 1,1-Dichloroethene	2.591	96	4982	1.08	ug/l	98
10) Iodomethane	2.732	142	8350	1.23	ug/l	96
11) Allyl Chloride	2.932	41	9876	1.06	ug/l	97
12) Acrylonitrile	3.327	53	2484	2.19	ug/l	# 88
13) Acetone	2.636	43	3674	4.83	ug/l	92
14) Carbon Disulfide	2.800	76	16983	1.13	ug/l	99
15) Methylene Chloride	3.054	84	6186	1.08	ug/l	98
16) trans-1,2-Dichloroethene	3.362	96	5735	1.16	ug/l	98
17) 1,1-Dichloroethane	3.880	63	11229	1.14	ug/l	99
18) 2-Butanone	4.735	43	8102	5.33	ug/l	100
19) Cyclohexane	5.395	56	10267m	1.06	ug/l	
20) Methylcyclohexane	6.767	83	8422	0.85	ug/l	98
21) 2,2-Dichloropropane	4.678	77	10462	1.12	ug/l	97
22) cis-1,2-Dichloroethene	4.674	96	6079	1.11	ug/l	98
23) Diethyl Ether	2.382	59	4364	1.16	ug/l	95
24) tert-Butyl Alcohol	3.211	59	2537	9.17	ug/l	# 82
25) Methyl tert-Butyl Ether	3.382	73	14808	1.09	ug/l	94
26) Bromochloromethane	4.986	128	3111	1.24	ug/l	92
27) Chloroform	5.099	83	11401	1.16	ug/l	98
28) 1,1,1-Trichloroethane	5.321	97	11140	1.22	ug/l	98
29) 1,1-Dichloropropene	5.533	75	7702	0.95	ug/l	97
30) Carbon Tetrachloride	5.533	117	8786	1.04	ug/l	94
31) Isopropyl Ether	4.012	45	20032	1.10	ug/l	100
32) Ethyl-t-butyl ether	4.523	59	16832	1.03	ug/l	94
33) Tert-Amyl methyl ether	5.960	73	13368	0.90	ug/l	# 91
34) Propionitrile	4.800	54	1588	4.17	ug/l	# 87
35) Benzene	5.784	78	21291	0.98	ug/l	98
36) 1,2-Dichloroethane	5.806	62	8177	1.03	ug/l	# 95
37) Trichloroethene	6.552	130	6537	1.11	ug/l	96
38) 1,2-Dichloropropane	6.803	63	5734	0.97	ug/l	96
39) Methacrylonitrile	4.996	41	2294	1.00	ug/l	# 83
40) Methyl acrylate	4.864	55	3535	1.15	ug/l	# 87
41) Tetrahydrofuran	5.076	42	2260	2.33	ug/l	95
42) 1-Chlorobutane	5.468	56	11687	0.98	ug/l	96
43) Dibromomethane	6.928	93	3434	1.08	ug/l	99
44) Bromodichloromethane	7.118	83	7562	0.98	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030921\
 Data File : VU042533.D
 Acq On : 09 Mar 2021 10:55
 Operator : SY/MD
 Sample : VSTDIC001
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTDIC001

Manual Integrations
 APPROVED

SAM
 3/12/2021 5:19:01 PM

Quant Time: Mar 10 06:01:47 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U030921DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Mar 10 05:59:54 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 4-Methyl-2-Pentanone	7.812	43	24123	4.63	ug/l	99
46) t-1,4-Dichloro-2-butene	10.844	75	2759m	1.67	ug/l	
47) Methyl methacrylate	6.980	69	6082	1.89	ug/l	96
48) Ethyl methacrylate	8.346	69	5845	0.86	ug/l	96
49) Toluene	7.976	92	13211	0.96	ug/l	100
50) t-1,3-Dichloropropene	8.221	75	6921	0.89	ug/l	100
51) cis-1,3-Dichloropropene	7.620	75	7743	0.89	ug/l	93
52) 1,1,2-Trichloroethane	8.411	97	4234	0.98	ug/l	94
53) 1,3-Dichloropropane	8.587	76	7977	1.02	ug/l	98
54) 2-Hexanone	8.706	43	17234	4.56	ug/l	98
55) Dibromochloromethane	8.819	129	5543	1.06	ug/l	96
56) 1,2-Dibromoethane	8.935	107	4419	1.03	ug/l	99
58) Tetrachloroethene	8.558	164	6210	1.18	ug/l	95
59) Chlorobenzene	9.455	112	13947	0.93	ug/l	99
60) 1,1,1,2-Tetrachloroethane	9.542	131	5417	0.96	ug/l	99
61) Pentachloroethane	11.436	117	4349	0.98	ug/l	100
62) Hexachloroethane	12.484	117	4294	0.97	ug/l	92
63) Ethyl Benzene	9.578	91	24305	0.90	ug/l	96
64) m/p-Xylenes	9.703	106	18042	1.78	ug/l	98
65) o-Xylene	10.108	106	9137	0.92	ug/l	95
66) Styrene	10.124	104	14686	0.87	ug/l	98
67) Bromoform	10.301	173	3172	0.95	ug/l #	97
69) Isopropylbenzene	10.494	105	24200	0.90	ug/l	98
70) 1,1,2,2-Tetrachloroethane	10.793	83	6014	1.05	ug/l #	95
71) 1,2,3-Trichloropropane	10.838	75	5613m	1.24	ug/l	
72) Bromobenzene	10.790	156	6791	0.99	ug/l	99
73) n-propylbenzene	10.915	120	6398	0.89	ug/l	99
74) 2-Chlorotoluene	10.996	126	5911	0.92	ug/l	99
75) 1,3,5-Trimethylbenzene	11.095	105	20391	0.86	ug/l	95
76) 4-Chlorotoluene	11.105	126	5828	0.89	ug/l	90
77) tert-Butylbenzene	11.430	119	21457	0.92	ug/l	100
78) 1,2,4-Trimethylbenzene	11.475	105	20867	0.88	ug/l	99
79) sec-Butylbenzene	11.652	105	26962	0.89	ug/l	100
80) Nitrobenzene	13.224	77	590m	3.72	ug/l	
81) p-Isopropyltoluene	11.803	119	22843	0.89	ug/l	98
82) 1,3-Dichlorobenzene	11.754	146	13283	1.02	ug/l	97
83) 1,4-Dichlorobenzene	11.844	146	13365	1.02	ug/l	98
84) n-Butylbenzene	12.217	91	20223	0.82	ug/l	97
85) 1,2-Dichlorobenzene	12.221	146	13092	1.03	ug/l	99
86) 1,2-Dibromo-3-Chloropr...	13.012	75	982	0.84	ug/l	89
87) 1,2,4-Trichlorobenzene	13.851	180	8251	0.86	ug/l	99
88) Hexachlorobutadiene	14.028	225	5686	1.03	ug/l	96
89) Naphthalene	14.098	128	12447	0.70	ug/l #	96
90) 1,2,3-Trichlorobenzene	14.340	180	7954	0.90	ug/l	97

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

