

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012423\
 Data File : VU052790.D
 Acq On : 24 Jan 2023 14:35
 Operator : JC/MD
 Sample : RL CHECK
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RL CHECK

Manual Integrations
 APPROVED

Reviewed By :Krupa Patel 01/31/2023
 Supervised By :Mahesh Dadoda 01/31/2023

Quant Time: Jan 25 01:37:53 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U012323DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jan 24 05:16:59 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Fluorobenzene	6.111	96	33533	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.629	95	10561	0.947	ug/l	0.00
Spiked Amount 1.000			Recovery	=	95.000%	
68) 1,2-Dichlorobenzene-d4	12.188	152	11318	0.986	ug/l	0.00
Spiked Amount 1.000			Recovery	=	99.000%	
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.385	85	5605	0.494	ug/l	95
3) Chloromethane	1.517	50	5300	0.492	ug/l	92
4) Vinyl Chloride	1.604	62	4669	0.449	ug/l	99
5) Bromomethane	1.858	94	2787	0.516	ug/l	96
6) Chloroethane	1.935	64	2936	0.469	ug/l	91
7) Trichlorofluoromethane	2.137	101	7695	0.508	ug/l	97
8) 1,1,2-Trichloro-1,2,2-...	2.581	101	4136	0.499	ug/l	96
9) 1,1-Dichloroethene	2.581	96	4034	0.516	ug/l	91
10) Iodomethane	2.723	142	4394	0.507	ug/l	100
11) Allyl Chloride	2.925	41	4663	0.499	ug/l	94
12) Acrylonitrile	3.350	53	1497m	0.902	ug/l	
13) Acetone	2.658	43	8536m	3.628	ug/l	
14) Carbon Disulfide	2.790	76	13336	0.509	ug/l	98
15) Methylene Chloride	3.044	84	6280	0.664	ug/l	92
16) trans-1,2-Dichloroethene	3.353	96	4251	0.488	ug/l	86
17) 1,1-Dichloroethane	3.867	63	7869	0.535	ug/l #	94
18) 2-Butanone	4.732	43	7005	2.563	ug/l	98
19) Cyclohexane	5.382	56	5413m	0.419	ug/l	
20) Methylcyclohexane	6.761	83	5336	0.370	ug/l	90
21) 2,2-Dichloropropane	4.658	77	7067	0.518	ug/l	99
22) cis-1,2-Dichloroethene	4.665	96	4645	0.510	ug/l	96
23) Diethyl Ether	2.375	59	2669	0.483	ug/l	89
24) tert-Butyl Alcohol	3.362	59	2645m	3.971	ug/l	
25) Methyl tert-Butyl Ether	3.362	73	8664	0.483	ug/l #	81
26) Bromochloromethane	4.970	128	2258	0.532	ug/l	95
27) Chloroform	5.086	83	10096	0.641	ug/l	97
28) 1,1,1-Trichloroethane	5.314	97	8016	0.551	ug/l	97
29) 1,1-Dichloropropene	5.523	75	5409	0.441	ug/l	97
30) Carbon Tetrachloride	5.523	117	6501	0.492	ug/l	98
31) Isopropyl Ether	3.996	45	9827	0.465	ug/l	91
32) Ethyl-t-butyl ether	4.501	59	9071	0.460	ug/l	100
33) Tert-Amyl methyl ether	5.938	73	7687	0.441	ug/l	97
34) Propionitrile	4.870	54	1370m	2.189	ug/l	
35) Benzene	5.774	78	16892	0.475	ug/l	96
36) 1,2-Dichloroethane	5.793	62	5124	0.479	ug/l	99
37) Trichloroethene	6.539	130	4656	0.480	ug/l	98
38) 1,2-Dichloropropane	6.790	63	4133	0.458	ug/l	94
39) Methacrylonitrile	4.989	41	1034	0.414	ug/l #	55
40) Methyl acrylate	4.861	55	1881m	0.424	ug/l	
41) Tetrahydrofuran	5.083	42	1041m	0.837	ug/l	
42) 1-Chlorobutane	5.456	56	6942	0.443	ug/l	92
43) Dibromomethane	6.919	93	1959	0.429	ug/l	89
44) Bromodichloromethane	7.105	83	6244	0.503	ug/l	93

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012423\
 Data File : VU052790.D
 Acq On : 24 Jan 2023 14:35
 Operator : JC/MD
 Sample : RL CHECK
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RL CHECK

Quant Time: Jan 25 01:37:53 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U012323DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jan 24 05:16:59 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Krupa Patel 01/31/2023
 Supervised By :Mahesh Dadoda 01/31/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 4-Methyl-2-Pentanone	7.790	43	9077	1.890	ug/l	98
46) t-1,4-Dichloro-2-butene	10.822	75	2534m	1.006	ug/l	
47) Methyl methacrylate	6.960	69	3016	0.742	ug/l	95
48) Ethyl methacrylate	8.333	69	2687	0.365	ug/l	99
49) Toluene	7.967	92	8532	0.395	ug/l	90
50) t-1,3-Dichloropropene	8.211	75	4237	0.392	ug/l	94
51) cis-1,3-Dichloropropene	7.603	75	5350	0.420	ug/l #	89
52) 1,1,2-Trichloroethane	8.398	97	2856	0.439	ug/l	90
53) 1,3-Dichloropropane	8.574	76	5259	0.468	ug/l	99
54) 2-Hexanone	8.687	43	8208	1.953	ug/l	99
55) Dibromochloromethane	8.806	129	4089	0.481	ug/l	98
56) 1,2-Dibromoethane	8.922	107	3012	0.487	ug/l	94
58) Tetrachloroethene	8.549	164	4152	0.474	ug/l	94
59) Chlorobenzene	9.446	112	10382	0.441	ug/l	95
60) 1,1,1,2-Tetrachloroethane	9.529	131	4371	0.480	ug/l	96
61) Pentachloroethane	11.420	117	3259	0.437	ug/l	90
62) Hexachloroethane	12.468	117	3471	0.461	ug/l	99
63) Ethyl Benzene	9.568	91	14468	0.380	ug/l	99
64) m/p-Xylenes	9.690	106	10816	0.712	ug/l	98
65) o-Xylene	10.095	106	5772	0.394	ug/l	96
66) Styrene	10.111	104	8438	0.707	ug/l	97
67) Bromoform	10.288	173	2197	0.438	ug/l #	95
69) Isopropylbenzene	10.481	105	13986	0.375	ug/l	100
70) 1,1,2,2-Tetrachloroethane	10.777	83	3575	0.441	ug/l	96
71) 1,2,3-Trichloropropane	10.822	75	3312m	0.472	ug/l	
72) Bromobenzene	10.780	156	4422	0.467	ug/l	97
73) n-propylbenzene	10.902	120	4045	0.748	ug/l	88
74) 2-Chlorotoluene	10.983	126	3590	0.381	ug/l	94
75) 1,3,5-Trimethylbenzene	11.082	105	10867	0.692	ug/l	97
76) 4-Chlorotoluene	11.095	126	3403	0.352	ug/l	91
77) tert-Butylbenzene	11.417	119	12749	0.393	ug/l	95
78) 1,2,4-Trimethylbenzene	11.465	105	11240	0.729	ug/l	97
79) sec-Butylbenzene	11.639	105	14747	0.710	ug/l	93
80) Nitrobenzene	13.224	77	796m	1.587	ug/l	
81) p-Isopropyltoluene	11.787	119	11359	0.732	ug/l	96
82) 1,3-Dichlorobenzene	11.741	146	8331	0.431	ug/l	96
83) 1,4-Dichlorobenzene	11.835	146	8210	0.423	ug/l	99
84) n-Butylbenzene	12.204	91	10275	0.807	ug/l	100
85) 1,2-Dichlorobenzene	12.211	146	7995	0.438	ug/l	99
86) 1,2-Dibromo-3-Chloropr...	13.002	75	523	0.348	ug/l #	88
87) 1,2,4-Trichlorobenzene	13.838	180	4345	0.397	ug/l	94
88) Hexachlorobutadiene	14.018	225	3762	0.490	ug/l	91
89) Naphthalene	14.085	128	7806	0.368	ug/l	97
90) 1,2,3-Trichlorobenzene	14.327	180	4122	0.397	ug/l	91

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

