Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090718\
Data File : VU026610.D
Acq On : 07 Sep 2018 09:02
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Sep 07 12:16:10 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U082018DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Mon Aug 20 13:35:41 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	72	0.00
2	T	Dichlorodifluoromethane	10.000	7.910	20.9	60	0.00
3	t	Chloromethane	10.000	11.111	-11.1	85	0.00
4	Rt	Vinyl Chloride	10.000	10.972	-9.7	82	0.00
5	T	Bromomethane	10.000	9.319	6.8	69	0.00
6	T	Chloroethane	10.000	11.291	-12.9	87	0.00
7	T	Trichlorofluoromethane	10.000	9.423	5.8	72	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.990	0.1	75	0.00
9	Rt	1,1-Dichloroethene	10.000	10.117	-1.2	77	0.00
10	t	Iodomethane	10.000	5.982	40.2#	43	0.00
11	t	Allyl Chloride	10.000	11.601	-16.0	87	0.00
12	t	Acrylonitrile	20.000	24.930	-24.6	98	0.01
13	T	Acetone	50.000	109.788	-119.6#	151	0.03
14	T	Carbon Disulfide	10.000	10.211	-2.1	78	0.00
15	RT	Methylene Chloride	10.000	11.376	-13.8	81	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.768	2.3	77	0.00
17	t	1,1-Dichloroethane	10.000	10.756	-7.6	82	0.00
18	T	2-Butanone	50.000	61.810	-23.6	91	0.02
19		Cyclohexane	10.000	9.867	1.3	67	0.00
20		Methylcyclohexane	10.000	9.832	1.7	65	0.00
21	T	2,2-Dichloropropane	10.000	7.836	21.6	59	0.00
22	RT	cis-1,2-Dichloroethene	10.000	8.225	17.8	60	0.00
23	t	Diethyl Ether	10.000	11.212	-12.1	85	0.00
24	t	tert-Butyl Alcohol	100.000	99.561	0.4	80	0.08
25	t	Methyl tert-Butyl Ether	10.000	10.424	-4.2	79	0.00
26	t	Bromochloromethane	10.000	7.740	22.6	56	0.00
27	t	Chloroform	10.000	7.974	20.3	60	0.00
28	RT	1,1,1-Trichloroethane	10.000	7.688	23.1	57	0.00
29	T	1,1-Dichloropropene	10.000	8.881	11.2	62	0.00
30	RT	Carbon Tetrachloride	10.000	7.376	26.2	54	0.00
31	t	Isopropyl Ether	10.000	9.336	6.6	69	0.00
32		Ethyl-t-butyl ether	10.000	9.149	8.5	66	0.00
33		Tert-Amyl methyl ether	10.000	9.201	8.0	62	0.00
34	t	Propionitrile	50.000	52.033	-4.1	77	0.03
35	RT	Benzene	10.000	8.840	11.6	65	0.00
36	RT	1,2-Dichloroethane	10.000	8.175	18.2	61	0.00
37	RT	Trichloroethene	10.000	8.165	18.4	61	0.00
38	Rt	1,2-Dichloropropane	10.000	9.160	8.4	66	0.00
39	t	Methacrylonitrile	10.000	9.380	6.2	66	0.00
40	t	Methyl acrylate	10.000	10.096	-1.0	72	0.00
41	t	Tetrahydrofuran	20.000	18.336	8.3	74	0.01
42	t	1-Chlorobutane	10.000	9.184	8.2	63	0.00
43	T	Dibromomethane	10.000	8.356	16.4	61	0.00
44	T	Bromodichloromethane	10.000	8.164	18.4	61	0.00
45	T	4-Methyl-2-Pentanone	50.000	52.138	-4.3	71	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	15.052	24.7	55	0.00

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		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
47	t	Methyl methacrylate	20.000	19.879	0.6	65	0.00
48	t	Ethyl methacrylate	10.000	10.432	-4.3	67	0.00
49	Rt	Toluene	10.000	9.232	7.7	62	0.00
50	T	t-1,3-Dichloropropene	10.000	8.853	11.5	62	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.072	9.3	64	0.00
52	RT	1,1,2-Trichloroethane	10.000	8.354	16.5	62	0.00
53	t	1,3-Dichloropropane	10.000	8.759	12.4	63	0.00
54	t	2-Hexanone	50.000	61.674	-23.3	84	0.00
55	t	Dibromochloromethane	10.000	7.714	22.9	56	0.00
56	T	1,2-Dibromoethane	10.000	8.116	18.8	59	0.00
57	S	4-Bromofluorobenzene	1.000	1.010	-1.0	69	0.00
58	RT	Tetrachloroethene	10.000	7.848	21.5	57	0.00
59	Rt	Chlorobenzene	10.000	8.683	13.2	61	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	8.051	19.5	59	0.00
61	t	Pentachloroethane	10.000	7.800	22.0	58	0.00
62	t	Hexachloroethane	10.000	8.087	19.1	56	0.00
63	Rt	Ethyl Benzene	10.000	9.540	4.6	62	0.00
64	RT	m/p-Xylenes	20.000	18.980	5.1	60	0.00
65	RT	o-Xylene	10.000	9.482	5.2	62	0.00
66	RT	Styrene	10.000	9.374	6.3	60	0.00
67	t	Bromoform	10.000	7.873	21.3	56	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.946	5.4	68	0.00
69	T	Isopropylbenzene	10.000	9.385	6.2	60	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	8.747	12.5	64	0.00
71	T	1,2,3-Trichloropropane	10.000	9.231	7.7	66	0.00
72	t	Bromobenzene	10.000	8.295	17.1	57	0.00
73	t	n-propylbenzene	10.000	9.158	8.4	57	0.00
74	t	2-Chlorotoluene	10.000	8.896	11.0	59	0.00
75	t	1,3,5-Trimethylbenzene	10.000	9.322	6.8	59	0.00
76	t	4-Chlorotoluene	10.000	8.912	10.9	59	0.00
77	t	tert-Butylbenzene	10.000	9.126	8.7	59	0.00
78	t	1,2,4-Trimethylbenzene	10.000	9.238	7.6	58	0.00
79	t	sec-Butylbenzene	10.000	9.274	7.3	59	0.00
80		Nitrobenzene	50.000	42.356	15.3	57	0.00
81	t	p-Isopropyltoluene	10.000	9.207	7.9	57	0.00
82	t	1,3-Dichlorobenzene	10.000	7.944	20.6	55	0.00
83	Rt	1,4-Dichlorobenzene	10.000	7.898	21.0	55	0.00
84	t	n-Butylbenzene	10.000	9.017	9.8	57	0.00
85	Rt	1,2-Dichlorobenzene	10.000	8.101	19.0	57	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	7.858	21.4	58	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	8.115	18.8	56	0.00
88	t	Hexachlorobutadiene	10.000	8.097	19.0	57	0.00
89	t	Naphthalene	10.000	8.993	10.1	57	0.00
90	t	1,2,3-Trichlorobenzene	10.000	8.281	17.2	55	0.00

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

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