

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\U032923\
 Data File : U053469.D
 Acq On : 29 Mar 2023 16:54
 Operator : JC/MD
 Sample : VSTDICV010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICVU032923

Quant Time: Mar 30 11:12:42 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U032923DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Mar 30 11:12:17 2023
 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	92	0.00
2 T	Dichlorodifluoromethane	0.361	0.250	30.7#	59	0.00
3 t	Chloromethane	0.368	0.265	28.0	66	0.00
4 Rt	Vinyl Chloride	0.389	0.315	19.0	70	0.00
5 T	Bromomethane	0.330	0.264	20.0	82	0.00
6 T	Chloroethane	0.249	0.197	20.9	74	0.00
7 T	Trichlorofluoromethane	0.583	0.483	17.2	75	0.00
8	1,1,2-Trichloro-1,2,2-trifl	0.289	0.264	8.7	84	0.00
9 Rt	1,1-Dichloroethene	0.264	0.243	8.0	84	0.00
10 t	Iodomethane	0.268	0.262	2.2	75	0.00
11 t	Allyl Chloride	0.298	0.310	-4.0	88	0.00
12 t	Acrylonitrile	0.051	0.118	-131.4#	201#	-0.01
13 T	Acetone	0.101	0.056	44.6#	52	-0.02
14 T	Carbon Disulfide	0.810	0.648	20.0	70	0.00
15 RT	Methylene Chloride	0.624	0.381	38.9#	98	0.00
16 RT	trans-1,2-Dichloroethene	0.273	0.266	2.6	85	0.00
17 t	1,1-Dichloroethane	0.498	0.495	0.6	84	0.00
18 T	2-Butanone	0.102	0.065	36.3#	54	-0.03
19	Cyclohexane	0.410	0.355	13.4	74	0.00
20	Methylcyclohexane	0.512	0.487	4.9	79	0.00
21 T	2,2-Dichloropropane	0.426	0.447	-4.9	87	0.00
22 RT	cis-1,2-Dichloroethene	0.297	0.311	-4.7	88	0.00
23 t	Diethyl Ether	0.210	0.184	12.4	78	0.00
24 t	tert-Butyl Alcohol	0.019	0.009	52.6#	39	-0.05
25 t	Methyl tert-Butyl Ether	0.586	0.619	-5.6	89	0.00
26 t	Bromochloromethane	0.120	0.087	27.5	62	0.00
27 t	Chloroform	0.517	0.518	-0.2	86	0.00
28 RT	1,1,1-Trichloroethane	0.446	0.445	0.2	83	0.00
29 T	1,1-Dichloropropene	0.386	0.383	0.8	83	0.00
30 RT	Carbon Tetrachloride	0.355	0.371	-4.5	84	0.00
31 t	Isopropyl Ether	0.685	0.710	-3.6	87	-0.01
32	Ethyl-t-butyl ether	0.704	0.000	100.0#	0#	-4.51#
33	Tert-Amyl methyl ether	0.626	0.000	100.0#	0#	-5.94#
34 t	Propionitrile	0.017	0.038	-123.5#	163	-0.03
35 RT	Benzene	1.129	1.126	0.3	84	0.00
36 RT	1,2-Dichloroethane	0.324	0.315	2.8	85	0.00
37 RT	Trichloroethene	0.284	0.287	-1.1	83	0.00
38 Rt	1,2-Dichloropropane	0.289	0.286	1.0	83	0.00
39 t	Methacrylonitrile	0.070	0.072	-2.9	88	-0.01
40 t	Methyl acrylate	0.128	0.131	-2.3	95	-0.02
41 t	Tetrahydrofuran	0.037	0.091	-145.9#	204#	-0.02
42 t	1-Chlorobutane	0.513	0.513	0.0	85	0.00
43 T	Dibromomethane	0.138	0.148	-7.2	89	0.00
44 T	Bromodichloromethane	0.345	0.370	-7.2	87	0.00
45 T	4-Methyl-2-Pentanone	0.143	0.142	0.7	82	0.00
46 t	t-1,4-Dichloro-2-butene	0.078	0.052	33.3#	51	0.00
47 t	Methyl methacrylate	0.127	0.063	50.4#	42	0.00
48 t	Ethyl methacrylate	0.246	0.263	-6.9	86	0.00
49 Rt	Toluene	0.692	0.712	-2.9	85	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
50 T t-1,3-Dichloropropene	0.319	0.371	-16.3	91	0.00
51 T cis-1,3-Dichloropropene	0.395	0.437	-10.6	88	0.00
52 RT 1,1,2-Trichloroethane	0.200	0.207	-3.5	88	0.00
53 t 1,3-Dichloropropane	0.345	0.359	-4.1	87	0.00
54 t 2-Hexanone	0.150	0.112	25.3	63	0.00
55 t Dibromochloromethane	0.203	0.219	-7.9	84	0.00
56 T 1,2-Dibromoethane	0.184	0.194	-5.4	87	0.00
57 S 4-Bromofluorobenzene	0.361	0.368	-1.9	89	0.00
58 RT Tetrachloroethene	0.265	0.277	-4.5	83	0.00
59 Rt Chlorobenzene	0.762	0.760	0.3	82	0.00
60 T 1,1,1,2-Tetrachloroethane	0.231	0.253	-9.5	86	0.00
61 t Pentachloroethane	0.163	0.173	-6.1	91	0.00
62 t Hexachloroethane	0.187	0.215	-15.0	84	0.00
63 Rt Ethyl Benzene	1.336	1.413	-5.8	85	0.00
64 RT m/p-Xylenes	0.511	0.552	-8.0	86	0.00
65 RT o-Xylene	0.500	0.532	-6.4	86	0.00
66 RT Styrene	0.768	0.864	-12.5	86	0.00
67 t Bromoform	0.096	0.107	-11.5	85	0.00
68 S 1,2-Dichlorobenzene-d4	0.362	0.374	-3.3	90	0.00
69 T Isopropylbenzene	1.309	1.412	-7.9	86	0.00
70 T 1,1,2,2-Tetrachloroethane	0.240	0.249	-3.8	86	0.00
71 T 1,2,3-Trichloropropane	0.182	0.188	-3.3	102	0.00
72 t Bromobenzene	0.274	0.297	-8.4	88	0.00
73 t n-propylbenzene	0.346	0.390	-12.7	87	0.00
74 t 2-Chlorotoluene	0.299	0.320	-7.0	87	0.00
75 t 1,3,5-Trimethylbenzene	1.099	1.241	-12.9	89	0.00
76 t 4-Chlorotoluene	0.310	0.339	-9.4	88	0.00
77 t tert-Butylbenzene	1.019	1.128	-10.7	89	0.00
78 t 1,2,4-Trimethylbenzene	1.119	1.222	-9.2	85	0.00
79 t sec-Butylbenzene	1.467	1.637	-11.6	87	0.00
80 Nitrobenzene	0.006	0.002	66.7#	22	0.00
81 t p-Isopropyltoluene	1.152	1.307	-13.5	87	0.00
82 t 1,3-Dichlorobenzene	0.584	0.621	-6.3	86	0.00
83 Rt 1,4-Dichlorobenzene	0.591	0.635	-7.4	88	0.00
84 t n-Butylbenzene	1.139	1.327	-16.5	88	0.00
85 Rt 1,2-Dichlorobenzene	0.554	0.594	-7.2	88	0.00
86 t 1,2-Dibromo-3-Chloropropane	0.037	0.041	-10.8	90	0.00
87 Rt 1,2,4-Trichlorobenzene	0.323	0.378	-17.0	90	0.00
88 t Hexachlorobutadiene	0.154	0.170	-10.4	86	0.00
89 t Naphthalene	0.650	0.787	-21.1	93	0.00
90 t 1,2,3-Trichlorobenzene	0.278	0.340	-22.3	93	0.00

(#) = Out of Range

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