

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU020819\
 Data File : VU029384.D
 Acq On : 08 Feb 2019 17:46
 Operator : JC/SP
 Sample : VU0208WBSD01
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VU0208WBSD01

Manual Integrations
 APPROVED

MMDadoda
 2/18/2019 12:15:56 PM

Quant Time: Feb 09 04:04:17 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U020519DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Feb 08 10:52:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----	-----	-----	-----	-----	-----	-----
1) Fluorobenzene	5.74	96	51149	1.00	ug/l	0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.31	95	19726	0.99	ug/l	0.00
Spiked Amount 1.000			Recovery	=	99.00%	
68) 1,2-Dichlorobenzene-d4	11.86	152	21911	1.07	ug/l	0.00
Spiked Amount 1.000			Recovery	=	107.00%	
Target Compounds						
					Ovalue	
2) Dichlorodifluoromethane	1.21	85	37525	1.966	ug/l	99
3) Chloromethane	1.33	50	31810	1.824	ug/l	94
4) Vinyl Chloride	1.40	62	39705	1.942	ug/l	96
5) Bromomethane	1.63	94	27617	2.428	ug/l	97
6) Chloroethane	1.70	64	23151	1.938	ug/l	99
7) Trichlorofluoromethane	1.88	101	63263	2.215	ug/l	96
8) 1,1,2-Trichloro-1,2,2-trif	2.29	101	34915	2.316	ug/l	89
9) 1,1-Dichloroethene	2.28	96	32620	2.211	ug/l	72
10) Iodomethane	2.41	142	29217	1.583	ug/l #	78
11) Allyl Chloride	2.59	41	47762	1.930	ug/l	91
12) Acrylonitrile	2.94	53	14225	4.301	ug/l	90
13) Acetone	2.32	43	28569	8.964	ug/l #	80
14) Carbon Disulfide	2.48	76	108878	2.090	ug/l	98
15) Methylene Chloride	2.70	84	38093	2.129	ug/l #	76
16) trans-1,2-Dichloroethene	2.98	96	36655	2.221	ug/l #	78
17) 1,1-Dichloroethane	3.44	63	50553	1.933	ug/l #	93
18) 2-Butanone	4.28	43	32366	9.086	ug/l	98
19) Cyclohexane	4.99	56	38074	1.570	ug/l #	85
20) Methylcyclohexane	6.41	83	52402	2.058	ug/l #	81
21) 2,2-Dichloropropane	4.22	77	47177	1.949	ug/l #	87
22) cis-1,2-Dichloroethene	4.22	96	33886	2.180	ug/l	74
23) Diethyl Ether	2.10	59	24252	2.047	ug/l	88
24) tert-Butyl Alcohol	2.84	59	26322	21.376	ug/l	100
25) Methyl tert-Butyl Ether	3.00	73	89509	2.154	ug/l #	85
26) Bromochloromethane	4.55	128	14636	2.132	ug/l	71
27) Chloroform	4.67	83	53109	2.089	ug/l	98
28) 1,1,1-Trichloroethane	4.91	97	48749	2.068	ug/l	95
29) 1,1-Dichloropropene	5.13	75	41134	2.010	ug/l	89
30) Carbon Tetrachloride	5.13	117	44106	2.067	ug/l	98
31) Isopropyl Ether	3.58	45	75946	1.925	ug/l	91
32) Ethyl-t-butyl ether	4.08	59	77348	1.956	ug/l	89
33) Tert-Amyl methyl ether	5.58	73	72150	2.018	ug/l	98
34) Propionitrile	4.34	54	9305	9.937	ug/l #	86
35) Benzene	5.39	78	116042	2.051	ug/l	99
36) 1,2-Dichloroethane	5.41	62	33704	1.969	ug/l #	90
37) Trichloroethene	6.18	130	36301	2.236	ug/l	87
38) 1,2-Dichloropropane	6.43	63	29764	2.022	ug/l	95
39) Methacrylonitrile	4.56	41	8470	1.766	ug/l #	82
40) Methyl acrylate	4.43	55	12874	1.779	ug/l #	84
41) Tetrahydrofuran	4.66	42	7936	3.286	ug/l #	69
42) 1-Chlorobutane	5.06	56	52491	1.915	ug/l	85

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU020819\
 Data File : VU029384.D
 Acq On : 08 Feb 2019 17:46
 Operator : JC/SP
 Sample : VU0208WBSD01
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleID :
 VU0208WBSD01

Manual Integrations
 APPROVED

MMDadoda
 2/18/2019 12:15:56 PM

Quant Time: Feb 09 04:04:17 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U020519DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Feb 08 10:52:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) Dibromomethane	6.56	93	16011	2.041	ug/l	87
44) Bromodichloromethane	6.76	83	40235	2.013	ug/l	95
45) 4-Methyl-2-Pentanone	7.46	43	82528	8.975	ug/l #	88
46) t-1,4-Dichloro-2-butene	10.52	75	14479m	3.232	ug/l	
47) Methyl methacrylate	6.63	69	28502	4.130	ug/l #	70
48) Ethyl methacrylate	8.02	69	27753	1.874	ug/l	77
49) Toluene	7.63	92	74408	2.055	ug/l	98
50) t-1,3-Dichloropropene	7.88	75	35794	1.827	ug/l	95
51) cis-1,3-Dichloropropene	7.27	75	45033	1.935	ug/l #	87
52) 1,1,2-Trichloroethane	8.07	97	20797	2.003	ug/l	92
53) 1,3-Dichloropropane	8.24	76	36611	2.007	ug/l	100
54) 2-Hexanone	8.37	43	56682	8.794	ug/l	85
55) Dibromochloromethane	8.48	129	29670	2.010	ug/l	99
56) 1,2-Dibromoethane	8.59	107	23025	2.161	ug/l	95
58) Tetrachloroethene	8.22	164	35180	2.328	ug/l	90
59) Chlorobenzene	9.12	112	88148	2.161	ug/l	99
60) 1,1,1,2-Tetrachloroethane	9.21	131	32517	2.158	ug/l	97
61) Pentachloroethane	11.10	117	23873	2.044	ug/l	88
62) Hexachloroethane	12.14	117	23708	1.860	ug/l	91
63) Ethyl Benzene	9.25	91	146782	2.072	ug/l	94
64) m/p-Xylenes	9.37	106	118510	4.305	ug/l	84
65) o-Xylene	9.78	106	56890	2.105	ug/l	87
66) Styrene	9.79	104	94497	2.098	ug/l	92
67) Bromoform	9.96	173	18636	2.030	ug/l	98
69) Isopropylbenzene	10.16	105	152159	2.108	ug/l	95
70) 1,1,2,2-Tetrachloroethane	10.46	83	27762	2.086	ug/l	98
71) 1,2,3-Trichloropropane	10.49	75	19887m	2.176	ug/l	
72) Bromobenzene	10.45	156	39473	2.256	ug/l	84
73) n-propylbenzene	10.59	120	44163	2.196	ug/l	74
74) 2-Chlorotoluene	10.66	126	37112	2.168	ug/l	75
75) 1,3,5-Trimethylbenzene	10.77	105	126464	2.032	ug/l	94
76) 4-Chlorotoluene	10.77	126	38808	2.184	ug/l	71
77) tert-Butylbenzene	11.10	119	130107	2.115	ug/l	91
78) 1,2,4-Trimethylbenzene	11.14	105	129899	2.062	ug/l	95
79) sec-Butylbenzene	11.32	105	169259	2.103	ug/l	95
80) Nitrobenzene	12.87	77	3401m	5.735	ug/l	
81) p-Isopropyltoluene	11.47	119	145095	2.127	ug/l	93
82) 1,3-Dichlorobenzene	11.41	146	76025	2.215	ug/l	96
83) 1,4-Dichlorobenzene	11.50	146	76560	2.239	ug/l	95
84) n-Butylbenzene	11.89	91	127814	2.041	ug/l	97
85) 1,2-Dichlorobenzene	11.88	146	71711	2.209	ug/l	97
86) 1,2-Dibromo-3-Chloropropan	12.66	75	4561	1.836	ug/l #	58
87) 1,2,4-Trichlorobenzene	13.50	180	49260	2.161	ug/l	99
88) Hexachlorobutadiene	13.69	225	30550	2.314	ug/l	98
89) Naphthalene	13.74	128	85693	2.216	ug/l	98
90) 1,2,3-Trichlorobenzene	13.99	180	49995	2.399	ug/l	97

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

