Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU031020\
Data File : VU037151.D
Acq On : 10 Mar 2020 14:08
Operator : JC/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Mar 10 14:48:04 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U030920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Mon Mar 09 15:46:40 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	92	0.00
2	T	Dichlorodifluoromethane	10.000	10.146	-1.5	95	0.00
3	t	Chloromethane	10.000	10.876	-8.8	100	0.00
4	Rt	Vinyl Chloride	10.000	10.286	-2.9	94	0.00
5	T	Bromomethane	10.000	10.601	-6.0	106	0.00
6	T	Chloroethane	10.000	10.030	-0.3	96	0.00
7	T	Trichlorofluoromethane	10.000	10.399	-4.0	99	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	10.220	-2.2	96	0.00
9	Rt	1,1-Dichloroethene	10.000	10.212	-2.1	97	0.00
10	t	Iodomethane	10.000	10.708	-7.1	101	0.00
11	t	Allyl Chloride	10.000	10.028	-0.3	97	0.00
12	t	Acrylonitrile	20.000	20.899	-4.5	100	0.00
13	T	Acetone	51.000	51.705	-1.4	111	0.00
14	T	Carbon Disulfide	10.000	10.085	-0.9	95	0.00
15	RT	Methylene Chloride	10.000	9.409	5.9	101	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.068	-0.7	95	0.00
17	t	1,1-Dichloroethane	10.000	9.798	2.0	94	0.00
18	T	2-Butanone	50.000	49.848	0.3	99	0.00
19		Cyclohexane	10.000	10.369	-3.7	93	0.00
20		Methylcyclohexane	10.000	10.534	-5.3	95	0.00
21	T	2,2-Dichloropropane	10.000	10.093	-0.9	98	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.373	-3.7	97	0.00
23	t	Diethyl Ether	10.000	10.140	-1.4	98	0.00
24	t	tert-Butyl Alcohol	100.000	110.876	-10.9	101	0.00
25	t	Methyl tert-Butyl Ether	10.000	10.216	-2.2	98	0.00
26	t	Bromochloromethane	10.000	10.476	-4.8	99	0.00
27	t	Chloroform	10.000	9.795	2.1	97	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.235	-2.3	98	0.00
29	T	1,1-Dichloropropene	10.000	10.254	-2.5	96	0.00
30	RT	Carbon Tetrachloride	10.000	10.299	-3.0	95	0.00
31	t	Isopropyl Ether	10.000	9.977	0.2	93	0.00
32		Ethyl-t-butyl ether	10.000	10.028	-0.3	92	0.00
33		Tert-Amyl methyl ether	10.000	10.316	-3.2	94	0.00
34	t	Propionitrile	50.000	55.613	-11.2	103	0.00
35	RT	Benzene	10.000	10.181	-1.8	97	0.00
36	RT	1,2-Dichloroethane	10.000	9.826	1.7	96	0.00
37	RT	Trichloroethene	10.000	9.766	2.3	96	0.00
38	Rt	1,2-Dichloropropane	10.000	9.930	0.7	94	0.00
39	t	Methacrylonitrile	10.000	10.429	-4.3	97	0.00
40	t	Methyl acrylate	10.000	10.264	-2.6	98	0.00
41	t	Tetrahydrofuran	20.000	20.000	0.0	94	0.00
42	t	1-Chlorobutane	10.000	10.208	-2.1	94	0.00
43	T	Dibromomethane	10.000	10.433	-4.3	100	0.00
44	T	Bromodichloromethane	10.000	10.242	-2.4	95	0.00
45	T	4-Methyl-2-Pentanone	50.000	51.763	-3.5	97	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	21.932	-9.7	98	0.00

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Max. RRF Dev : 30% Max. Rel. Area : 200%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
47	t	Methyl methacrylate	20.000	22.355	-11.8	97	0.00
48	t	Ethyl methacrylate	10.000	10.876	-8.8	96	0.00
49	Rt	Toluene	10.000	10.695	-7.0	95	0.00
50	T	t-1,3-Dichloropropene	10.000	10.713	-7.1	98	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.162	-1.6	94	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.475	-4.7	98	0.00
53	t	1,3-Dichloropropane	10.000	10.294	-2.9	97	0.00
54	t	2-Hexanone	50.000	53.703	-7.4	97	0.00
55	t	Dibromochloromethane	10.000	10.653	-6.5	97	0.00
56	T	1,2-Dibromoethane	10.000	10.530	-5.3	93	0.00
57	S	4-Bromofluorobenzene	1.000	1.106	-10.6	102	0.00
58	RT	Tetrachloroethene	10.000	10.055	-0.5	95	0.00
59	Rt	Chlorobenzene	10.000	10.317	-3.2	96	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.251	-2.5	93	0.00
61	t	Pentachloroethane	10.000	10.549	-5.5	96	0.00
62	t	Hexachloroethane	10.000	10.535	-5.4	95	0.00
63	Rt	Ethyl Benzene	10.000	10.968	-9.7	96	0.00
64	RT	m/p-Xylenes	20.000	22.332	-11.7	95	0.00
65	RT	o-Xylene	10.000	10.884	-8.8	95	0.00
66	RT	Styrene	10.000	11.587	-15.9	97	0.00
67	t	Bromoform	10.000	11.126	-11.3	97	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.046	-4.6	95	0.00
69	T	Isopropylbenzene	10.000	11.027	-10.3	97	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.558	-5.6	98	0.00
71	T	1,2,3-Trichloropropane	10.000	10.058	-0.6	89	0.00
72	t	Bromobenzene	10.000	10.434	-4.3	100	0.00
73	t	n-propylbenzene	10.000	11.415	-14.1	97	0.00
74	t	2-Chlorotoluene	10.000	10.917	-9.2	96	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.462	-14.6	98	0.00
76	t	4-Chlorotoluene	10.000	11.399	-14.0	99	0.00
77	t	tert-Butylbenzene	10.000	10.768	-7.7	95	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.479	-14.8	96	0.00
79	t	sec-Butylbenzene	10.000	11.077	-10.8	97	0.00
80		Nitrobenzene	50.000	51.006	-2.0	87	0.00
81	t	p-Isopropyltoluene	10.000	11.492	-14.9	96	0.00
82	t	1,3-Dichlorobenzene	10.000	10.717	-7.2	98	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.543	-5.4	97	0.00
84	t	n-Butylbenzene	10.000	11.568	-15.7	96	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.363	-3.6	97	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.731	-7.3	94	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.179	-11.8	93	0.00
88	t	Hexachlorobutadiene	10.000	10.402	-4.0	97	0.00
89	t	Naphthalene	10.000	9.923	0.8	91	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.138	-11.4	95	0.00

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

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