Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU022122\
Data File : VU047223.D
Acq On : 21 Feb 2022 10:18
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC010

Quant Time: Feb 22 03:57:19 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U021422DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Tue Feb 15 03:17:28 2022
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	0.333	0.333	0.0	93	0.00
3	t	Chloromethane	0.342	0.509	-48.8#	139	0.00
4	Rt	Vinyl Chloride	0.338	0.331	2.1	92	0.00
5	T	Bromomethane	0.221	0.094	57.5#	43	0.00
6	T	Chloroethane	0.194	0.195	-0.5	95	0.00
7	T	Trichlorofluoromethane	0.398	0.431	-8.3	101	0.00
8		1,1,2-Trichloro-1,2,2-trifl	0.230	0.239	-3.9	97	0.00
9	Rt	1,1-Dichloroethene	0.236	0.239	-1.3	98	0.00
10	t	Iodomethane	0.223	0.106	52.5#	39	0.00
11	t	Allyl Chloride	0.368	0.333	9.5	85	0.00
12	t	Acrylonitrile	0.069	0.064	7.2	101	0.01
13	T	Acetone	0.040	0.053	-32.5#	140	0.03
14	T	Carbon Disulfide	0.800	0.766	4.3	90	0.00
15	RT	Methylene Chloride	0.265	0.282	-6.4	103	0.00
16	RT	trans-1,2-Dichloroethene	0.249	0.255	-2.4	96	0.00
17	t	1,1-Dichloroethane	0.457	0.453	0.9	93	0.00
18	T	2-Butanone	0.070	0.082	-17.1	116	0.02
19		Cyclohexane	0.423	0.409	3.3	90	0.00
20		Methylcyclohexane	0.467	0.473	-1.3	94	0.00
21	T	2,2-Dichloropropane	0.408	0.424	-3.9	96	0.00
22	RT	cis-1,2-Dichloroethene	0.287	0.286	0.3	95	0.00
23	t	Diethyl Ether	0.191	0.192	-0.5	95	0.00
24	t	tert-Butyl Alcohol	0.012	0.013	-8.3	96	0.10
25	t	Methyl tert-Butyl Ether	0.663	0.662	0.2	95	0.00
26	t	Bromochloromethane	0.129	0.118	8.5	85	0.00
27	t	Chloroform	0.464	0.462	0.4	96	0.00
28	RT	1,1,1-Trichloroethane	0.402	0.411	-2.2	96	0.00
29	T	1,1-Dichloropropene	0.368	0.373	-1.4	95	0.00
30	RT	Carbon Tetrachloride	0.353	0.360	-2.0	94	0.00
31	t	Isopropyl Ether	0.731	0.722	1.2	92	0.00
32		Ethyl-t-butyl ether	0.743	0.753	-1.3	94	0.00
33		Tert-Amyl methyl ether	0.684	0.697	-1.9	95	0.00
34	t	Propionitrile	0.019	0.025	-31.6#	135	0.02
35	RT	Benzene	1.030	1.043	-1.3	94	0.00
36	RT	1,2-Dichloroethane	0.312	0.318	-1.9	95	0.00
37	RT	Trichloroethene	0.285	0.298	-4.6	98	0.00
38	Rt	1,2-Dichloropropane	0.273	0.271	0.7	93	0.00
39	t	Methacrylonitrile	0.094	0.100	-6.4	104	0.00
40	t	Methyl acrylate	0.159	0.186	-17.0	115	0.00
41	t	Tetrahydrofuran	0.048	0.051	-6.2	116	0.01
42	t	1-Chlorobutane	0.494	0.493	0.2	93	0.00
43	T	Dibromomethane	0.147	0.152	-3.4	96	0.00
44	T	Bromodichloromethane	0.348	0.353	-1.4	93	0.00
45	T	4-Methyl-2-Pentanone	0.199	0.191	4.0	88	0.00
46	t	t-1,4-Dichloro-2-butene	0.087	0.080	8.0	104	0.00
47	t	Methyl methacrylate	0.154	0.154	0.0	96	0.00
48	t	Ethyl methacrylate	0.313	0.314	-0.3	90	0.00
49	Rt	Toluene	0.664	0.677	-2.0	93	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
50	T	t-1,3-Dichloropropene	0.376	0.382	-1.6	92	0.00
51	T	cis-1,3-Dichloropropene	0.422	0.426	-0.9	92	0.00
52	RT	1,1,2-Trichloroethane	0.210	0.214	-1.9	96	0.00
53	t	1,3-Dichloropropane	0.375	0.368	1.9	90	0.00
54	t	2-Hexanone	0.148	0.139	6.1	88	0.00
55	t	Dibromochloromethane	0.255	0.256	-0.4	90	0.00
56	T	1,2-Dibromoethane	0.212	0.217	-2.4	93	0.00
57	S	4-Bromofluorobenzene	0.387	0.374	3.4	94	0.00
58	RT	Tetrachloroethene	0.263	0.272	-3.4	96	0.00
59	Rt	Chlorobenzene	0.741	0.748	-0.9	94	0.00
60	T	1,1,1,2-Tetrachloroethane	0.255	0.264	-3.5	94	0.00
61	t	Pentachloroethane	0.193	0.211	-9.3	96	0.00
62	t	Hexachloroethane	0.209	0.194	7.2	82	0.00
63	Rt	Ethyl Benzene	1.302	1.302	0.0	92	0.00
64	RT	m/p-Xylenes	0.496	0.500	-0.8	92	0.00
65	RT	o-Xylene	0.481	0.482	-0.2	91	0.00
66	RT	Styrene	0.816	0.818	-0.2	91	0.00
67	t	Bromoform	0.162	0.152	6.2	84	0.00
68	S	1,2-Dichlorobenzene-d4	0.403	0.384	4.7	92	0.00
69	T	Isopropylbenzene	1.293	1.297	-0.3	92	0.00
70	T	1,1,2,2-Tetrachloroethane	0.284	0.288	-1.4	94	0.00
71	T	1,2,3-Trichloropropane	0.239	0.227	5.0	83	0.00
72	t	Bromobenzene	0.316	0.320	-1.3	93	0.00
73	t	n-propylbenzene	0.356	0.365	-2.5	93	0.00
74	t	2-Chlorotoluene	0.304	0.310	-2.0	93	0.00
75	t	1,3,5-Trimethylbenzene	1.061	1.076	-1.4	92	0.00
76	t	4-Chlorotoluene	0.319	0.329	-3.1	94	0.00
77	t	tert-Butylbenzene	1.064	1.093	-2.7	93	0.00
78	t	1,2,4-Trimethylbenzene	1.082	1.091	-0.8	92	0.00
79	t	sec-Butylbenzene	1.413	1.431	-1.3	91	0.00
80		Nitrobenzene	0.014	0.007	50.0#	46	0.00
81	t	p-Isopropyltoluene	1.196	1.216	-1.7	92	0.00
82	t	1,3-Dichlorobenzene	0.615	0.626	-1.8	93	0.00
83	Rt	1,4-Dichlorobenzene	0.619	0.626	-1.1	94	0.00
84	t	n-Butylbenzene	1.148	1.173	-2.2	92	0.00
85	Rt	1,2-Dichlorobenzene	0.592	0.595	-0.5	93	0.00
86	t	1,2-Dibromo-3-Chloropropane	0.052	0.051	1.9	91	0.00
87	Rt	1,2,4-Trichlorobenzene	0.465	0.474	-1.9	93	0.00
88	t	Hexachlorobutadiene	0.238	0.249	-4.6	96	0.00
89	t	Naphthalene	0.952	0.923	3.0	90	0.00
90	t	1,2,3-Trichlorobenzene	0.417	0.429	-2.9	92	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0