Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR080625\
Data File : VR026721.D
Acq On : 6 Aug 2025 17:03
Operator : SY/MD
Sample : VSTDCCC050
Misc : 5mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_R
LabSampleId :
VSTDCCC050

Quant Time: Aug 07 01:16:52 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_R\methods\82R080425W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 05 02:27:49 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	50#	-0.02
2	T	Dichlorodifluoromethane	0.364	0.321	11.8	45#	0.00
3	P	Chloromethane	0.542	0.493	9.0	48#	-0.01
4	C	Vinyl Chloride	0.554	0.601	-8.5#	58	-0.01
5	T	Bromomethane	0.270	0.303	-12.2	59	-0.01
6	T	Chloroethane	0.356	0.401	-12.6	62	-0.01
7	T	Trichlorofluoromethane	0.998	1.304	-30.7#	67	0.00
8	T	Diethyl Ether	0.443	0.417	5.9	48#	-0.02
9	T	1,1,2-Trichlorotrifluoroeth	0.538	0.484	10.0	46#	-0.02
10	T	Methyl Iodide	0.519	0.486	6.4	46#	-0.02
11	T	Tert butyl alcohol	0.099	0.085	14.1	43#	-0.02
12	CM	1,1-Dichloroethene	0.532	0.486	8.6#	46#	-0.02
13	T	Acrolein	0.145	0.121	16.6	41#	0.00
14	T	Allyl chloride	1.046	1.023	2.2	53	-0.02
15	T	Acrylonitrile	0.393	0.419	-6.6	51	-0.02
16	T	Acetone	0.297	0.289	2.7	46#	-0.02
17	T	Carbon Disulfide	0.787	0.595	24.4	38#	-0.02
18	T	Methyl Acetate	1.240	1.414	-14.0	56	-0.02
19	T	Methyl tert-butyl Ether	2.129	1.979	7.0	47#	-0.02
20	T	Methylene Chloride	0.705	0.642	8.9	47#	-0.02
21	T	trans-1,2-Dichloroethene	0.585	0.521	10.9	45#	-0.02
22	T	Diisopropyl ether	2.452	2.435	0.7	50#	-0.02
23	T	Vinyl Acetate	1.398	1.476	-5.6	52	-0.02
24	P	1,1-Dichloroethane	1.242	1.141	8.1	47#	-0.03
25	T	2-Butanone	0.460	0.502	-9.1	52	-0.02
26	T	2,2-Dichloropropane	0.858	0.696	18.9	40#	-0.02
27	T	cis-1,2-Dichloroethene	0.752	0.716	4.8	47#	-0.02
28	T	Bromochloromethane	0.447	0.412	7.8	51	-0.02
29	C	Chloroform	1.076	1.081	-0.5#	50	-0.02
30	T	Cyclohexane	1.003	0.887	11.6	48#	-0.03
31	T	1,1,1-Trichloroethane	0.777	0.768	1.2	49#	-0.02
32	S	1,2-Dichloroethane-d4	0.633	0.625	1.3	49#	-0.02
33	I	1,4-Difluorobenzene	1.000	1.000	0.0	50#	-0.01
34	S	Dibromofluoromethane	0.309	0.310	-0.3	48#	-0.01
35	T	1,1-Dichloropropene	0.390	0.395	-1.3	50#	-0.02
36	T	Ethyl Acetate	0.467	0.469	-0.4	49#	-0.01
37	T	Carbon Tetrachloride	0.318	0.326	-2.5	50#	-0.03
38	T	Methylcyclohexane	0.447	0.448	-0.2	50#	-0.02
39	TM	Benzene	1.405	1.414	-0.6	50#	-0.02
40	T	Methacrylonitrile	0.290	0.303	-4.5	53	-0.02
41	TM	1,2-Dichloroethane	0.357	0.372	-4.2	53	-0.02
42	T	Isopropyl Acetate	0.751	0.805	-7.2	52	-0.01
43	TM	Trichloroethene	0.270	0.283	-4.8	54	-0.02
44	C	1,2-Dichloropropane	0.332	0.348	-4.8#	53	-0.02
45	T	Dibromomethane	0.185	0.199	-7.6	55	-0.02
46	T	Bromodichloromethane	0.358	0.400	-11.7	54	-0.02
47	T	Methyl methacrylate	0.329	0.390	-18.5	56	-0.01
48	T	1,4-Dioxane	0.005	0.006	-20.0	58	-0.02

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Quant Title : SW846 8260
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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	S	Toluene-d8	1.108	1.209	-9.1	52	-0.02
50	T	4-Methyl-2-Pentanone	0.392	0.504	-28.6#	61	-0.02
51	CM	Toluene	0.717	0.821	-14.5#	57	0.00
52	T	t-1,3-Dichloropropene	0.351	0.369	-5.1	50	-0.02
53	T	cis-1,3-Dichloropropene	0.421	0.457	-8.6	52	-0.01
54	T	1,1,2-Trichloroethane	0.262	0.321	-22.5	60	0.00
55	T	Ethyl methacrylate	0.433	0.521	-20.3	56	-0.01
56	T	1,3-Dichloropropane	0.446	0.515	-15.5	56	-0.02
57	T	2-Chloroethyl Vinyl ether	0.024	0.031	-29.2#	57	-0.02
58	T	2-Hexanone	0.264	0.342	-29.5#	59	-0.01
59	T	Dibromochloromethane	0.222	0.264	-18.9	57	-0.01
60	T	1,2-Dibromoethane	0.225	0.273	-21.3	60	-0.02
61	S	4-Bromofluorobenzene	0.315	0.344	-9.2	53	-0.01
62	I	Chlorobenzene-d5	1.000	1.000	0.0	56	-0.02
63	T	Tetrachloroethene	0.273	0.275	-0.7	58	-0.02
64	PM	Chlorobenzene	1.041	1.092	-4.9	60	-0.02
65	T	1,1,1,2-Tetrachloroethane	0.284	0.303	-6.7	57	-0.02
66	C	Ethyl Benzene	1.737	1.852	-6.6#	59	-0.02
67	T	m/p-Xylenes	0.694	0.730	-5.2	59	-0.02
68	T	o-Xylene	0.674	0.718	-6.5	59	-0.02
69	T	Styrene	1.126	1.224	-8.7	60	-0.02
70	P	Bromoform	0.161	0.188	-16.8	60	-0.02
71	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	59	-0.02
72	T	Isopropylbenzene	3.564	3.499	1.8	59	-0.01
73	T	N-amyl acetate	1.413	1.433	-1.4	59	-0.02
74	P	1,1,2,2-Tetrachloroethane	1.046	1.074	-2.7	61	-0.01
75	T	1,2,3-Trichloropropane	0.829	0.846	-2.1	62	-0.01
76	T	Bromobenzene	0.825	0.821	0.5	61	0.00
77	T	n-propylbenzene	4.252	4.206	1.1	59	-0.01
78	T	2-Chlorotoluene	2.412	2.267	6.0	57	-0.02
79	T	1,3,5-Trimethylbenzene	2.787	2.737	1.8	58	-0.02
80	T	trans-1,4-Dichloro-2-butene	0.285	0.257	9.8	49#	-0.02
81	T	4-Chlorotoluene	2.464	2.361	4.2	59	-0.02
82	T	tert-Butylbenzene	2.535	2.503	1.3	58	-0.01
83	T	1,2,4-Trimethylbenzene	2.759	2.849	-3.3	59	-0.02
84	T	sec-Butylbenzene	3.759	3.854	-2.5	60	-0.02
85	T	p-Isopropyltoluene	3.065	3.227	-5.3	60	-0.01
86	T	1,3-Dichlorobenzene	1.617	1.649	-2.0	60	-0.02
87	T	1,4-Dichlorobenzene	1.623	1.668	-2.8	62	-0.02
88	T	n-Butylbenzene	2.825	2.844	-0.7	57	-0.01
89	T	Hexachloroethane	0.440	0.434	1.4	56	-0.02
90	T	1,2-Dichlorobenzene	1.519	1.582	-4.1	63	-0.02
91	T	1,2-Dibromo-3-Chloropropane	0.139	0.146	-5.0	64	-0.01
92	T	1,2,4-Trichlorobenzene	0.908	0.962	-5.9	65	-0.02
93	T	Hexachlorobutadiene	0.295	0.289	2.0	63	-0.02
94	T	Naphthalene	2.783	3.003	-7.9	64	-0.02
95	T	1,2,3-Trichlorobenzene	0.862	0.939	-8.9	64	-0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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