Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091725\
Data File : VX047601.D
Acq On : 17 Sep 2025 08:51
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: Sep 18 01:54:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 17 06:39:58 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	117	-0.01
2	T	Dichlorodifluoromethane	0.518	0.588	-13.5	138	0.00
3	P	Chloromethane	0.680	0.692	-1.8	123	0.00
4	C	Vinyl Chloride	0.666	0.744	-11.7#	132	0.00
5	T	Bromomethane	0.422	0.456	-8.1	126	0.00
6	T	Chloroethane	0.445	0.480	-7.9	127	0.00
7	T	Trichlorofluoromethane	0.989	1.145	-15.8	135	0.00
8	T	Diethyl Ether	0.405	0.426	-5.2	122	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.578	0.682	-18.0	137	0.00
10	T	Methyl Iodide	0.902	0.907	-0.6	117	0.00
11	T	Tert butyl alcohol	0.089	0.088	1.1	115	0.00
12	CM	1,1-Dichloroethene	0.594	0.660	-11.1#	128	0.00
13	T	Acrolein	0.082	0.092	-12.2	133	0.00
14	T	Allyl chloride	1.226	1.281	-4.5	124	0.00
15	T	Acrylonitrile	0.328	0.341	-4.0	115	-0.01
16	T	Acetone	0.346	0.406	-17.3	150	0.00
17	T	Carbon Disulfide	1.626	1.690	-3.9	127	0.00
18	T	Methyl Acetate	0.770	0.866	-12.5	116	0.00
19	T	Methyl tert-butyl Ether	2.169	2.249	-3.7	117	0.00
20	T	Methylene Chloride	0.732	0.734	-0.3	118	0.00
21	T	trans-1,2-Dichloroethene	0.640	0.685	-7.0	124	0.00
22	T	Diisopropyl ether	2.376	2.478	-4.3	117	-0.01
23	T	Vinyl Acetate	1.901	1.942	-2.2	114	0.00
24	P	1,1-Dichloroethane	1.263	1.336	-5.8	121	0.00
25	T	2-Butanone	0.432	0.454	-5.1	121	0.00
26	T	2,2-Dichloropropane	1.050	1.141	-8.7	125	0.00
27	T	cis-1,2-Dichloroethene	0.783	0.811	-3.6	118	-0.01
28	T	Bromochloromethane	0.551	0.577	-4.7	117	-0.01
29	T	Tetrahydrofuran	0.260	0.255	1.9	109	0.00
30	C	Chloroform	1.276	1.326	-3.9#	118	0.00
31	T	Cyclohexane	1.052	1.124	-6.8	128	0.00
32	T	1,1,1-Trichloroethane	1.082	1.164	-7.6	121	0.00
33	S	1,2-Dichloroethane-d4	0.796	0.768	3.5	117	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	113	0.00
35	S	Dibromofluoromethane	0.335	0.352	-5.1	120	-0.01
36	T	1,1-Dichloropropene	0.490	0.515	-5.1	120	-0.01
37	T	Ethyl Acetate	0.525	0.532	-1.3	113	-0.01
38	T	Carbon Tetrachloride	0.532	0.581	-9.2	122	-0.01
39	T	Methylcyclohexane	0.556	0.646	-16.2	129	0.00
40	TM	Benzene	1.488	1.564	-5.1	116	0.00
41	T	Methacrylonitrile	0.271	0.280	-3.3	111	0.00
42	TM	1,2-Dichloroethane	0.566	0.577	-1.9	112	0.00
43	T	Isopropyl Acetate	0.860	0.869	-1.0	110	0.00
44	TM	Trichloroethene	0.360	0.386	-7.2	118	0.00
45	C	1,2-Dichloropropane	0.381	0.397	-4.2#	113	0.00
46	T	Dibromomethane	0.277	0.281	-1.4	111	0.00
47	T	Bromodichloromethane	0.572	0.605	-5.8	113	0.00
48	T	Methyl methacrylate	0.429	0.430	-0.2	108	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	0.004	0.005	-25.0	110	0.00
50	S	Toluene-d8	1.160	1.191	-2.7	117	0.00
51	T	4-Methyl-2-Pentanone	0.477	0.481	-0.8	107	0.00
52	CM	Toluene	0.917	0.945	-3.1#	114	0.00
53	T	t-1,3-Dichloropropene	0.584	0.604	-3.4	110	0.00
54	T	cis-1,3-Dichloropropene	0.624	0.650	-4.2	113	0.00
55	T	1,1,2-Trichloroethane	0.354	0.354	0.0	109	0.00
56	T	Ethyl methacrylate	0.555	0.581	-4.7	110	0.00
57	T	1,3-Dichloropropane	0.607	0.620	-2.1	110	0.00
58	T	2-Chloroethyl Vinyl ether	0.227	0.279	-22.9	116	0.00
59	T	2-Hexanone	0.343	0.342	0.3	108	0.00
60	T	Dibromochloromethane	0.407	0.425	-4.4	111	0.00
61	T	1,2-Dibromoethane	0.360	0.366	-1.7	109	0.00
62	S	4-Bromofluorobenzene	0.440	0.448	-1.8	119	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	110	0.00
64	T	Tetrachloroethene	0.326	0.353	-8.3	119	0.00
65	PM	Chlorobenzene	1.136	1.193	-5.0	111	0.00
66	T	1,1,1,2-Tetrachloroethane	0.392	0.420	-7.1	111	0.00
67	C	Ethyl Benzene	1.960	2.137	-9.0#	115	0.00
68	T	m/p-Xylenes	0.729	0.793	-8.8	114	0.00
69	T	o-Xylene	0.706	0.757	-7.2	111	0.00
70	T	Styrene	1.236	1.311	-6.1	111	0.00
71	P	Bromoform	0.293	0.312	-6.5	109	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	108	0.00
73	T	Isopropylbenzene	3.801	4.316	-13.5	115	0.00
74	T	N-amyl acetate	1.733	1.844	-6.4	107	0.00
75	P	1,1,2,2-Tetrachloroethane	1.150	1.208	-5.0	107	0.00
76	T	1,2,3-Trichloropropane	1.012	0.972	4.0	93	0.00
77	T	Bromobenzene	0.936	1.005	-7.4	110	0.00
78	T	n-propylbenzene	4.494	5.065	-12.7	114	0.00
79	T	2-Chlorotoluene	2.752	2.995	-8.8	111	0.00
80	T	1,3,5-Trimethylbenzene	3.123	3.476	-11.3	113	0.00
81	T	trans-1,4-Dichloro-2-butene	0.397	0.420	-5.8	110	0.00
82	T	4-Chlorotoluene	3.251	3.550	-9.2	112	0.00
83	T	tert-Butylbenzene	3.197	3.598	-12.5	116	0.00
84	T	1,2,4-Trimethylbenzene	3.166	3.460	-9.3	112	0.00
85	T	sec-Butylbenzene	3.797	4.370	-15.1	118	0.00
86	T	p-Isopropyltoluene	3.215	3.633	-13.0	115	0.00
87	T	1,3-Dichlorobenzene	1.739	1.839	-5.8	110	0.00
88	T	1,4-Dichlorobenzene	1.785	1.872	-4.9	111	0.00
89	T	n-Butylbenzene	2.975	3.457	-16.2	119	0.00
90	T	Hexachloroethane	0.564	0.646	-14.5	118	0.00
91	T	1,2-Dichlorobenzene	1.657	1.766	-6.6	110	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.233	0.247	-6.0	105	0.00
93	T	1,2,4-Trichlorobenzene	1.077	1.192	-10.7	112	0.00
94	T	Hexachlorobutadiene	0.366	0.428	-16.9	121	0.00
95	T	Naphthalene	3.279	3.403	-3.8	104	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	1.002	1.099	-9.7	109	0.00

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