

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101922\
 Data File : VU051455.D
 Acq On : 19 Oct 2022 13:50
 Operator : JC/MD
 Sample : N4955-09 0.25PPB
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL-WATER-03-QT4-2022

Quant Time: Oct 20 07:40:45 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U101722DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Oct 20 07:37:57 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Krupa Patel 10/20/2022
 Supervised By :Mahesh Dadoda 10/25/2022

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

Internal Standards						
1) Fluorobenzene	6.112	96	46183	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.632	95	12327	0.761	ug/l	0.00
Spiked Amount 1.000			Recovery	=	76.000%	
68) 1,2-Dichlorobenzene-d4	12.192	152	14493	0.801	ug/l	0.00
Spiked Amount 1.000			Recovery	=	80.000%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.533	50	6716	0.264	ug/l	97
4) Vinyl Chloride	1.601	62	4437	0.202	ug/l	90
5) Bromomethane	1.855	94	3441	0.281	ug/l	91
6) Chloroethane	1.932	64	3923	0.313	ug/l	91
7) Trichlorofluoromethane	2.138	101	5134	0.209	ug/l	92
9) 1,1-Dichloroethene	2.578	96	3356	0.237	ug/l	91
10) Iodomethane	2.720	142	3759	0.205	ug/l	92
11) Allyl Chloride	2.922	41	5459	0.253	ug/l	91
12) Acrylonitrile	3.324	53	2348	0.563	ug/l	97
13) Acetone	2.639	43	5363m	1.786	ug/l	
14) Carbon Disulfide	2.790	76	11639	0.230	ug/l	# 93
15) Methylene Chloride	3.041	84	11008	0.530	ug/l	94
16) trans-1,2-Dichloroethene	3.347	96	3877	0.239	ug/l	90
17) 1,1-Dichloroethane	3.864	63	6929	0.237	ug/l	# 95
18) 2-Butanone	4.732	43	3697	1.015	ug/l	75
21) 2,2-Dichloropropane	4.662	77	4714	0.231	ug/l	93
22) cis-1,2-Dichloroethene	4.671	96	3139	0.204	ug/l	97
23) Diethyl Ether	2.376	59	2906	0.235	ug/l	93
24) tert-Butyl Alcohol	3.234	59	3582m	2.760	ug/l	
25) Methyl tert-Butyl Ether	3.366	73	8639	0.234	ug/l	93
26) Bromochloromethane	4.970	128	1630	0.217	ug/l	91
27) Chloroform	5.089	83	7313	0.247	ug/l	99
28) 1,1,1-Trichloroethane	5.317	97	5202	0.220	ug/l	96
30) Carbon Tetrachloride	5.517	117	4102	0.206	ug/l	96
31) Isopropyl Ether	4.002	45	7063	0.204	ug/l	86
32) Ethyl-t-butyl ether	4.510	59	6402	0.203	ug/l	96
34) Propionitrile	4.829	54	1195m	0.975	ug/l	
36) 1,2-Dichloroethane	5.800	62	3956	0.211	ug/l	98
38) 1,2-Dichloropropane	6.790	63	3550	0.213	ug/l	# 82
41) Tetrahydrofuran	5.080	42	923	0.420	ug/l	# 74
43) Dibromomethane	6.919	93	1903	0.220	ug/l	96
44) Bromodichloromethane	7.108	83	4410	0.211	ug/l	91
45) 4-Methyl-2-Pentanone	7.800	43	7373	0.896	ug/l	# 89
46) t-1,4-Dichloro-2-butene	10.825	75	2101m	0.458	ug/l	
47) Methyl methacrylate	6.967	69	1702	0.258	ug/l	89
52) 1,1,2-Trichloroethane	8.401	97	2731	0.221	ug/l	# 88
54) 2-Hexanone	8.690	43	4976	0.879	ug/l	87
56) 1,2-Dibromoethane	8.928	107	2177	0.201	ug/l	93
62) Hexachloroethane	12.475	117	2568	0.217	ug/l	87
64) m/p-Xylenes	9.694	106	6108	0.277	ug/l	99
67) Bromoform	10.288	173	1590	0.211	ug/l	# 93
70) 1,1,2,2-Tetrachloroethane	10.780	83	3675	0.220	ug/l	97
80) Nitrobenzene	13.224	77	852	1.031	ug/l	# 75

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

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