Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090920\
Data File : VU040069.D
Acq On : 09 Sep 2020 18:42
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Sep 10 04:43:55 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U090920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Thu Sep 10 04:09:54 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	10.062	-0.6	101	0.00
3	t	Chloromethane	10.000	9.656	3.4	100	0.00
4	Rt	Vinyl Chloride	10.000	10.040	-0.4	105	0.00
5	T	Bromomethane	10.000	9.744	2.6	96	0.00
6	T	Chloroethane	10.000	9.070	9.3	98	0.00
7	T	Trichlorofluoromethane	10.000	10.307	-3.1	103	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.847	1.5	100	0.00
9	Rt	1,1-Dichloroethene	10.000	9.810	1.9	99	0.00
10	t	Iodomethane	10.000	10.265	-2.7	104	0.00
11	t	Allyl Chloride	10.000	9.863	1.4	102	0.00
12	t	Acrylonitrile	20.000	21.703	-8.5	112	0.00
13	T	Acetone	50.000	50.008	-0.0	102	0.00
14	T	Carbon Disulfide	10.000	9.864	1.4	100	0.00
15	RT	Methylene Chloride	10.000	10.149	-1.5	99	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.402	-4.0	103	0.00
17	t	1,1-Dichloroethane	10.000	9.863	1.4	103	0.00
18	T	2-Butanone	50.000	53.133	-6.3	110	0.02
19		Cyclohexane	10.000	10.131	-1.3	105	0.00
20		Methylcyclohexane	10.000	10.500	-5.0	99	0.00
21	T	2,2-Dichloropropane	10.000	9.462	5.4	102	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.910	0.9	101	0.00
23	t	Diethyl Ether	10.000	10.257	-2.6	106	0.00
24	t	tert-Butyl Alcohol	100.000	94.798	5.2	103	0.04
25	t	Methyl tert-Butyl Ether	10.000	10.103	-1.0	102	0.00
26	t	Bromochloromethane	10.000	10.077	-0.8	103	0.00
27	t	Chloroform	10.000	10.198	-2.0	106	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.003	-0.0	101	0.00
29	T	1,1-Dichloropropene	10.000	10.094	-0.9	102	0.00
30	RT	Carbon Tetrachloride	10.000	10.093	-0.9	101	0.00
31	t	Isopropyl Ether	10.000	10.097	-1.0	103	0.00
32		Ethyl-t-butyl ether	10.000	10.152	-1.5	102	0.00
33		Tert-Amyl methyl ether	10.000	10.388	-3.9	102	0.00
34	t	Propionitrile	50.000	55.600	-11.2	113	0.01
35	RT	Benzene	10.000	10.363	-3.6	102	0.00
36	RT	1,2-Dichloroethane	10.000	10.283	-2.8	103	0.00
37	RT	Trichloroethene	10.000	9.868	1.3	98	0.00
38	Rt	1,2-Dichloropropane	10.000	10.348	-3.5	103	0.00
39	t	Methacrylonitrile	10.000	9.517	4.8	105	0.00
40	t	Methyl acrylate	10.000	10.679	-6.8	111	0.00
41	t	Tetrahydrofuran	20.000	19.919	0.4	113	0.00
42	t	1-Chlorobutane	10.000	10.202	-2.0	102	0.00
43	T	Dibromomethane	10.000	10.107	-1.1	104	0.00
44	T	Bromodichloromethane	10.000	10.104	-1.0	103	0.00
45	T	4-Methyl-2-Pentanone	50.000	55.702	-11.4	107	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	25.613	-28.1	121	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	22.162	-10.8	104	0.00
48	t	Ethyl methacrylate	10.000	11.274	-12.7	104	0.00
49	Rt	Toluene	10.000	10.895	-8.9	106	0.00
50	T	t-1,3-Dichloropropene	10.000	10.952	-9.5	105	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.332	-3.3	100	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.287	-2.9	103	0.00
53	t	1,3-Dichloropropane	10.000	10.532	-5.3	104	0.00
54	t	2-Hexanone	50.000	54.827	-9.7	105	0.00
55	t	Dibromochloromethane	10.000	10.709	-7.1	105	0.00
56	T	1,2-Dibromoethane	10.000	10.688	-6.9	107	0.00
57	S	4-Bromofluorobenzene	1.000	0.940	6.0	89	0.00
58	RT	Tetrachloroethene	10.000	10.042	-0.4	100	0.00
59	Rt	Chlorobenzene	10.000	10.460	-4.6	102	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.398	-4.0	101	0.00
61	t	Pentachloroethane	10.000	10.131	-1.3	103	0.00
62	t	Hexachloroethane	10.000	10.408	-4.1	103	0.00
63	Rt	Ethyl Benzene	10.000	11.028	-10.3	101	0.00
64	RT	m/p-Xylenes	20.000	22.765	-13.8	102	0.00
65	RT	o-Xylene	10.000	11.022	-10.2	101	0.00
66	RT	Styrene	10.000	11.362	-13.6	100	0.00
67	t	Bromoform	10.000	11.199	-12.0	107	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.974	2.6	96	0.00
69	T	Isopropylbenzene	10.000	11.032	-10.3	101	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.269	-2.7	102	0.00
71	T	1,2,3-Trichloropropane	10.000	10.379	-3.8	103	0.00
72	t	Bromobenzene	10.000	10.442	-4.4	102	0.00
73	t	n-propylbenzene	10.000	10.905	-9.0	99	0.00
74	t	2-Chlorotoluene	10.000	10.832	-8.3	102	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.279	-12.8	100	0.00
76	t	4-Chlorotoluene	10.000	10.666	-6.7	100	0.00
77	t	tert-Butylbenzene	10.000	11.072	-10.7	102	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.224	-12.2	103	0.00
79	t	sec-Butylbenzene	10.000	10.897	-9.0	99	0.00
80		Nitrobenzene	50.000	54.871	-9.7	99	0.00
81	t	p-Isopropyltoluene	10.000	11.312	-13.1	101	0.00
82	t	1,3-Dichlorobenzene	10.000	10.336	-3.4	102	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.383	-3.8	101	0.00
84	t	n-Butylbenzene	10.000	11.552	-15.5	102	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.279	-2.8	100	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.524	-5.2	98	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.116	-11.2	101	0.00
88	t	Hexachlorobutadiene	10.000	10.619	-6.2	100	0.00
89	t	Naphthalene	10.000	9.768	2.3	99	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.021	-10.2	100	0.00

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

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