Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU091120\
Data File : VU040108.D
Acq On : 11 Sep 2020 17:23
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Sep 12 05:20:21 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U091120DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Fri Sep 11 13:27:06 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	102	0.00
2	T	Dichlorodifluoromethane	10.000	9.691	3.1	98	0.00
3	t	Chloromethane	10.000	9.676	3.2	99	0.00
4	Rt	Vinyl Chloride	10.000	9.442	5.6	99	0.00
5	T	Bromomethane	10.000	14.324	-43.2#	118	0.00
6	T	Chloroethane	10.000	9.445	5.5	102	0.00
7	T	Trichlorofluoromethane	10.000	9.303	7.0	96	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.852	1.5	98	0.00
9	Rt	1,1-Dichloroethene	10.000	9.736	2.6	102	0.00
10	t	Iodomethane	10.000	0.000	100.0#	0	0.00
11	t	Allyl Chloride	10.000	0.000	100.0#	0	0.12
12	t	Acrylonitrile	20.000	0.000	100.0#	0	0.04
13	T	Acetone	50.000	47.573	4.9	95	0.03
14	T	Carbon Disulfide	10.000	8.847	11.5	97	0.00
15	RT	Methylene Chloride	10.000	7.999	20.0	96	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.293	-2.9	101	0.00
17	t	1,1-Dichloroethane	10.000	9.838	1.6	100	0.00
18	T	2-Butanone	50.000	52.064	-4.1	100	0.02
19		Cyclohexane	10.000	10.659	-6.6	103	0.00
20		Methylcyclohexane	10.000	10.960	-9.6	101	0.00
21	T	2,2-Dichloropropane	10.000	0.000	100.0#	0	-4.68#
22	RT	cis-1,2-Dichloroethene	10.000	10.576	-5.8	103	0.00
23	t	Diethyl Ether	10.000	0.000	100.0#	0	-2.39#
24	t	tert-Butyl Alcohol	100.000	0.000	100.0#	0	0.00
25	t	Methyl tert-Butyl Ether	10.000	10.350	-3.5	101	0.00
26	t	Bromochloromethane	10.000	9.795	2.1	99	0.00
27	t	Chloroform	10.000	9.884	1.2	100	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.800	2.0	100	0.00
29	T	1,1-Dichloropropene	10.000	0.000	100.0#	0	-5.54#
30	RT	Carbon Tetrachloride	10.000	9.964	0.4	99	0.00
31	t	Isopropyl Ether	10.000	0.000	100.0#	0	-4.02#
32		Ethyl-t-butyl ether	10.000	0.000	100.0#	0	0.16
33		Tert-Amyl methyl ether	10.000	0.000	100.0#	0	0.15
34	t	Propionitrile	50.000	0.000	100.0#	0	0.00
35	RT	Benzene	10.000	10.109	-1.1	98	0.00
36	RT	1,2-Dichloroethane	10.000	9.674	3.3	98	0.00
37	RT	Trichloroethene	10.000	10.050	-0.5	100	0.00
38	Rt	1,2-Dichloropropane	10.000	9.593	4.1	94	0.00
39	t	Methacrylonitrile	10.000	0.000	100.0#	0	-5.00#
40	t	Methyl acrylate	10.000	0.000	100.0#	0	-0.14
41	t	Tetrahydrofuran	20.000	0.000	100.0#	0	-5.09#
42	t	1-Chlorobutane	10.000	0.000	100.0#	0	-0.07
43	T	Dibromomethane	10.000	0.000	100.0#	0	0.19
44	T	Bromodichloromethane	10.000	9.844	1.6	98	0.00
45	T	4-Methyl-2-Pentanone	50.000	52.647	-5.3	97	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	0.000	100.0#	0	-10.83#

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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	0.000	100.0#	0	-6.98#
48	t	Ethyl methacrylate	10.000	0.000	100.0#	0	-8.34#
49	Rt	Toluene	10.000	10.602	-6.0	99	0.00
50	T	t-1,3-Dichloropropene	10.000	10.063	-0.6	95	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.341	-3.4	98	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.366	6.3	96	0.00
53	t	1,3-Dichloropropane	10.000	0.000	100.0#	0	-8.58#
54	t	2-Hexanone	50.000	53.591	-7.2	97	0.00
55	t	Dibromochloromethane	10.000	9.692	3.1	96	0.00
56	T	1,2-Dibromoethane	10.000	9.493	5.1	95	0.00
57	S	4-Bromofluorobenzene	1.000	0.957	4.3	98	0.00
58	RT	Tetrachloroethene	10.000	10.060	-0.6	98	0.00
59	Rt	Chlorobenzene	10.000	10.277	-2.8	99	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	0.000	100.0#	0	-9.53#
61	t	Pentachloroethane	10.000	0.000	100.0#	0	-11.42#
62	t	Hexachloroethane	10.000	0.000	100.0#	0	-12.47#
63	Rt	Ethyl Benzene	10.000	10.922	-9.2	99	0.00
64	RT	m/p-Xylenes	20.000	22.153	-10.8	100	0.00
65	RT	o-Xylene	10.000	11.209	-12.1	100	0.00
66	RT	Styrene	10.000	11.266	-12.7	99	0.00
67	t	Bromoform	10.000	10.009	-0.1	97	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.989	1.1	92	0.00
69	T	Isopropylbenzene	10.000	10.968	-9.7	98	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.627	3.7	95	0.00
71	T	1,2,3-Trichloropropane	10.000	9.669	3.3	96	0.00
72	t	Bromobenzene	10.000	0.000	100.0#	0	-10.78#
73	t	n-propylbenzene	10.000	0.000	100.0#	0	-10.90#
74	t	2-Chlorotoluene	10.000	0.000	100.0#	0	-10.98#
75	t	1,3,5-Trimethylbenzene	10.000	11.413	-14.1	97	0.00
76	t	4-Chlorotoluene	10.000	0.000	100.0#	0	-11.10#
77	t	tert-Butylbenzene	10.000	0.000	100.0#	0	-11.42#
78	t	1,2,4-Trimethylbenzene	10.000	11.535	-15.4	98	0.00
79	t	sec-Butylbenzene	10.000	0.000	100.0#	0	-11.64#
80		Nitrobenzene	50.000	0.000	100.0#	0	-0.21
81	t	p-Isopropyltoluene	10.000	0.000	100.0#	0	-11.79#
82	t	1,3-Dichlorobenzene	10.000	10.307	-3.1	99	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.452	-4.5	98	0.00
84	t	n-Butylbenzene	10.000	0.000	100.0#	0	-0.01
85	Rt	1,2-Dichlorobenzene	10.000	10.383	-3.8	100	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.114	-1.1	94	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.622	-16.2	100	0.00
88	t	Hexachlorobutadiene	10.000	0.000	100.0#	0	-14.01#
89	t	Naphthalene	10.000	12.493	-24.9	95	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.481	-14.8	99	0.00

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

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