

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U092220\
 Data File : VU040249.D
 Acq On : 22 Sep 2020 16:04
 Operator : SY/MD
 Sample : VU0922WBS01
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VU0922WBS01

Quant Time: Sep 23 08:20:21 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U091120DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Sep 11 13:27:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	6.12	96	49330	1.00	ug/l	0.00

System Monitoring Compounds

57) 4-Bromofluorobenzene	10.63	95	19335	0.94	ug/l	0.00
Spiked Amount 1.000			Recovery	=	94.00%	
68) 1,2-Dichlorobenzene-d4	12.19	152	19241	1.04	ug/l	0.00
Spiked Amount 1.000			Recovery	=	104.00%	

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.39	85	42710	1.957	100
3) Chloromethane	1.53	50	41729	2.390	100
4) Vinyl Chloride	1.61	62	40142	2.036	100
5) Bromomethane	1.86	94	20680	3.347	93
6) Chloroethane	1.94	64	25787	2.284	97
7) Trichlorofluoromethane	2.15	101	55809	2.006	97
8) 1,1,2-Trichloro-1,2,2-trif	2.59	101	23155	1.863	98
9) 1,1-Dichloroethene	2.59	96	17929	1.884	97
13) Acetone	2.64	43	69713	9.668	100
14) Carbon Disulfide	2.80	76	29596	1.778	99
15) Methylene Chloride	3.06	84	45118	2.822	95
16) trans-1,2-Dichloroethene	3.37	96	20780	2.032	93
17) 1,1-Dichloroethane	3.89	63	51187	2.038	100
18) 2-Butanone	4.73	43	103254	9.481	98
19) Cyclohexane	5.40	56	33231	2.131	96
20) Methylcyclohexane	6.77	83	33981	2.151	98
22) cis-1,2-Dichloroethene	4.68	96	26885	2.085	89
25) Methyl tert-Butyl Ether	3.38	73	71332	2.070	99
26) Bromochloromethane	4.99	128	11771	1.879	96
27) Chloroform	5.11	83	54917	1.999	96
28) 1,1,1-Trichloroethane	5.33	97	46263	2.004	100
30) Carbon Tetrachloride	5.53	117	38413	2.026	96
35) Benzene	5.78	78	98153	1.992	97
36) 1,2-Dichloroethane	5.81	62	38767	1.957	99
37) Trichloroethene	6.55	130	26893	2.124	95
38) 1,2-Dichloropropane	6.80	63	31999	2.098	98
44) Bromodichloromethane	7.11	83	43305	2.078	99
45) 4-Methyl-2-Pentanone	7.81	43	231352	9.307	99
49) Toluene	7.97	92	64058	2.050	98
50) t-1,3-Dichloropropene	8.22	75	37647	1.928	100
51) cis-1,3-Dichloropropene	7.62	75	39076	1.945	98
52) 1,1,2-Trichloroethane	8.41	97	23073	1.904	97
54) 2-Hexanone	8.70	43	159882	9.099	97
55) Dibromochloromethane	8.82	129	27271	1.946	100
56) 1,2-Dibromoethane	8.93	107	21857	1.963	98
58) Tetrachloroethene	8.56	164	20507	2.179	92
59) Chlorobenzene	9.45	112	72771	1.980	99
63) Ethyl Benzene	9.57	91	112428	1.842	100
64) m/p-Xylenes	9.69	106	44631	3.851	95
65) o-Xylene	10.10	106	42345	1.904	99
66) Styrene	10.11	104	79584	2.005	97

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

67) Bromoform	10.29	173	15025	2.048	ug/l	95
69) Isopropylbenzene	10.48	105	118624	1.925	ug/l	99
70) 1,1,2,2-Tetrachloroethane	10.78	83	29198	1.851	ug/l	92
71) 1,2,3-Trichloropropane	10.82	75	22519	1.867	ug/l	83
75) 1,3,5-Trimethylbenzene	11.09	105	98014	1.825	ug/l	99
77) tert-Butylbenzene	11.47	119	12534	1.946	ug/l #	55
78) 1,2,4-Trimethylbenzene	11.47	105	95594	1.790	ug/l	99
82) 1,3-Dichlorobenzene	11.74	146	63169	2.034	ug/l	98
83) 1,4-Dichlorobenzene	11.83	146	64125	2.112	ug/l	99
85) 1,2-Dichlorobenzene	12.21	146	61043	2.109	ug/l	98
86) 1,2-Dibromo-3-Chloropropan	12.99	75	5075	1.890	ug/l	90
87) 1,2,4-Trichlorobenzene	13.83	180	36506	2.195	ug/l	98
89) Naphthalene	14.08	128	52443	1.708	ug/l	98
90) 1,2,3-Trichlorobenzene	14.32	180	33136	2.097	ug/l	96

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

