Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU020719\
Data File : VU029364.D
Acq On : 07 Feb 2019 09:54
Operator : JC/SP
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Feb 08 12:57:45 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	85	0.00
2	T	Dichlorodifluoromethane	10.000	9.552	4.5	82	0.00
3	t	Chloromethane	10.000	9.069	9.3	80	0.00
4	Rt	Vinyl Chloride	10.000	9.570	4.3	80	0.00
5	T	Bromomethane	10.000	9.758	2.4	91	0.00
6	T	Chloroethane	10.000	9.843	1.6	84	0.00
7	T	Trichlorofluoromethane	10.000	10.580	-5.8	89	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	10.884	-8.8	91	0.00
9	Rt	1,1-Dichloroethene	10.000	10.591	-5.9	90	0.00
10	t	Iodomethane	10.000	11.187	-11.9	95	0.00
11	t	Allyl Chloride	10.000	9.768	2.3	82	0.00
12	t	Acrylonitrile	20.000	19.653	1.7	84	0.00
13	T	Acetone	51.000	49.189	3.6	85	0.00
14	T	Carbon Disulfide	10.000	10.104	-1.0	86	0.00
15	RT	Methylene Chloride	10.000	9.500	5.0	88	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.421	-4.2	88	0.00
17	t	1,1-Dichloroethane	10.000	9.671	3.3	83	0.00
18	T	2-Butanone	50.000	48.540	2.9	76	0.00
19		Cyclohexane	10.000	9.014	9.9	81	0.00
20		Methylcyclohexane	10.000	10.137	-1.4	84	0.00
21	T	2,2-Dichloropropane	10.000	9.822	1.8	84	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.484	-4.8	87	0.00
23	t	Diethyl Ether	10.000	9.861	1.4	83	0.00
24	t	tert-Butyl Alcohol	100.000	97.129	2.9	83	0.00
25	t	Methyl tert-Butyl Ether	10.000	10.235	-2.3	86	0.00
26	t	Bromochloromethane	10.000	10.598	-6.0	89	0.00
27	t	Chloroform	10.000	10.253	-2.5	85	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.336	-3.4	86	0.00
29	T	1,1-Dichloropropene	10.000	9.811	1.9	82	0.00
30	RT	Carbon Tetrachloride	10.000	10.625	-6.3	88	0.00
31	t	Isopropyl Ether	10.000	9.423	5.8	78	0.00
32		Ethyl-t-butyl ether	10.000	9.579	4.2	80	0.00
33		Tert-Amyl methyl ether	10.000	10.091	-0.9	83	0.00
34	t	Propionitrile	50.000	49.526	0.9	75	0.00
35	RT	Benzene	10.000	10.158	-1.6	84	0.00
36	RT	1,2-Dichloroethane	10.000	9.616	3.8	79	0.00
37	RT	Trichloroethene	10.000	10.934	-9.3	91	0.00
38	Rt	1,2-Dichloropropane	10.000	9.921	0.8	82	0.00
39	t	Methacrylonitrile	10.000	9.098	9.0	75	0.00
40	t	Methyl acrylate	10.000	9.595	4.0	78	0.00
41	t	Tetrahydrofuran	20.000	17.699	11.5	74	0.00
42	t	1-Chlorobutane	10.000	9.727	2.7	80	0.00
43	T	Dibromomethane	10.000	10.113	-1.1	86	0.00
44	T	Bromodichloromethane	10.000	10.181	-1.8	85	0.00
45	T	4-Methyl-2-Pentanone	50.000	47.580	4.8	76	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	20.142	-0.7	79	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	20.449	-2.2	82	0.00
48	t	Ethyl methacrylate	10.000	10.094	-0.9	80	0.00
49	Rt	Toluene	10.000	10.403	-4.0	86	0.00
50	T	t-1,3-Dichloropropene	10.000	9.919	0.8	80	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.031	-0.3	82	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.339	-3.4	85	0.00
53	t	1,3-Dichloropropane	10.000	10.134	-1.3	83	0.00
54	t	2-Hexanone	50.000	48.032	3.9	75	0.00
55	t	Dibromochloromethane	10.000	10.620	-6.2	87	0.00
56	T	1,2-Dibromoethane	10.000	10.282	-2.8	86	0.00
57	S	4-Bromofluorobenzene	1.000	1.130	-13.0	96	0.00
58	RT	Tetrachloroethene	10.000	11.166	-11.7	92	0.00
59	Rt	Chlorobenzene	10.000	10.661	-6.6	88	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.796	-8.0	89	0.00
61	t	Pentachloroethane	10.000	10.935	-9.4	89	0.00
62	t	Hexachloroethane	10.000	10.944	-9.4	88	0.00
63	Rt	Ethyl Benzene	10.000	10.333	-3.3	85	0.00
64	RT	m/p-Xylenes	20.000	21.292	-6.5	86	0.00
65	RT	o-Xylene	10.000	10.669	-6.7	86	0.00
66	RT	Styrene	10.000	10.626	-6.3	85	0.00
67	t	Bromoform	10.000	11.108	-11.1	89	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.110	-11.0	90	0.00
69	T	Isopropylbenzene	10.000	10.609	-6.1	86	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.196	-2.0	82	0.00
71	T	1,2,3-Trichloropropane	10.000	10.722	-7.2	83	0.00
72	t	Bromobenzene	10.000	11.256	-12.6	90	0.00
73	t	n-propylbenzene	10.000	11.046	-10.5	89	0.00
74	t	2-Chlorotoluene	10.000	11.018	-10.2	90	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.554	-5.5	86	0.00
76	t	4-Chlorotoluene	10.000	11.021	-10.2	89	0.00
77	t	tert-Butylbenzene	10.000	10.763	-7.6	87	0.00
78	t	1,2,4-Trimethylbenzene	10.000	10.625	-6.3	85	0.00
79	t	sec-Butylbenzene	10.000	10.596	-6.0	85	0.00
80		Nitrobenzene	50.000	50.406	-0.8	67	0.00
81	t	p-Isopropyltoluene	10.000	10.871	-8.7	86	0.00
82	t	1,3-Dichlorobenzene	10.000	11.203	-12.0	90	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.126	-11.3	89	0.00
84	t	n-Butylbenzene	10.000	10.604	-6.0	83	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.927	-9.3	87	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	9.959	0.4	81	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.630	-16.3	89	0.00
88	t	Hexachlorobutadiene	10.000	12.093	-20.9	97	0.00
89	t	Naphthalene	10.000	11.233	-12.3	83	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.455	-14.6	89	0.00

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

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