Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU020819\
Data File : VU029372.D
Acq On : 08 Feb 2019 10:27
Operator : JC/SP
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Feb 09 03:25:40 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U020519DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Fri Feb 08 10:52:46 2019
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	96	0.00
2	T	Dichlorodifluoromethane	10.000	9.589	4.1	93	0.00
3	t	Chloromethane	10.000	8.721	12.8	88	0.00
4	Rt	Vinyl Chloride	10.000	9.634	3.7	92	0.00
5	T	Bromomethane	10.000	10.272	-2.7	109	0.00
6	T	Chloroethane	10.000	9.850	1.5	96	0.00
7	T	Trichlorofluoromethane	10.000	10.705	-7.1	102	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	11.008	-10.1	105	0.00
9	Rt	1,1-Dichloroethene	10.000	10.750	-7.5	103	0.00
10	t	Iodomethane	10.000	10.088	-0.9	97	0.00
11	t	Allyl Chloride	10.000	9.603	4.0	91	0.00
12	t	Acrylonitrile	20.000	20.137	-0.7	97	0.00
13	T	Acetone	51.000	49.854	2.2	98	0.00
14	T	Carbon Disulfide	10.000	10.117	-1.2	97	0.00
15	RT	Methylene Chloride	10.000	9.692	3.1	102	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.724	-7.2	103	0.00
17	t	1,1-Dichloroethane	10.000	9.932	0.7	97	0.00
18	T	2-Butanone	50.000	48.920	2.2	86	0.00
19		Cyclohexane	10.000	7.761	22.4	79	0.00
20		Methylcyclohexane	10.000	10.222	-2.2	96	0.00
21	T	2,2-Dichloropropane	10.000	10.206	-2.1	99	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.624	-6.2	101	0.00
23	t	Diethyl Ether	10.000	9.968	0.3	96	0.00
24	t	tert-Butyl Alcohol	100.000	101.389	-1.4	99	0.00
25	t	Methyl tert-Butyl Ether	10.000	10.349	-3.5	99	0.00
26	t	Bromochloromethane	10.000	10.883	-8.8	104	0.00
27	t	Chloroform	10.000	10.415	-4.1	98	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.388	-3.9	98	0.00
29	T	1,1-Dichloropropene	10.000	9.922	0.8	94	0.00
30	RT	Carbon Tetrachloride	10.000	10.813	-8.1	101	0.00
31	t	Isopropyl Ether	10.000	9.310	6.9	87	0.00
32		Ethyl-t-butyl ether	10.000	9.609	3.9	91	0.00
33		Tert-Amyl methyl ether	10.000	10.118	-1.2	95	0.00
34	t	Propionitrile	50.000	52.230	-4.5	90	0.00
35	RT	Benzene	10.000	10.117	-1.2	95	0.00
36	RT	1,2-Dichloroethane	10.000	9.808	1.9	91	0.00
37	RT	Trichloroethene	10.000	11.131	-11.3	105	0.00
38	Rt	1,2-Dichloropropane	10.000	9.759	2.4	92	0.00
39	t	Methacrylonitrile	10.000	9.264	7.4	86	0.00
40	t	Methyl acrylate	10.000	9.903	1.0	92	0.00
41	t	Tetrahydrofuran	20.000	17.808	11.0	84	0.00
42	t	1-Chlorobutane	10.000	9.531	4.7	89	0.00
43	T	Dibromomethane	10.000	10.261	-2.6	99	0.00
44	T	Bromodichloromethane	10.000	10.175	-1.8	96	0.00
45	T	4-Methyl-2-Pentanone	50.000	47.133	5.7	86	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	19.688	1.6	88	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	20.647	-3.2	94	0.00
48	t	Ethyl methacrylate	10.000	10.359	-3.6	93	0.00
49	Rt	Toluene	10.000	10.516	-5.2	98	0.00
50	T	t-1,3-Dichloropropene	10.000	10.149	-1.5	93	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.117	-1.2	94	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.568	-5.7	98	0.00
53	t	1,3-Dichloropropane	10.000	10.311	-3.1	96	0.00
54	t	2-Hexanone	50.000	48.649	2.7	86	0.00
55	t	Dibromochloromethane	10.000	11.011	-10.1	102	0.00
56	T	1,2-Dibromoethane	10.000	10.356	-3.6	99	0.00
57	S	4-Bromofluorobenzene	1.000	1.133	-13.3	109	0.00
58	RT	Tetrachloroethene	10.000	11.419	-14.2	106	0.00
59	Rt	Chlorobenzene	10.000	11.053	-10.5	103	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	11.207	-12.1	105	0.00
61	t	Pentachloroethane	10.000	10.975	-9.7	101	0.00
62	t	Hexachloroethane	10.000	11.257	-12.6	103	0.00
63	Rt	Ethyl Benzene	10.000	10.555	-5.5	98	0.00
64	RT	m/p-Xylenes	20.000	21.785	-8.9	100	0.00
65	RT	o-Xylene	10.000	10.824	-8.2	99	0.00
66	RT	Styrene	10.000	10.869	-8.7	99	0.00
67	t	Bromoform	10.000	11.384	-13.8	104	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.113	-11.3	102	0.00
69	T	Isopropylbenzene	10.000	10.894	-8.9	100	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.586	-5.9	97	0.00
71	T	1,2,3-Trichloropropane	10.000	10.869	-8.7	96	0.00
72	t	Bromobenzene	10.000	11.566	-15.7	105	0.00
73	t	n-propylbenzene	10.000	11.387	-13.9	104	0.00
74	t	2-Chlorotoluene	10.000	11.203	-12.0	104	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.737	-7.4	99	0.00
76	t	4-Chlorotoluene	10.000	11.337	-13.4	103	0.00
77	t	tert-Butylbenzene	10.000	11.037	-10.4	101	0.00
78	t	1,2,4-Trimethylbenzene	10.000	10.808	-8.1	98	0.00
79	t	sec-Butylbenzene	10.000	10.870	-8.7	99	0.00
80		Nitrobenzene	50.000	53.028	-6.1	79	0.00
81	t	p-Isopropyltoluene	10.000	11.192	-11.9	101	0.00
82	t	1,3-Dichlorobenzene	10.000	11.604	-16.0	105	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.637	-16.4	105	0.00
84	t	n-Butylbenzene	10.000	10.961	-9.6	97	0.00
85	Rt	1,2-Dichlorobenzene	10.000	11.393	-13.9	103	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	9.526	4.7	88	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	12.307	-23.1	107	0.00
88	t	Hexachlorobutadiene	10.000	12.467	-24.7	113	0.00
89	t	Naphthalene	10.000	11.623	-16.2	98	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.805	-18.0	105	0.00

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

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