

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101321\
 Data File : VU045305.D
 Acq On : 13 Oct 2021 12:54
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICVVU101321

Quant Time: Oct 14 06:30:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U101321DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Oct 14 05:46:27 2021
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 i	Fluorobenzene	1.000	1.000	0.0	100	0.00
2 T	Dichlorodifluoromethane	0.382	0.231	39.5#	61	0.00
3 t	Chloromethane	0.398	0.320	19.6	84	0.00
4 Rt	Vinyl Chloride	0.413	0.338	18.2	83	0.00
5 T	Bromomethane	0.227	0.211	7.0	97	0.00
6 T	Chloroethane	0.268	0.227	15.3	90	0.00
7 T	Trichlorofluoromethane	0.481	0.434	9.8	92	0.00
8	1,1,2-Trichloro-1,2,2-trifl	0.278	0.260	6.5	94	0.00
9 Rt	1,1-Dichloroethene	0.260	0.254	2.3	98	0.00
10 t	Iodomethane	0.272	0.280	-2.9	92	0.00
11 t	Allyl Chloride	0.382	0.403	-5.5	104	0.00
12 t	Acrylonitrile	0.066	0.151	-128.8#	221#	0.00
13 T	Acetone	0.079	0.069	12.7	90	0.00
14 T	Carbon Disulfide	0.793	0.763	3.8	97	0.00
15 RT	Methylene Chloride	0.315	0.307	2.5	103	0.00
16 RT	trans-1,2-Dichloroethene	0.279	0.286	-2.5	102	0.00
17 t	1,1-Dichloroethane	0.524	0.544	-3.8	104	0.00
18 T	2-Butanone	0.101	0.079	21.8	77	0.00
19	Cyclohexane	0.485	0.491	-1.2	100	0.00
20	Methylcyclohexane	0.482	0.509	-5.6	102	0.00
21 T	2,2-Dichloropropane	0.446	0.460	-3.1	103	0.00
22 RT	cis-1,2-Dichloroethene	0.307	0.328	-6.8	106	0.00
23 t	Diethyl Ether	0.246	0.221	10.2	91	0.00
24 t	tert-Butyl Alcohol	0.019	0.009	52.6#	53	0.00
25 t	Methyl tert-Butyl Ether	0.726	0.717	1.2	99	0.00
26 t	Bromochloromethane	0.131	0.110	16.0	85	0.00
27 t	Chloroform	0.537	0.540	-0.6	104	0.00
28 RT	1,1,1-Trichloroethane	0.440	0.455	-3.4	103	0.00
29 T	1,1-Dichloropropene	0.402	0.412	-2.5	101	0.00
30 RT	Carbon Tetrachloride	0.341	0.361	-5.9	105	0.00
31 t	Isopropyl Ether	0.830	0.904	-8.9	107	0.00
32	Ethyl-t-butyl ether	0.836	0.000	100.0#	0#	-4.51#
33	Tert-Amyl methyl ether	0.740	0.000	100.0#	0#	-5.95#
34 t	Propionitrile	0.025	0.048	-92.0#	188	0.00
35 RT	Benzene	1.162	1.180	-1.5	102	0.00
36 RT	1,2-Dichloroethane	0.371	0.362	2.4	99	0.00
37 RT	Trichloroethene	0.283	0.290	-2.5	101	0.00
38 Rt	1,2-Dichloropropane	0.301	0.301	0.0	99	0.00
39 t	Methacrylonitrile	0.099	0.096	3.0	97	0.00
40 t	Methyl acrylate	0.165	0.173	-4.8	96	0.00
41 t	Tetrahydrofuran	0.057	0.127	-122.8#	233#	0.00
42 t	1-Chlorobutane	0.577	0.591	-2.4	102	0.00
43 T	Dibromomethane	0.150	0.153	-2.0	102	0.00
44 T	Bromodichloromethane	0.373	0.399	-7.0	104	0.00
45 T	4-Methyl-2-Pentanone	0.204	0.189	7.4	90	0.00
46 t	t-1,4-Dichloro-2-butene	0.072	0.066	8.3	82	0.00
47 t	Methyl methacrylate	0.152	0.074	51.3#	47	0.00
48 t	Ethyl methacrylate	0.300	0.305	-1.7	95	0.00
49 Rt	Toluene	0.715	0.746	-4.3	102	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
50 T	t-1,3-Dichloropropene	0.365	0.396	-8.5	104	0.00
51 T	cis-1,3-Dichloropropene	0.427	0.471	-10.3	106	0.00
52 RT	1,1,2-Trichloroethane	0.216	0.215	0.5	99	0.00
53 t	1,3-Dichloropropane	0.402	0.387	3.7	98	0.00
54 t	2-Hexanone	0.158	0.142	10.1	88	0.00
55 t	Dibromochloromethane	0.227	0.233	-2.6	99	0.00
56 T	1,2-Dibromoethane	0.203	0.204	-0.5	98	0.00
57 S	4-Bromofluorobenzene	0.367	0.387	-5.4	101	0.00
58 RT	Tetrachloroethene	0.255	0.265	-3.9	105	0.00
59 Rt	Chlorobenzene	0.741	0.756	-2.0	101	0.00
60 T	1,1,1,2-Tetrachloroethane	0.246	0.255	-3.7	102	0.00
61 t	Pentachloroethane	0.179	0.202	-12.8	106	0.00
62 t	Hexachloroethane	0.188	0.213	-13.3	106	0.00
63 Rt	Ethyl Benzene	1.331	1.460	-9.7	105	0.00
64 RT	m/p-Xylenes	0.505	0.564	-11.7	106	0.00
65 RT	o-Xylene	0.493	0.546	-10.8	107	0.00
66 RT	Styrene	0.807	0.913	-13.1	105	0.00
67 t	Bromoform	0.119	0.126	-5.9	101	0.00
68 S	1,2-Dichlorobenzene-d4	0.344	0.351	-2.0	97	0.00
69 T	Isopropylbenzene	1.296	1.450	-11.9	106	0.00
70 T	1,1,2,2-Tetrachloroethane	0.277	0.267	3.6	96	0.00
71 T	1,2,3-Trichloropropane	0.228	0.229	-0.4	109	0.00
72 t	Bromobenzene	0.286	0.305	-6.6	104	0.00
73 t	n-propylbenzene	0.339	0.394	-16.2	109	0.00
74 t	2-Chlorotoluene	0.293	0.323	-10.2	106	0.00
75 t	1,3,5-Trimethylbenzene	1.081	1.259	-16.5	109	0.00
76 t	4-Chlorotoluene	0.299	0.333	-11.4	106	0.00
77 t	tert-Butylbenzene	1.049	1.168	-11.3	106	0.00
78 t	1,2,4-Trimethylbenzene	1.104	1.250	-13.2	105	0.00
79 t	sec-Butylbenzene	1.447	1.628	-12.5	107	0.00
80	Nitrobenzene	0.008	0.003	62.5#	29	0.00
81 t	p-Isopropyltoluene	1.160	1.330	-14.7	106	0.00
82 t	1,3-Dichlorobenzene	0.572	0.626	-9.4	107	0.00
83 Rt	1,4-Dichlorobenzene	0.565	0.625	-10.6	108	0.00
84 t	n-Butylbenzene	1.106	1.310	-18.4	107	0.00
85 Rt	1,2-Dichlorobenzene	0.543	0.576	-6.1	104	0.00
86 t	1,2-Dibromo-3-Chloropropane	0.046	0.045	2.2	93	0.00
87 Rt	1,2,4-Trichlorobenzene	0.335	0.389	-16.1	107	0.00
88 t	Hexachlorobutadiene	0.157	0.180	-14.6	110	0.00
89 t	Naphthalene	0.658	0.747	-13.5	101	0.00
90 t	1,2,3-Trichlorobenzene	0.311	0.346	-11.3	103	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0