Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX080218\
Data File : VX003819.D
Acq On : 02 Aug 2018 12:03
Operator : JC/MD
Sample : VSTDCCC020
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_X
LabSampleId : VSTDCCC020

Quant Time: Aug 03 03:25:57 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X072718W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
QLast Update : Fri Jul 27 16:17:43 2018
Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 30% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	Bromochloromethane	1.000	1.000	0.0	99	0.00
2	M	Dichlorodifluoromethane	1.585	2.085	-31.5#	144	0.00
3	M	Chloromethane	2.133	2.572	-20.6	132	0.00
4	M	Vinyl Chloride	2.243	1.964	12.4	96	0.00
5	M	Bromomethane	1.341	1.045	22.1	91	0.00
6	M	Chloroethane	1.468	1.240	15.5	92	0.00
7	M	Trichlorofluoromethane	3.017	2.633	12.7	97	0.00
8	T	Diethyl Ether	1.293	1.083	16.2	93	0.00
9		1,1,2-Trichlorotrifluoroeth	1.829	1.612	11.9	96	0.00
10	M	1,1-Dichloroethene	1.720	1.570	8.7	101	0.00
11		Methyl Iodide	2.384	1.654	30.6#	77	0.00
12		Methyl Acetate	3.182	0.958	69.9#	32	0.43
13	M	Acrolein	0.248	0.293	-18.1	114	0.00
14	M	Acrylonitrile	1.254	0.991	21.0	85	0.00
15	M	Acetone	0.340	0.293	13.8	93	0.00
16	M	Carbon Disulfide	4.776	4.169	12.7	98	0.00
17		Allyl chloride	3.560	2.861	19.6	88	0.00
18	M	Methylene Chloride	2.037	1.854	9.0	98	0.00
19	M	trans-1,2-Dichloroethene	1.859	1.565	15.8	91	0.00
20	T	Diisopropyl ether	6.537	5.420	17.1	89	0.00
21	M	1,1-Dichloroethane	3.701	3.110	16.0	91	0.00
22	M	cis-1,2-Dichloroethene	1.877	1.926	-2.6	105	0.00
23	M	tert-Butyl Alcohol	0.527	0.416	21.1	84	-0.01
24	M	Methyl tert-Butyl Ether	6.169	5.218	15.4	93	-0.01
25	M	Chloroform	3.205	3.103	3.2	98	0.00
26		Cyclohexane	2.780	2.673	3.8	100	-0.01
27	s	1,2-Dichloroethane-d4	2.202	2.001	9.1	95	0.00
28	I	1,4-Difluorobenzene	1.000	1.000	0.0	92	0.00
29		1,1-Dichloropropene	0.443	0.460	-3.8	97	0.00
30	M	2-Butanone	0.283	0.294	-3.9	93	-0.01
31		2,2-Dichloropropane	0.429	0.489	-14.0	104	-0.01
32	M	1,1,1-Trichloroethane	0.507	0.522	-3.0	96	0.00
33	M	Carbon Tetrachloride	0.440	0.464	-5.5	97	-0.01
34	M	Benzene	1.361	1.433	-5.3	96	-0.01
35		Methacrylonitrile	0.294	0.312	-6.1	98	0.00
36	M	1,2-Dichloroethane	0.489	0.478	2.2	91	0.00
37	M	Trichloroethene	0.355	0.357	-0.6	93	0.00
38		Methylcyclohexane	0.503	0.494	1.8	93	0.00
39	M	1,2-Dichloropropane	0.360	0.357	0.8	92	0.00
40		Dibromomethane	0.233	0.233	0.0	92	0.00
41	M	Bromodichloromethane	0.445	0.429	3.6	92	0.00
42	M	Vinyl Acetate	1.087	0.913	16.0	84	-0.01
43		Ethyl Acetate	0.528	0.566	-7.2	97	0.00
44		Isopropyl Acetate	0.811	0.760	6.3	88	0.00
45	T	1,4-Dioxane	0.009	0.009#	0.0	87	0.00

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Quant Time: Aug 03 03:25:57 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\624X072718W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
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Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
46		Methyl methacrylate	0.418	0.398	4.8	89	0.00
47		n-amyl Acetate	0.765	0.672	12.2	89	0.00
48	M	t-1,3-Dichloropropene	0.540	0.537	0.6	95	0.00
49	T	cis-1,3-Dichloropropene	0.506	0.488	3.6	98	0.00
50	M	1,1,2-Trichloroethane	0.370	0.357	3.5	94	0.00
51		Ethyl methacrylate	0.556	0.501	9.9	92	0.00
52		1,3-Dichloropropane	0.615	0.590	4.1	94	0.00
53	M	Dibromochloromethane	0.363	0.353	2.8	98	0.00
54	M	1,2-Dibromoethane	0.386	0.372	3.6	95	0.00
55	M	2-Chloroethyl vinyl ether	0.287	0.264	8.0	88	0.00
56	M	Bromoform	0.282	0.269	4.6	100	0.00
57	I	Chlorobenzene-d5	1.000	1.000	0.0	94	0.00
58	M	4-Methyl-2-Pentanone	0.566	0.549	3.0	86	0.00
59	M	2-Hexanone	0.451	0.417	7.5	86	0.00
60	S	4-Bromofluorobenzene	0.483	0.485	-0.4	96	0.00
61	M	Tetrachloroethene	0.349	0.362	-3.7	100	0.00
62	M	Toluene	1.557	1.559	-0.1	92	0.00
63	S	Toluene-d8	1.351	1.339	0.9	90	0.00
64	M	Chlorobenzene	1.006	1.017	-1.1	94	0.00
65		1,1,1,2-Tetrachloroethane	0.356	0.364	-2.2	96	0.00
66	M	Ethyl Benzene	1.713	1.680	1.9	93	0.00
67	M	m/p-Xylenes	0.660	0.666	-0.9	95	0.00
68	M	o-Xylene	0.631	0.635	-0.6	95	0.00
69	M	Styrene	1.069	1.061	0.7	94	0.00
70		Isopropylbenzene	1.738	1.701	2.1	94	0.00
71	M	1,1,2,2-Tetrachloroethane	0.621	0.595	4.2	91	0.00
72		1,2,3-Trichloropropane	0.516	0.538	-4.3	99	0.00
73		Bromobenzene	0.468	0.477	-1.9	100	0.00
74		n-propylbenzene	2.048	2.002	2.2	94	0.00
75		2-Chlorotoluene	1.207	1.197	0.8	94	0.00
76		1,3,5-Trimethylbenzene	1.475	1.454	1.4	95	0.00
77		t-1,4-Dichloro-2-butene	0.176	0.171	2.8	98	0.00
78		4-Chlorotoluene	1.412	1.395	1.2	95	0.00
79		tert-butylbenzene	1.518	1.517	0.1	96	0.00
80		1,2,4-Trimethylbenzene	1.510	1.493	1.1	94	0.00
81		sec-Butylbenzene	1.718	1.699	1.1	96	0.00
82		p-Isopropyltoluene	1.518	1.517	0.1	96	0.00
83	M	1,3-Dichlorobenzene	0.856	0.886	-3.5	99	0.00
84	M	1,4-Dichlorobenzene	0.861	0.903	-4.9	100	0.00
85		n-Butylbenzene	1.347	1.320	2.0	98	0.00
86	T	Hexachloroethane	0.245	0.224	8.6	90	0.00
87	M	1,2-Dichlorobenzene	0.859	0.889	-3.5	99	0.00
88		1,2-Dibromo-3-Chloropropane	0.131	0.124	5.3	95	0.00
89		1,2,4-Trichlorobenzene	0.603	0.661	-9.6	108	0.00
90		Hexachlorobutadiene	0.273	0.304	-11.4	106	0.00

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91 M	Naphthalene	1.843	1.921	-4.2	100	0.00
92	1,2,3-Trichlorobenzene	0.614	0.675	-9.9	105	0.00

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