

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U020519\
 Data File : VU029346.D
 Acq On : 05 Feb 2019 14:24
 Operator : JC/SP
 Sample : VSTDIC0.5
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTDIC0.5

Manual Integrations
 APPROVED

sam
 2/8/2019 2:26:31 PM

Quant Time: Feb 08 10:49:40 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U020519DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Feb 08 10:20:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----	-----	-----	-----	-----	-----	-----
1) Fluorobenzene	5.74	96	53469	1.00	ug/l	0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.31	95	20354	0.98	ug/l	0.00
Spiked Amount 1.000			Recovery	=	98.00%	
68) 1,2-Dichlorobenzene-d4	11.86	152	20775	0.97	ug/l	0.00
Spiked Amount 1.000			Recovery	=	97.00%	
Target Compounds						
					Ovalue	
2) Dichlorodifluoromethane	1.21	85	10762	0.539	ug/l	92
3) Chloromethane	1.33	50	9457	0.519	ug/l	96
4) Vinyl Chloride	1.40	62	10829	0.507	ug/l	98
5) Bromomethane	1.63	94	7833	0.659	ug/l	91
6) Chloroethane	1.70	64	6710	0.537	ug/l	98
7) Trichlorofluoromethane	1.88	101	14878	0.498	ug/l	98
8) 1,1,2-Trichloro-1,2,2-trif	2.29	101	8093	0.513	ug/l	91
9) 1,1-Dichloroethene	2.28	96	7914	0.513	ug/l	69
10) Iodomethane	2.41	142	5636m	0.285	ug/l	
11) Allyl Chloride	2.59	41	13307	0.514	ug/l	95
12) Acrylonitrile	2.94	53	3713	1.074	ug/l	91
13) Acetone	2.32	43	9165m	2.764	ug/l	
14) Carbon Disulfide	2.47	76	28542	0.524	ug/l	99
15) Methylene Chloride	2.70	84	11977	0.640	ug/l #	78
16) trans-1,2-Dichloroethene	2.98	96	8950	0.519	ug/l #	81
17) 1,1-Dichloroethane	3.44	63	12420	0.454	ug/l #	93
18) 2-Butanone	4.29	43	7821	2.100	ug/l	89
19) Cyclohexane	4.99	56	14192m	0.589	ug/l	
20) Methylcyclohexane	6.41	83	12758	0.479	ug/l	88
21) 2,2-Dichloropropane	4.22	77	12880	0.509	ug/l #	91
22) cis-1,2-Dichloroethene	4.22	96	7772	0.478	ug/l	87
23) Diethyl Ether	2.10	59	6531	0.527	ug/l #	88
24) tert-Butyl Alcohol	2.84	59	7727	6.003	ug/l #	94
25) Methyl tert-Butyl Ether	3.00	73	21554	0.496	ug/l #	84
26) Bromochloromethane	4.55	128	3573	0.498	ug/l	73
27) Chloroform	4.67	83	13061	0.491	ug/l	94
28) 1,1,1-Trichloroethane	4.91	97	11852	0.481	ug/l	91
29) 1,1-Dichloropropene	5.13	75	10880	0.509	ug/l	99
30) Carbon Tetrachloride	5.13	117	10650	0.477	ug/l	96
31) Isopropyl Ether	3.58	45	19469	0.472	ug/l	97
32) Ethyl-t-butyl ether	4.07	59	20488	0.496	ug/l	92
33) Tert-Amyl methyl ether	5.58	73	17739	0.475	ug/l	97
35) Benzene	5.39	78	29179	0.493	ug/l	94
36) 1,2-Dichloroethane	5.41	62	8650	0.484	ug/l #	84
37) Trichloroethene	6.19	130	8257	0.486	ug/l	82
38) 1,2-Dichloropropane	6.43	63	7669	0.498	ug/l	92
41) Tetrahydrofuran	4.66	42	2346	0.895	ug/l #	70
42) 1-Chlorobutane	5.06	56	14066	0.491	ug/l	89
43) Dibromomethane	6.56	93	4524	0.552	ug/l	92
44) Bromodichloromethane	6.76	83	9994	0.478	ug/l	88
45) 4-Methyl-2-Pentanone	7.47	43	21949	2.284	ug/l #	92

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU020519\
 Data File : VU029346.D
 Acq On : 05 Feb 2019 14:24
 Operator : JC/SP
 Sample : VSTDIC0.5
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTDIC0.5

Manual Integrations
 APPROVED

sam
 2/8/2019 2:26:31 PM

Quant Time: Feb 08 10:49:40 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U020519DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Feb 08 10:20:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

46) t-1,4-Dichloro-2-butene	10.52	75	4433m	0.950	ug/l	
47) Methyl methacrylate	6.63	69	6459	0.895	ug/l #	82
48) Ethyl methacrylate	8.03	69	7049	0.455	ug/l	77
49) Toluene	7.64	92	18560	0.490	ug/l	94
50) t-1,3-Dichloropropene	7.89	75	9562	0.467	ug/l	97
51) cis-1,3-Dichloropropene	7.27	75	11817	0.486	ug/l #	87
52) 1,1,2-Trichloroethane	8.07	97	5446	0.502	ug/l	96
53) 1,3-Dichloropropane	8.25	76	9082	0.476	ug/l	93
54) 2-Hexanone	8.37	43	14797	2.196	ug/l	85
55) Dibromochloromethane	8.48	129	7392	0.479	ug/l	98
56) 1,2-Dibromoethane	8.59	107	5749	0.516	ug/l	93
58) Tetrachloroethene	8.22	164	7720	0.489	ug/l	92
59) Chlorobenzene	9.12	112	20882	0.490	ug/l	99
60) 1,1,1,2-Tetrachloroethane	9.21	131	7508	0.477	ug/l	95
61) Pentachloroethane	11.10	117	5791	0.474	ug/l	92
62) Hexachloroethane	12.14	117	6121	0.459	ug/l	97
63) Ethyl Benzene	9.25	91	35859	0.484	ug/l	99
64) m/p-Xylenes	9.38	106	26920	0.936	ug/l	90
65) o-Xylene	9.78	106	13699	0.485	ug/l	87
66) Styrene	9.79	104	22722	0.483	ug/l	95
67) Bromoform	9.96	173	4592	0.478	ug/l	94
69) Isopropylbenzene	10.17	105	34256	0.454	ug/l	94
70) 1,1,2,2-Tetrachloroethane	10.46	83	6646	0.478	ug/l #	94
71) 1,2,3-Trichloropropane	10.49	75	4376m	0.449	ug/l	
72) Bromobenzene	10.45	156	8770	0.479	ug/l	89
73) n-propylbenzene	10.59	120	9477	0.451	ug/l	94
74) 2-Chlorotoluene	10.66	126	8480	0.474	ug/l	87
75) 1,3,5-Trimethylbenzene	10.77	105	30743	0.473	ug/l	97
76) 4-Chlorotoluene	10.77	126	9349	0.503	ug/l	72
77) tert-Butylbenzene	11.10	119	30139	0.469	ug/l	91
78) 1,2,4-Trimethylbenzene	11.15	105	30912	0.469	ug/l	93
79) sec-Butylbenzene	11.32	105	38573	0.458	ug/l	94
81) p-Isopropyltoluene	11.48	119	33436	0.469	ug/l	94
82) 1,3-Dichlorobenzene	11.42	146	17085	0.476	ug/l	98
83) 1,4-Dichlorobenzene	11.51	146	17505	0.490	ug/l	94
84) n-Butylbenzene	11.89	91	30216	0.462	ug/l	96
85) 1,2-Dichlorobenzene	11.88	146	16575	0.488	ug/l	98
86) 1,2-Dibromo-3-Chloropropan	12.66	75	1238	0.477	ug/l	89
87) 1,2,4-Trichlorobenzene	13.50	180	10504	0.441	ug/l	99
88) Hexachlorobutadiene	13.69	225	6740	0.488	ug/l	99
89) Naphthalene	13.75	128	16069	0.398	ug/l #	96
90) 1,2,3-Trichlorobenzene	13.99	180	10426	0.479	ug/l	94

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

