Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050520\
Data File : VU038015.D
Acq On : 05 May 2020 14:15
Operator : JC/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: May 06 01:21:25 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U050420DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Tue May 05 02:20:02 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	99	0.00
2	T	Dichlorodifluoromethane	10.000	9.648	3.5	97	0.00
3	t	Chloromethane	10.000	9.240	7.6	97	0.00
4	Rt	Vinyl Chloride	10.000	9.637	3.6	97	0.00
5	T	Bromomethane	10.000	8.027	19.7	98	0.00
6	T	Chloroethane	10.000	9.571	4.3	97	0.00
7	T	Trichlorofluoromethane	10.000	9.754	2.5	98	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.761	2.4	99	0.00
9	Rt	1,1-Dichloroethene	10.000	9.627	3.7	100	0.00
10	t	Iodomethane	10.000	9.358	6.4	101	0.00
11	t	Allyl Chloride	10.000	9.691	3.1	98	0.00
12	t	Acrylonitrile	20.000	18.417	7.9	93	0.00
13	T	Acetone	51.000	49.021	3.9	101	0.00
14	T	Carbon Disulfide	10.000	9.411	5.9	97	0.00
15	RT	Methylene Chloride	10.000	10.301	-3.0	105	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.682	3.2	99	0.00
17	t	1,1-Dichloroethane	10.000	9.633	3.7	99	0.00
18	T	2-Butanone	50.000	47.164	5.7	97	0.00
19		Cyclohexane	10.000	9.638	3.6	98	0.00
20		Methylcyclohexane	10.000	10.004	-0.0	99	0.00
21	T	2,2-Dichloropropane	10.000	9.754	2.5	99	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.964	0.4	101	0.00
23	t	Diethyl Ether	10.000	9.904	1.0	101	0.00
24	t	tert-Butyl Alcohol	100.000	79.226	20.8	87	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.567	4.3	98	0.00
26	t	Bromochloromethane	10.000	9.530	4.7	98	0.00
27	t	Chloroform	10.000	9.656	3.4	100	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.828	1.7	99	0.00
29	T	1,1-Dichloropropene	10.000	9.684	3.2	98	0.00
30	RT	Carbon Tetrachloride	10.000	9.612	3.9	96	0.00
31	t	Isopropyl Ether	10.000	9.674	3.3	98	0.00
32		Ethyl-t-butyl ether	10.000	9.476	5.2	95	0.00
33		Tert-Amyl methyl ether	10.000	9.464	5.4	96	0.00
34	t	Propionitrile	50.000	43.752	12.5	91	0.00
35	RT	Benzene	10.000	9.748	2.5	100	0.00
36	RT	1,2-Dichloroethane	10.000	9.653	3.5	99	0.00
37	RT	Trichloroethene	10.000	9.313	6.9	97	0.00
38	Rt	1,2-Dichloropropane	10.000	9.700	3.0	100	0.00
39	t	Methacrylonitrile	10.000	9.122	8.8	97	0.00
40	t	Methyl acrylate	10.000	9.585	4.1	98	0.00
41	t	Tetrahydrofuran	20.000	17.548	12.3	96	0.00
42	t	1-Chlorobutane	10.000	9.791	2.1	99	0.00
43	T	Dibromomethane	10.000	9.550	4.5	100	0.00
44	T	Bromodichloromethane	10.000	9.963	0.4	101	0.00
45	T	4-Methyl-2-Pentanone	50.000	47.129	5.7	96	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	16.386	18.1	74	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	19.131	4.3	96	0.00
48	t	Ethyl methacrylate	10.000	9.705	2.9	95	0.00
49	Rt	Toluene	10.000	9.898	1.0	100	0.00
50	T	t-1,3-Dichloropropene	10.000	9.619	3.8	95	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.735	2.7	97	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.432	5.7	97	0.00
53	t	1,3-Dichloropropane	10.000	9.693	3.1	99	0.00
54	t	2-Hexanone	50.000	46.265	7.5	94	0.00
55	t	Dibromochloromethane	10.000	9.675	3.2	96	0.00
56	T	1,2-Dibromoethane	10.000	9.385	6.2	96	0.00
57	S	4-Bromofluorobenzene	1.000	0.957	4.3	94	0.00
58	RT	Tetrachloroethene	10.000	9.616	3.8	97	0.00
59	Rt	Chlorobenzene	10.000	9.693	3.1	98	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	9.640	3.6	96	0.00
61	t	Pentachloroethane	10.000	9.660	3.4	94	0.00
62	t	Hexachloroethane	10.000	9.080	9.2	91	0.00
63	Rt	Ethyl Benzene	10.000	9.906	0.9	98	0.00
64	RT	m/p-Xylenes	20.000	20.063	-0.3	99	0.00
65	RT	o-Xylene	10.000	9.909	0.9	99	0.00
66	RT	Styrene	10.000	10.112	-1.1	98	0.00
67	t	Bromoform	10.000	9.613	3.9	94	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.944	5.6	96	0.00
69	T	Isopropylbenzene	10.000	9.909	0.9	98	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.567	4.3	97	0.00
71	T	1,2,3-Trichloropropane	10.000	9.510	4.9	92	0.00
72	t	Bromobenzene	10.000	9.594	4.1	95	0.00
73	t	n-propylbenzene	10.000	9.885	1.2	96	0.00
74	t	2-Chlorotoluene	10.000	9.801	2.0	98	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.021	-0.2	98	0.00
76	t	4-Chlorotoluene	10.000	9.796	2.0	97	0.00
77	t	tert-Butylbenzene	10.000	9.805	2.0	97	0.00
78	t	1,2,4-Trimethylbenzene	10.000	9.981	0.2	98	0.00
79	t	sec-Butylbenzene	10.000	9.941	0.6	98	0.00
80		Nitrobenzene	50.000	38.622	22.8	80	0.00
81	t	p-Isopropyltoluene	10.000	9.987	0.1	97	0.00
82	t	1,3-Dichlorobenzene	10.000	9.578	4.2	97	0.00
83	Rt	1,4-Dichlorobenzene	10.000	9.428	5.7	96	0.00
84	t	n-Butylbenzene	10.000	10.210	-2.1	98	0.00
85	Rt	1,2-Dichlorobenzene	10.000	9.584	4.2	98	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	8.932	10.7	94	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	9.345	6.5	95	0.00
88	t	Hexachlorobutadiene	10.000	8.917	10.8	88	0.00
89	t	Naphthalene	10.000	9.424	5.8	94	0.00
90	t	1,2,3-Trichlorobenzene	10.000	9.350	6.5	94	0.00

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

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