Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102220\
Data File : VU040924.D
Acq On : 23 Oct 2020 01:49
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Oct 23 07:47:51 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U102220DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Fri Oct 23 07:07:33 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	9.064	9.4	100	0.00
3	t	Chloromethane	10.000	9.026	9.7	99	0.00
4	Rt	Vinyl Chloride	10.000	9.286	7.1	98	0.00
5	T	Bromomethane	10.000	8.520	14.8	104	0.00
6	T	Chloroethane	10.000	9.101	9.0	100	0.00
7	T	Trichlorofluoromethane	10.000	9.483	5.2	100	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.344	6.6	99	0.00
9	Rt	1,1-Dichloroethene	10.000	9.470	5.3	99	0.00
10	t	Iodomethane	10.000	10.106	-1.1	102	0.00
11	t	Allyl Chloride	10.000	9.314	6.9	99	0.00
12	t	Acrylonitrile	20.000	19.455	2.7	102	0.00
13	T	Acetone	50.000	46.289	7.4	103	0.00
14	T	Carbon Disulfide	10.000	9.390	6.1	98	0.00
15	RT	Methylene Chloride	10.000	8.715	12.9	100	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.393	6.1	99	0.00
17	t	1,1-Dichloroethane	10.000	9.393	6.1	100	0.00
18	T	2-Butanone	50.000	48.634	2.7	101	0.00
19		Cyclohexane	10.000	9.507	4.9	99	0.00
20		Methylcyclohexane	10.000	10.359	-3.6	95	0.00
21	T	2,2-Dichloropropane	10.000	8.264	17.4	91	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.271	7.3	102	0.00
23	t	Diethyl Ether	10.000	9.441	5.6	98	0.00
24	t	tert-Butyl Alcohol	100.000	94.488	5.5	95	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.518	4.8	100	0.00
26	t	Bromochloromethane	10.000	9.630	3.7	102	0.00
27	t	Chloroform	10.000	9.354	6.5	100	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.312	6.9	99	0.00
29	T	1,1-Dichloropropene	10.000	9.510	4.9	99	0.00
30	RT	Carbon Tetrachloride	10.000	9.230	7.7	99	0.00
31	t	Isopropyl Ether	10.000	9.331	6.7	99	0.00
32		Ethyl-t-butyl ether	10.000	9.432	5.7	99	0.00
33		Tert-Amyl methyl ether	10.000	10.154	-1.5	99	0.00
34	t	Propionitrile	50.000	46.159	7.7	99	0.00
35	RT	Benzene	10.000	9.575	4.3	98	0.00
36	RT	1,2-Dichloroethane	10.000	9.444	5.6	101	0.00
37	RT	Trichloroethene	10.000	9.462	5.4	98	0.00
38	Rt	1,2-Dichloropropane	10.000	9.579	4.2	98	0.00
39	t	Methacrylonitrile	10.000	9.377	6.2	99	0.00
40	t	Methyl acrylate	10.000	9.776	2.2	101	0.00
41	t	Tetrahydrofuran	20.000	18.481	7.6	101	0.00
42	t	1-Chlorobutane	10.000	9.363	6.4	97	0.00
43	T	Dibromomethane	10.000	9.524	4.8	102	0.00
44	T	Bromodichloromethane	10.000	9.635	3.7	100	0.00
45	T	4-Methyl-2-Pentanone	50.000	53.948	-7.9	100	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	19.247	3.8	96	0.00

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Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
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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	21.953	-9.8	100	0.00
48	t	Ethyl methacrylate	10.000	10.813	-8.1	98	0.00
49	Rt	Toluene	10.000	10.378	-3.8	96	0.00
50	T	t-1,3-Dichloropropene	10.000	9.994	0.1	97	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.611	3.9	94	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.024	-0.2	103	0.00
53	t	1,3-Dichloropropane	10.000	10.029	-0.3	101	0.00
54	t	2-Hexanone	50.000	53.815	-7.6	100	0.00
55	t	Dibromochloromethane	10.000	9.685	3.1	101	0.00
56	T	1,2-Dibromoethane	10.000	9.980	0.2	101	0.00
57	S	4-Bromofluorobenzene	1.000	1.034	-3.4	101	0.00
58	RT	Tetrachloroethene	10.000	9.532	4.7	100	0.00
59	Rt	Chlorobenzene	10.000	9.913	0.9	97	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	9.711	2.9	100	0.00
61	t	Pentachloroethane	10.000	9.396	6.0	99	0.00
62	t	Hexachloroethane	10.000	9.975	0.3	99	0.00
63	Rt	Ethyl Benzene	10.000	10.795	-7.9	97	0.00
64	RT	m/p-Xylenes	20.000	19.683	1.6	99	0.00
65	RT	o-Xylene	10.000	11.060	-10.6	97	0.00
66	RT	Styrene	10.000	9.878	1.2	99	0.00
67	t	Bromoform	10.000	9.815	1.9	100	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.964	3.6	99	0.00
69	T	Isopropylbenzene	10.000	10.948	-9.5	97	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.905	1.0	101	0.00
71	T	1,2,3-Trichloropropane	10.000	10.723	-7.2	99	0.00
72	t	Bromobenzene	10.000	10.098	-1.0	97	0.00
73	t	n-propylbenzene	10.000	9.733	2.7	97	0.00
74	t	2-Chlorotoluene	10.000	10.486	-4.9	97	0.00
75	t	1,3,5-Trimethylbenzene	10.000	9.606	3.9	97	0.00
76	t	4-Chlorotoluene	10.000	11.091	-10.9	99	0.00
77	t	tert-Butylbenzene	10.000	10.799	-8.0	98	0.00
78	t	1,2,4-Trimethylbenzene	10.000	9.753	2.5	99	0.00
79	t	sec-Butylbenzene	10.000	11.050	-10.5	98	0.00
80		Nitrobenzene	50.000	51.151	-2.3	96	0.00
81	t	p-Isopropyltoluene	10.000	9.609	3.9	97	0.00
82	t	1,3-Dichlorobenzene	10.000	9.732	2.7	97	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.003	-0.0	101	0.00
84	t	n-Butylbenzene	10.000	11.097	-11.0	97	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.029	-0.3	99	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	9.701	3.0	101	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	10.054	-0.5	94	0.00
88	t	Hexachlorobutadiene	10.000	9.629	3.7	96	0.00
89	t	Naphthalene	10.000	8.983	10.2	97	0.00
90	t	1,2,3-Trichlorobenzene	10.000	10.934	-9.3	97	0.00

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

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