Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030920\
Data File : VU037141.D
Acq On : 09 Mar 2020 19:25
Operator : JC/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Mar 09 20:06:31 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U030920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Mon Mar 09 18:12:43 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	93	0.00
2	T	Dichlorodifluoromethane	10.000	10.447	-4.5	100	0.00
3	t	Chloromethane	10.000	10.451	-4.5	98	0.00
4	Rt	Vinyl Chloride	10.000	10.414	-4.1	97	0.00
5	T	Bromomethane	10.000	10.843	-8.4	111	0.00
6	T	Chloroethane	10.000	9.978	0.2	97	0.00
7	T	Trichlorofluoromethane	10.000	10.427	-4.3	101	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	10.422	-4.2	99	0.00
9	Rt	1,1-Dichloroethene	10.000	10.160	-1.6	98	0.00
10	t	Iodomethane	10.000	13.616	-36.2#	102	0.00
11	t	Allyl Chloride	10.000	10.166	-1.7	100	0.00
12	t	Acrylonitrile	20.000	20.392	-2.0	99	0.00
13	T	Acetone	51.000	48.178	5.5	105	0.00
14	T	Carbon Disulfide	10.000	10.108	-1.1	97	0.00
15	RT	Methylene Chloride	10.000	10.335	-3.4	113	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.318	-3.2	99	0.00
17	t	1,1-Dichloroethane	10.000	10.185	-1.9	99	0.00
18	T	2-Butanone	50.000	50.779	-1.6	103	0.00
19		Cyclohexane	10.000	10.704	-7.0	98	0.00
20		Methylcyclohexane	10.000	10.700	-7.0	98	0.00
21	T	2,2-Dichloropropane	10.000	9.669	3.3	95	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.786	-7.9	102	0.00
23	t	Diethyl Ether	10.000	10.149	-1.5	100	0.00
24	t	tert-Butyl Alcohol	100.000	107.559	-7.6	100	0.00
25	t	Methyl tert-Butyl Ether	10.000	10.450	-4.5	102	0.00
26	t	Bromochloromethane	10.000	10.564	-5.6	102	0.00
27	t	Chloroform	10.000	10.123	-1.2	102	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.427	-4.3	101	0.00
29	T	1,1-Dichloropropene	10.000	10.476	-4.8	100	0.00
30	RT	Carbon Tetrachloride	10.000	10.636	-6.4	100	0.00
31	t	Isopropyl Ether	10.000	10.445	-4.5	99	0.00
32		Ethyl-t-butyl ether	10.000	10.514	-5.1	99	0.00
33		Tert-Amyl methyl ether	10.000	10.777	-7.8	100	0.00
34	t	Propionitrile	50.000	54.560	-9.1	103	0.00
35	RT	Benzene	10.000	10.413	-4.1	100	0.00
36	RT	1,2-Dichloroethane	10.000	10.305	-3.0	102	0.00
37	RT	Trichloroethene	10.000	10.065	-0.6	100	0.00
38	Rt	1,2-Dichloropropane	10.000	10.383	-3.8	100	0.00
39	t	Methacrylonitrile	10.000	10.498	-5.0	99	0.00
40	t	Methyl acrylate	10.000	10.977	-9.8	107	0.00
41	t	Tetrahydrofuran	20.000	20.810	-4.0	100	0.00
42	t	1-Chlorobutane	10.000	10.699	-7.0	100	0.00
43	T	Dibromomethane	10.000	10.730	-7.3	104	0.00
44	T	Bromodichloromethane	10.000	10.658	-6.6	101	0.00
45	T	4-Methyl-2-Pentanone	50.000	52.746	-5.5	100	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	21.957	-9.8	100	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	23.056	-15.3	101	0.00
48	t	Ethyl methacrylate	10.000	11.066	-10.7	100	0.00
49	Rt	Toluene	10.000	11.052	-10.5	99	0.00
50	T	t-1,3-Dichloropropene	10.000	10.811	-8.1	100	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.477	-4.8	99	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.622	-6.2	101	0.00
53	t	1,3-Dichloropropane	10.000	10.503	-5.0	100	0.00
54	t	2-Hexanone	50.000	54.915	-9.8	101	0.00
55	t	Dibromochloromethane	10.000	10.756	-7.6	100	0.00
56	T	1,2-Dibromoethane	10.000	11.113	-11.1	100	0.00
57	S	4-Bromofluorobenzene	1.000	1.114	-11.4	104	0.00
58	RT	Tetrachloroethene	10.000	10.557	-5.6	102	0.00
59	Rt	Chlorobenzene	10.000	10.427	-4.3	98	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.351	-3.5	95	0.00
61	t	Pentachloroethane	10.000	10.620	-6.2	98	0.00
62	t	Hexachloroethane	10.000	10.965	-9.6	101	0.00
63	Rt	Ethyl Benzene	10.000	11.133	-11.3	99	0.00
64	RT	m/p-Xylenes	20.000	22.625	-13.1	98	0.00
65	RT	o-Xylene	10.000	11.191	-11.9	99	0.00
66	RT	Styrene	10.000	11.703	-17.0	99	0.00
67	t	Bromoform	10.000	11.209	-12.1	99	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.070	-7.0	99	0.00
69	T	Isopropylbenzene	10.000	11.131	-11.3	99	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.629	-6.3	100	0.00
71	T	1,2,3-Trichloropropane	10.000	10.159	-1.6	91	0.00
72	t	Bromobenzene	10.000	10.453	-4.5	101	0.00
73	t	n-propylbenzene	10.000	11.502	-15.0	99	0.00
74	t	2-Chlorotoluene	10.000	11.165	-11.6	100	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.611	-16.1	101	0.00
76	t	4-Chlorotoluene	10.000	11.384	-13.8	101	0.00
77	t	tert-Butylbenzene	10.000	11.071	-10.7	100	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.923	-19.2	101	0.00
79	t	sec-Butylbenzene	10.000	11.314	-13.1	101	0.00
80		Nitrobenzene	50.000	55.059	-10.1	96	0.00
81	t	p-Isopropyltoluene	10.000	11.697	-17.0	100	0.00
82	t	1,3-Dichlorobenzene	10.000	10.827	-8.3	101	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.799	-8.0	101	0.00
84	t	n-Butylbenzene	10.000	11.931	-19.3	101	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.602	-6.0	101	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.948	-9.5	97	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	12.216	-22.2	103	0.00
88	t	Hexachlorobutadiene	10.000	10.353	-3.5	98	0.00
89	t	Naphthalene	10.000	12.819	-28.2	101	0.00
90	t	1,2,3-Trichlorobenzene	10.000	12.047	-20.5	104	0.00

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

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