Evaluate Continuing Calibration Report

Data Path : W:\HPCHEM1\MSVOA_U\Data\VU051418\
Data File : VU023847.D
Acq On : 14 May 2018 10:48
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 15 04:14:09 2018
Quant Method : W:\HPCHEM1\MSVOA_U\METHOD\524U043018DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Thu May 03 10:23:54 2018
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	77	0.00
2	T	Dichlorodifluoromethane	10.000	10.690	-6.9	84	0.00
3	t	Chloromethane	10.000	10.565	-5.6	88	0.00
4	Rt	Vinyl Chloride	10.000	11.007	-10.1	87	0.00
5	T	Bromomethane	10.000	9.271	7.3	74	0.00
6	T	Chloroethane	10.000	11.407	-14.1	91	0.00
7	T	Trichlorofluoromethane	10.000	11.413	-14.1	89	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	11.213	-12.1	89	0.00
9	Rt	1,1-Dichloroethene	10.000	11.062	-10.6	87	0.00
10	t	Iodomethane	10.000	7.140	28.6	52	0.00
11	t	Allyl Chloride	10.000	11.371	-13.7	89	0.00
12	t	Acrylonitrile	20.000	25.607	-28.0	97	0.00
13	T	Acetone	50.000	143.120	-186.2#	240	-0.02
14	T	Carbon Disulfide	10.000	11.243	-12.4	90	0.00
15	RT	Methylene Chloride	10.000	10.493	-4.9	90	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.868	-8.7	87	0.00
17	t	1,1-Dichloroethane	10.000	10.580	-5.8	84	0.00
18	T	2-Butanone	50.000	87.262	-74.5#	137	-0.02
19		Cyclohexane	10.000	9.499	5.0	75	0.00
20		Methylcyclohexane	10.000	10.264	-2.6	76	0.00
21	T	2,2-Dichloropropane	10.000	10.057	-0.6	80	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.653	-6.5	84	0.00
23	t	Diethyl Ether	10.000	11.370	-13.7	89	0.00
24	t	tert-Butyl Alcohol	100.000	147.071	-47.1#	118	-0.04
25	t	Methyl tert-Butyl Ether	10.000	11.130	-11.3	87	0.00
26	t	Bromochloromethane	10.000	11.540	-15.4	93	0.00
27	t	Chloroform	10.000	10.579	-5.8	84	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.583	-5.8	84	0.00
29	T	1,1-Dichloropropene	10.000	10.501	-5.0	82	0.00
30	RT	Carbon Tetrachloride	10.000	10.980	-9.8	85	0.00
31	t	Isopropyl Ether	10.000	9.639	3.6	76	0.00
32		Ethyl-t-butyl ether	10.000	9.550	4.5	74	0.00
33		Tert-Amyl methyl ether	10.000	9.664	3.4	74	0.00
34	t	Propionitrile	50.000	56.948	-13.9	89	-0.03
35	RT	Benzene	10.000	9.937	0.6	75	0.00
36	RT	1,2-Dichloroethane	10.000	9.585	4.1	75	0.00
37	RT	Trichloroethene	10.000	9.445	5.5	74	0.00
38	Rt	1,2-Dichloropropane	10.000	9.796	2.0	75	0.00
39	t	Methacrylonitrile	10.000	9.884	1.2	80	-0.01
40	t	Methyl acrylate	10.000	9.854	1.5	75	0.00
41	t	Tetrahydrofuran	20.000	20.079	-0.4	86	-0.02
42	t	1-Chlorobutane	10.000	10.315	-3.1	81	0.00
43	T	Dibromomethane	10.000	10.384	-3.8	81	0.00
44	T	Bromodichloromethane	10.000	10.329	-3.3	80	0.00
45	T	4-Methyl-2-Pentanone	50.000	56.800	-13.6	83	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	25.666	-28.3	101	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	22.001	-10.0	80	0.00
48	t	Ethyl methacrylate	10.000	10.775	-7.8	76	0.00
49	Rt	Toluene	10.000	10.478	-4.8	77	0.00
50	T	t-1,3-Dichloropropene	10.000	10.870	-8.7	78	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.462	-4.6	77	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.129	-1.3	78	0.00
53	t	1,3-Dichloropropane	10.000	10.212	-2.1	77	0.00
54	t	2-Hexanone	50.000	78.682	-57.4#	112	0.00
55	t	Dibromochloromethane	10.000	10.877	-8.8	80	0.00
56	T	1,2-Dibromoethane	10.000	10.393	-3.9	79	0.00
57	S	4-Bromofluorobenzene	1.000	1.144	-14.4	82	0.00
58	RT	Tetrachloroethene	10.000	10.064	-0.6	76	0.00
59	Rt	Chlorobenzene	10.000	10.413	-4.1	78	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.457	-4.6	79	0.00
61	t	Pentachloroethane	10.000	11.434	-14.3	91	0.00
62	t	Hexachloroethane	10.000	12.016	-20.2	92	0.00
63	Rt	Ethyl Benzene	10.000	10.950	-9.5	78	0.00
64	RT	m/p-Xylenes	20.000	22.425	-12.1	80	0.00
65	RT	o-Xylene	10.000	11.227	-12.3	80	0.00
66	RT	Styrene	10.000	11.568	-15.7	80	0.00
67	t	Bromoform	10.000	11.806	-18.1	87	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.243	-24.3	93	0.00
69	T	Isopropylbenzene	10.000	11.195	-12.0	80	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	11.187	-11.9	85	0.00
71	T	1,2,3-Trichloropropane	10.000	9.806	1.9	77	0.00
72	t	Bromobenzene	10.000	11.153	-11.5	82	0.00
73	t	n-propylbenzene	10.000	11.826	-18.3	84	0.00
74	t	2-Chlorotoluene	10.000	11.506	-15.1	84	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.903	-19.0	84	0.00
76	t	4-Chlorotoluene	10.000	11.624	-16.2	84	0.00
77	t	tert-Butylbenzene	10.000	11.941	-19.4	86	0.00
78	t	1,2,4-Trimethylbenzene	10.000	12.096	-21.0	85	0.00
79	t	sec-Butylbenzene	10.000	11.891	-18.9	86	0.00
80		Nitrobenzene	50.000	46.179	7.6	77	0.00
81	t	p-Isopropyltoluene	10.000	12.529	-25.3	89	0.00
82	t	1,3-Dichlorobenzene	10.000	11.480	-14.8	87	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.666	-16.7	87	0.00
84	t	n-Butylbenzene	10.000	12.281	-22.8	87	0.00
85	Rt	1,2-Dichlorobenzene	10.000	11.603	-16.0	88	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	13.113	-31.1#	101	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.923	-19.2	89	0.00
88	t	Hexachlorobutadiene	10.000	12.448	-24.5	96	0.00
89	t	Naphthalene	10.000	11.893	-18.9	81	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.755	-17.6	88	0.00

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

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