Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU072920\
Data File : VU039715.D
Acq On : 29 Jul 2020 20:54
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Jul 30 02:44:08 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U072920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Thu Jul 30 01:45:12 2020
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	i	Fluorobenzene	1.000	1.000	0.0	101	0.00
2	T	Dichlorodifluoromethane	10.000	9.084	9.2	92	0.00
3	t	Chloromethane	10.000	9.266	7.3	99	0.00
4	Rt	Vinyl Chloride	10.000	9.319	6.8	98	0.00
5	T	Bromomethane	10.000	8.461	15.4	98	0.00
6	T	Chloroethane	10.000	9.045	9.6	98	0.00
7	T	Trichlorofluoromethane	10.000	9.265	7.3	96	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.245	7.6	97	0.00
9	Rt	1,1-Dichloroethene	10.000	9.316	6.8	97	0.00
10	t	Iodomethane	10.000	11.923	-19.2	103	0.00
11	t	Allyl Chloride	10.000	9.248	7.5	98	0.00
12	t	Acrylonitrile	20.000	19.961	0.2	107	0.00
13	T	Acetone	51.000	42.219	17.2	91	0.00
14	T	Carbon Disulfide	10.000	9.284	7.2	97	0.00
15	RT	Methylene Chloride	10.000	8.369	16.3	99	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.206	7.9	97	0.00
17	t	1,1-Dichloroethane	10.000	9.248	7.5	97	0.00
18	T	2-Butanone	50.000	46.882	6.2	101	0.00
19		Cyclohexane	10.000	9.204	8.0	99	0.00
20		Methylcyclohexane	10.000	9.577	4.2	93	0.00
21	T	2,2-Dichloropropane	10.000	8.450	15.5	91	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.425	5.7	99	0.00
23	t	Diethyl Ether	10.000	9.221	7.8	98	0.00
24	t	tert-Butyl Alcohol	100.000	94.628	5.4	100	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.471	5.3	98	0.00
26	t	Bromochloromethane	10.000	9.092	9.1	97	0.00
27	t	Chloroform	10.000	9.127	8.7	97	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.356	6.4	99	0.00
29	T	1,1-Dichloropropene	10.000	9.251	7.5	98	0.00
30	RT	Carbon Tetrachloride	10.000	9.307	6.9	97	0.00
31	t	Isopropyl Ether	10.000	9.256	7.4	97	0.00
32		Ethyl-t-butyl ether	10.000	9.412	5.9	99	0.00
33		Tert-Amyl methyl ether	10.000	9.392	6.1	96	0.00
34	t	Propionitrile	50.000	48.815	2.4	103	0.00
35	RT	Benzene	10.000	9.025	9.7	96	0.00
36	RT	1,2-Dichloroethane	10.000	9.025	9.7	95	0.00
37	RT	Trichloroethene	10.000	9.197	8.0	95	0.00
38	Rt	1,2-Dichloropropane	10.000	9.533	4.7	96	0.00
39	t	Methacrylonitrile	10.000	8.766	12.3	103	0.00
40	t	Methyl acrylate	10.000	9.251	7.5	99	0.00
41	t	Tetrahydrofuran	20.000	17.871	10.6	104	0.00
42	t	1-Chlorobutane	10.000	9.470	5.3	100	0.00
43	T	Dibromomethane	10.000	9.440	5.6	100	0.00
44	T	Bromodichloromethane	10.000	9.431	5.7	94	0.00
45	T	4-Methyl-2-Pentanone	50.000	49.602	0.8	98	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	19.549	2.3	109	0.00

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Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
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		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
47	t	Methyl methacrylate	20.000	19.986	0.1	96	0.00
48	t	Ethyl methacrylate	10.000	9.944	0.6	96	0.00
49	Rt	Toluene	10.000	9.633	3.7	95	0.00
50	T	t-1,3-Dichloropropene	10.000	9.938	0.6	92	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.473	5.3	92	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.511	4.9	98	0.00
53	t	1,3-Dichloropropane	10.000	9.347	6.5	96	0.00
54	t	2-Hexanone	50.000	49.695	0.6	98	0.00
55	t	Dibromochloromethane	10.000	9.923	0.8	97	0.00
56	T	1,2-Dibromoethane	10.000	9.623	3.8	97	0.00
57	S	4-Bromofluorobenzene	1.000	1.088	-8.8	106	0.00
58	RT	Tetrachloroethene	10.000	9.305	7.0	97	0.00
59	Rt	Chlorobenzene	10.000	9.582	4.2	95	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	9.660	3.4	96	0.00
61	t	Pentachloroethane	10.000	9.601	4.0	93	0.00
62	t	Hexachloroethane	10.000	9.981	0.2	94	0.00
63	Rt	Ethyl Benzene	10.000	9.939	0.6	93	0.00
64	RT	m/p-Xylenes	20.000	20.163	-0.8	94	0.00
65	RT	o-Xylene	10.000	10.049	-0.5	95	0.00
66	RT	Styrene	10.000	10.289	-2.9	95	0.00
67	t	Bromoform	10.000	10.118	-1.2	95	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.026	-2.6	101	0.00
69	T	Isopropylbenzene	10.000	9.972	0.3	94	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.870	1.3	99	0.00
71	T	1,2,3-Trichloropropane	10.000	8.617	13.8	78	0.00
72	t	Bromobenzene	10.000	9.666	3.3	96	0.00
73	t	n-propylbenzene	10.000	9.987	0.1	94	0.00
74	t	2-Chlorotoluene	10.000	9.830	1.7	96	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.266	-2.7	94	0.00
76	t	4-Chlorotoluene	10.000	10.208	-2.1	97	0.00
77	t	tert-Butylbenzene	10.000	10.064	-0.6	94	0.00
78	t	1,2,4-Trimethylbenzene	10.000	10.500	-5.0	96	0.00
79	t	sec-Butylbenzene	10.000	10.048	-0.5	94	0.00
80		Nitrobenzene	50.000	48.807	2.4	101	0.00
81	t	p-Isopropyltoluene	10.000	10.075	-0.7	93	0.00
82	t	1,3-Dichlorobenzene	10.000	9.452	5.5	94	0.00
83	Rt	1,4-Dichlorobenzene	10.000	9.516	4.8	96	0.00
84	t	n-Butylbenzene	10.000	9.831	1.7	93	0.00
85	Rt	1,2-Dichlorobenzene	10.000	9.545	4.6	95	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.631	-6.3	100	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	9.481	5.2	91	0.00
88	t	Hexachlorobutadiene	10.000	9.091	9.1	92	0.00
89	t	Naphthalene	10.000	10.064	-0.6	92	0.00
90	t	1,2,3-Trichlorobenzene	10.000	9.523	4.8	93	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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