Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120325\
Data File : VN088435.D
Acq On : 03 Dec 2025 17:18
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
LabSampleId :
VSTDCCC050

Quant Time: Dec 04 00:23:51 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
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	103	0.00
2	T	Dichlorodifluoromethane	0.751	0.834	-11.1	116	0.00
3	P	Chloromethane	0.807	0.793	1.7	102	0.00
4	C	Vinyl Chloride	0.821	0.866	-5.5#	110	0.00
5	T	Bromomethane	0.395	0.388	1.8	101	0.00
6	T	Chloroethane	0.508	0.540	-6.3	110	0.00
7	T	Trichlorofluoromethane	1.153	1.295	-12.3	122	0.00
8	T	Diethyl Ether	0.433	0.451	-4.2	104	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.614	0.675	-9.9	123	0.00
10	T	Methyl Iodide	0.608	0.577	5.1	90	0.00
11	T	Tert butyl alcohol	0.150	0.144	4.0	92	0.00
12	CM	1,1-Dichloroethene	0.610	0.640	-4.9#	116	0.00
13	T	Acrolein	0.144	0.186	-29.2#	127	0.00
14	T	Allyl chloride	1.131	1.171	-3.5	105	0.00
15	T	Acrylonitrile	0.434	0.459	-5.8	101	0.00
16	T	Acetone	0.341	0.467	-37.0#	143	0.00
17	T	Carbon Disulfide	1.993	2.020	-1.4	105	0.00
18	T	Methyl Acetate	1.014	1.105	-9.0	108	0.00
19	T	Methyl tert-butyl Ether	2.285	2.472	-8.2	101	0.00
20	T	Methylene Chloride	0.729	0.742	-1.8	101	0.00
21	T	trans-1,2-Dichloroethene	0.676	0.722	-6.8	108	0.00
22	T	Diisopropyl ether	2.414	2.672	-10.7	107	0.00
23	T	Vinyl Acetate	1.945	2.432	-25.0#	117	0.00
24	P	1,1-Dichloroethane	1.367	1.460	-6.8	105	0.00
25	T	2-Butanone	0.548	0.663	-21.0	115	0.00
26	T	2,2-Dichloropropane	1.180	1.229	-4.2	105	0.00
27	T	cis-1,2-Dichloroethene	0.785	0.835	-6.4	104	0.00
28	T	Bromochloromethane	0.681	0.725	-6.5	101	0.00
29	T	Tetrahydrofuran	0.393	0.428	-8.9	103	0.00
30	C	Chloroform	1.343	1.487	-10.7#	110	0.00
31	T	Cyclohexane	1.239	1.264	-2.0	115	0.00
32	T	1,1,1-Trichloroethane	1.181	1.303	-10.3	116	0.00
33	S	1,2-Dichloroethane-d4	0.942	1.003	-6.5	103	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	99	0.00
35	S	Dibromofluoromethane	0.346	0.369	-6.6	104	0.00
36	T	1,1-Dichloropropene	0.535	0.573	-7.1	115	0.00
37	T	Ethyl Acetate	0.703	0.761	-8.3	104	0.00
38	T	Carbon Tetrachloride	0.532	0.610	-14.7	121	0.00
39	T	Methylcyclohexane	0.572	0.637	-11.4	119	0.00
40	TM	Benzene	1.570	1.681	-7.1	107	0.00
41	T	Methacrylonitrile	0.348	0.382	-9.8	106	0.00
42	TM	1,2-Dichloroethane	0.599	0.673	-12.4	108	0.00
43	T	Isopropyl Acetate	1.074	1.151	-7.2	102	0.00
44	TM	Trichloroethene	0.371	0.378	-1.9	104	0.00
45	C	1,2-Dichloropropane	0.402	0.436	-8.5#	106	0.00
46	T	Dibromomethane	0.287	0.315	-9.8	106	0.00
47	T	Bromodichloromethane	0.586	0.645	-10.1	108	0.00
48	T	Methyl methacrylate	0.490	0.551	-12.4	106	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	0.006	0.006	0.0	92	0.00
50	S	Toluene-d8	1.289	1.353	-5.0	104	0.00
51	T	4-Methyl-2-Pentanone	0.633	0.729	-15.2	107	0.00
52	CM	Toluene	0.906	1.017	-12.3#	110	0.00
53	T	t-1,3-Dichloropropene	0.601	0.656	-9.2	103	0.00
54	T	cis-1,3-Dichloropropene	0.632	0.686	-8.5	102	0.00
55	T	1,1,2-Trichloroethane	0.351	0.388	-10.5	108	0.00
56	T	Ethyl methacrylate	0.554	0.647	-16.8	104	0.00
57	T	1,3-Dichloropropane	0.633	0.693	-9.5	106	0.00
58	T	2-Chloroethyl Vinyl ether	0.319	0.366	-14.7	99	0.00
59	T	2-Hexanone	0.399	0.492	-23.3	111	0.00
60	T	Dibromochloromethane	0.400	0.448	-12.0	108	0.00
61	T	1,2-Dibromoethane	0.355	0.386	-8.7	103	0.00
62	S	4-Bromofluorobenzene	0.442	0.490	-10.9	108	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	102	0.00
64	T	Tetrachloroethene	0.396	0.368	7.1	99	0.00
65	PM	Chlorobenzene	1.123	1.188	-5.8	109	0.00
66	T	1,1,1,2-Tetrachloroethane	0.379	0.411	-8.4	109	0.00
67	C	Ethyl Benzene	1.963	2.148	-9.4#	113	0.00
68	T	m/p-Xylenes	0.728	0.811	-11.4	112	0.00
69	T	o-Xylene	0.694	0.755	-8.8	110	0.00
70	T	Styrene	1.127	1.297	-15.1	111	0.00
71	P	Bromoform	0.277	0.311	-12.3	111	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	107	0.00
73	T	Isopropylbenzene	3.697	3.944	-6.7	117	0.00
74	T	N-amyl acetate	1.259	1.399	-11.1	144	0.00
75	P	1,1,2,2-Tetrachloroethane	1.195	1.199	-0.3	114	0.00
76	T	1,2,3-Trichloropropane	1.171	1.017	13.2	91	0.00
77	T	Bromobenzene	0.846	0.860	-1.7	107	0.00
78	T	n-propylbenzene	4.523	4.843	-7.1	117	0.00
79	T	2-Chlorotoluene	2.762	2.870	-3.9	114	0.00
80	T	1,3,5-Trimethylbenzene	3.063	3.307	-8.0	116	0.00
81	T	trans-1,4-Dichloro-2-butene	0.413	0.419	-1.5	104	0.00
82	T	4-Chlorotoluene	2.743	2.998	-9.3	113	0.00
83	T	tert-Butylbenzene	2.604	2.804	-7.7	120	0.00
84	T	1,2,4-Trimethylbenzene	3.045	3.337	-9.6	115	0.00
85	T	sec-Butylbenzene	3.741	4.063	-8.6	121	0.00
86	T	p-Isopropyltoluene	3.068	3.328	-8.5	117	0.00
87	T	1,3-Dichlorobenzene	1.647	1.691	-2.7	113	0.00
88	T	1,4-Dichlorobenzene	1.725	1.687	2.2	108	0.00
89	T	n-Butylbenzene	2.979	3.229	-8.4	121	0.00
90	T	Hexachloroethane	0.560	0.613	-9.5	123	0.00
91	T	1,2-Dichlorobenzene	1.550	1.584	-2.2	112	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.290	0.295	-1.7	110	0.00
93	T	1,2,4-Trichlorobenzene	0.921	0.896	2.7	108	0.00
94	T	Hexachlorobutadiene	0.331	0.333	-0.6	120	0.00
95	T	Naphthalene	3.190	3.235	-1.4	105	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.896	0.864	3.6	106	0.00

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