Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081122\
Data File : VN073838.D
Acq On : 11 Aug 2022 20:35
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
LabSampleId :
VSTDCCC050

Quant Time: Aug 12 06:10:45 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 10 05:09:32 2022
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	Pentafluorobenzene	1.000	1.000	0.0	102	0.00
2	T	Dichlorodifluoromethane	0.485	0.508	-4.7	114	0.00
3	P	Chloromethane	0.699	0.629	10.0	104	0.00
4	C	Vinyl Chloride	0.848	0.749	11.7#	96	0.00
5	T	Bromomethane	0.492	0.422	14.2	94	-0.01
6	T	Chloroethane	0.592	0.514	13.2	102	-0.02
7	T	Trichlorofluoromethane	0.874	1.038	-18.8	124	-0.01
8	T	Diethyl Ether	0.328	0.328	0.0	108	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.484	0.481	0.6	110	0.00
10	T	Methyl Iodide	0.492	0.560	-13.8	107	-0.01
11	T	Tert butyl alcohol	0.131	0.150	-14.5	116	0.01
12	CM	1,1-Dichloroethene	0.474	0.476	-0.4#	111	0.00
13	T	Acrolein	0.130	0.127	2.3	95	0.00
14	T	Allyl chloride	0.758	0.752	0.8	105	-0.01
15	T	Acrylonitrile	0.364	0.392	-7.7	111	0.00
16	T	Acetone	0.327	0.331	-1.2	115	0.00
17	T	Carbon Disulfide	1.258	1.258	0.0	109	0.00
18	T	Methyl Acetate	0.814	0.858	-5.4	115	0.00
19	T	Methyl tert-butyl Ether	1.621	1.735	-7.0	110	0.00
20	T	Methylene Chloride	0.691	0.596	13.7	111	0.00
21	T	trans-1,2-Dichloroethene	0.514	0.526	-2.3	109	-0.01
22	T	Diisopropyl ether	1.636	1.670	-2.1	106	0.00
23	T	Vinyl Acetate	1.372	1.501	-9.4	108	0.00
24	P	1,1-Dichloroethane	1.007	0.987	2.0	109	0.00
25	T	2-Butanone	0.508	0.534	-5.1	110	0.00
26	T	2,2-Dichloropropane	0.774	0.592	23.5	83	0.00
27	T	cis-1,2-Dichloroethene	0.633	0.637	-0.6	107	0.00
28	T	Bromochloromethane	0.423	0.393	7.1	96	0.00
29	T	Tetrahydrofuran	0.321	0.348	-8.4	108	0.00
30	C	Chloroform	1.049	1.035	1.3#	107	0.00
31	T	Cyclohexane	0.907	0.847	6.6	103	0.00
32	T	1,1,1-Trichloroethane	0.871	0.895	-2.8	109	0.00
33	S	1,2-Dichloroethane-d4	0.609	0.578	5.1	103	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	100	0.00
35	S	Dibromofluoromethane	0.314	0.325	-3.5	107	0.00
36	T	1,1-Dichloropropene	0.447	0.466	-4.3	105	0.00
37	T	Ethyl Acetate	0.589	0.662	-12.4	112	0.00
38	T	Carbon Tetrachloride	0.454	0.489	-7.7	111	0.00
39	T	Methylcyclohexane	0.535	0.573	-7.1	103	0.00
40	TM	Benzene	1.453	1.506	-3.6	106	0.00
41	T	Methacrylonitrile	0.267	0.294	-10.1	105	0.00
42	TM	1,2-Dichloroethane	0.490	0.506	-3.3	108	0.00
43	T	Isopropyl Acetate	0.796	0.893	-12.2	108	0.00
44	TM	Trichloroethene	0.368	0.387	-5.2	107	0.00
45	C	1,2-Dichloropropane	0.363	0.378	-4.1#	107	0.00
46	T	Dibromomethane	0.252	0.268	-6.3	108	0.00
47	T	Bromodichloromethane	0.484	0.515	-6.4	109	0.00
48	T	Methyl methacrylate	0.348	0.401	-15.2	111	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	T	1,4-Dioxane	0.010	0.012	-20.0	116	0.00
50	S	Toluene-d8	1.236	1.206	2.4	98	0.00
51	T	4-Methyl-2-Pentanone	0.571	0.648	-13.5	108	0.00
52	CM	Toluene	0.888	0.964	-8.6#	105	0.00
53	T	t-1,3-Dichloropropene	0.478	0.524	-9.6	104	0.00
54	T	cis-1,3-Dichloropropene	0.539	0.573	-6.3	102	0.00
55	T	1,1,2-Trichloroethane	0.370	0.401	-8.4	108	0.00
56	T	Ethyl methacrylate	0.510	0.629	-23.3	108	0.00
57	T	1,3-Dichloropropane	0.628	0.669	-6.5	107	0.00
58	T	2-Chloroethyl Vinyl ether	0.229	0.281	-22.7	109	0.00
59	T	2-Hexanone	0.433	0.513	-18.5	109	0.00
60	T	Dibromochloromethane	0.368	0.424	-15.2	111	0.00
61	T	1,2-Dibromoethane	0.381	0.419	-10.0	108	0.00
62	S	4-Bromofluorobenzene	0.382	0.421	-10.2	104	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	99	0.00
64	T	Tetrachloroethene	0.397	0.393	1.0	105	0.00
65	PM	Chlorobenzene	1.062	1.103	-3.9	107	0.00
66	T	1,1,1,2-Tetrachloroethane	0.381	0.419	-10.0	110	0.00
67	C	Ethyl Benzene	1.792	1.942	-8.4#	104	0.00
68	T	m/p-Xylenes	0.683	0.763	-11.7	106	0.00
69	T	o-Xylene	0.666	0.754	-13.2	105	0.00
70	T	Styrene	1.066	1.259	-18.1	107	0.00
71	P	Bromoform	0.314	0.366	-16.6	112	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
73	T	Isopropylbenzene	3.660	3.763	-2.8	105	0.00
74	T	N-amyl acetate	1.383	1.526	-10.3	111	0.00
75	P	1,1,2,2-Tetrachloroethane	1.412	1.320	6.5	108	0.00
76	T	1,2,3-Trichloropropane	1.135	1.009	11.1	109	0.00
77	T	Bromobenzene	0.953	0.959	-0.6	108	0.00
78	T	n-propylbenzene	4.118	4.288	-4.1	104	0.00
79	T	2-Chlorotoluene	2.570	2.582	-0.5	106	0.00
80	T	1,3,5-Trimethylbenzene	3.012	3.193	-6.0	107	0.00
81	T	trans-1,4-Dichloro-2-butene	0.386	0.383	0.8	103	0.00
82	T	4-Chlorotoluene	2.436	2.472	-1.5	105	0.00
83	T	tert-Butylbenzene	2.668	2.907	-9.0	108	0.00
84	T	1,2,4-Trimethylbenzene	2.962	3.177	-7.3	106	0.00
85	T	sec-Butylbenzene	3.687	3.919	-6.3	104	0.00
86	T	p-Isopropyltoluene	2.907	3.298	-13.5	106	0.00
87	T	1,3-Dichlorobenzene	1.740	1.734	0.3	108	0.00
88	T	1,4-Dichlorobenzene	1.741	1.690	2.9	106	0.00
89	T	n-Butylbenzene	2.365	2.566	-8.5	104	0.00
90	T	Hexachloroethane	0.577	0.596	-3.3	109	0.00
91	T	1,2-Dichlorobenzene	1.763	1.760	0.2	108	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.284	0.303	-6.7	114	0.00
93	T	1,2,4-Trichlorobenzene	0.876	0.966	-10.3	106	0.00
94	T	Hexachlorobutadiene	0.476	0.511	-7.4	113	0.00
95	T	Naphthalene	2.841	3.358	-18.2	107	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.921	1.036	-12.5	110	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6