Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090123\
Data File : VN079117.D
Acq On : 01 Sep 2023 12:35
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_N
LabSampleId :
VSTDCCC050

Quant Time: Sep 01 23:09:20 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 16 02:42:02 2023
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	142	0.00
2	T	Dichlorodifluoromethane	0.694	0.654	5.8	139	0.00
3	P	Chloromethane	0.969	0.773	20.2	118	0.00
4	C	Vinyl Chloride	0.849	0.923	-8.7#	158#	0.00
5	T	Bromomethane	0.488	0.340	30.3#	106	0.00
6	T	Chloroethane	0.594	0.355	40.2#	90	0.00
7	T	Trichlorofluoromethane	1.035	1.056	-2.0	147	0.00
8	T	Diethyl Ether	0.392	0.401	-2.3	151#	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.551	0.581	-5.4	152#	0.00
10	T	Methyl Iodide	0.617	0.756	-22.5	164#	-0.01
11	T	Tert butyl alcohol	0.131	0.104	20.6	116	-0.02
12	CM	1,1-Dichloroethene	0.537	0.567	-5.6#	151#	0.00
13	T	Acrolein	0.067	0.065	3.0	131	-0.01
14	T	Allyl chloride	1.068	0.897	16.0	115	0.00
15	T	Acrylonitrile	0.354	0.333	5.9	133	0.00
16	T	Acetone	0.321	0.320	0.3	139	-0.01
17	T	Carbon Disulfide	1.503	1.504	-0.1	146	-0.01
18	T	Methyl Acetate	1.356	1.110	18.1	119	-0.01
19	T	Methyl tert-butyl Ether	1.934	2.012	-4.0	146	-0.02
20	T	Methylene Chloride	0.684	0.704	-2.9	152#	0.00
21	T	trans-1,2-Dichloroethene	0.602	0.636	-5.6	150	0.00
22	T	Diisopropyl ether	2.260	2.197	2.8	135	-0.01
23	T	Vinyl Acetate	1.488	1.361	8.5	125	0.00
24	P	1,1-Dichloroethane	1.235	1.263	-2.3	147	0.00
25	T	2-Butanone	0.515	0.449	12.8	123	-0.01
26	T	2,2-Dichloropropane	0.998	1.057	-5.9	149	-0.01
27	T	cis-1,2-Dichloroethene	0.732	0.771	-5.3	153#	-0.01
28	T	Bromochloromethane	0.655	0.588	10.2	126	0.00
29	T	Tetrahydrofuran	0.341	0.282	17.3	118	0.00
30	C	Chloroform	1.234	1.289	-4.5#	148	0.00
31	T	Cyclohexane	1.068	0.986	7.7	141	0.00
32	T	1,1,1-Trichloroethane	1.052	1.095	-4.1	148	0.00
33	S	1,2-Dichloroethane-d4	0.778	0.726	6.7	134	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	134	0.00
35	S	Dibromofluoromethane	0.337	0.375	-11.3	147	0.00
36	T	1,1-Dichloropropene	0.506	0.573	-13.2	147	0.00
37	T	Ethyl Acetate	0.566	0.539	4.8	122	0.00
38	T	Carbon Tetrachloride	0.507	0.566	-11.6	146	0.00
39	T	Methylcyclohexane	0.466	0.545	-17.0	151#	0.00
40	TM	Benzene	1.496	1.710	-14.3	150#	0.00
41	T	Methacrylonitrile	0.309	0.285	7.8	118	-0.01
42	TM	1,2-Dichloroethane	0.559	0.576	-3.0	137	0.00
43	T	Isopropyl Acetate	1.001	0.867	13.4	115	0.00
44	TM	Trichloroethene	0.356	0.415	-16.6	155#	0.00
45	C	1,2-Dichloropropane	0.398	0.444	-11.6#	145	0.00
46	T	Dibromomethane	0.273	0.290	-6.2	144	0.00
47	T	Bromodichloromethane	0.540	0.605	-12.0	148	0.00
48	T	Methyl methacrylate	0.454	0.417	8.1	122	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	T	1,4-Dioxane	0.008	0.007	12.5	123	0.00
50	S	Toluene-d8	1.250	1.317	-5.4	135	0.00
51	T	4-Methyl-2-Pentanone	0.594	0.533	10.3	117	0.00
52	CM	Toluene	0.919	1.046	-13.8#	148	0.00
53	T	t-1,3-Dichloropropene	0.579	0.652	-12.6	143	0.00
54	T	cis-1,3-Dichloropropene	0.630	0.713	-13.2	148	0.00
55	T	1,1,2-Trichloroethane	0.368	0.421	-14.4	151#	0.00
56	T	Ethyl methacrylate	0.571	0.648	-13.5	143	0.00
57	T	1,3-Dichloropropane	0.659	0.723	-9.7	146	0.00
58	T	2-Chloroethyl Vinyl ether	0.274	0.213	22.3	101	0.00
59	T	2-Hexanone	0.430	0.393	8.6	117	0.00
60	T	Dibromochloromethane	0.388	0.445	-14.7	150#	0.00
61	T	1,2-Dibromoethane	0.374	0.412	-10.2	146	0.00
62	S	4-Bromofluorobenzene	0.408	0.485	-18.9	149	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	132	0.00
64	T	Tetrachloroethene	0.314	0.381	-21.3	165#	0.00
65	PM	Chlorobenzene	1.064	1.204	-13.2	147	0.00
66	T	1,1,1,2-Tetrachloroethane	0.402	0.465	-15.7	148	0.00
67	C	Ethyl Benzene	1.935	2.231	-15.3#	150#	0.00
68	T	m/p-Xylenes	0.724	0.825	-14.0	145	0.00
69	T	o-Xylene	0.704	0.802	-13.9	143	0.00
70	T	Styrene	1.174	1.363	-16.1	144	0.00
71	P	Bromoform	0.292	0.348	-19.2	152#	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	129	0.00
73	T	Isopropylbenzene	4.169	4.552	-9.2	145	0.00
74	T	N-amyl acetate	2.104	1.789	15.0	111	0.00
75	P	1,1,2,2-Tetrachloroethane	1.505	1.491	0.9	138	0.00
76	T	1,2,3-Trichloropropane	1.205	1.169	3.0	134	0.00
77	T	Bromobenzene	1.002	1.098	-9.6	146	0.00
78	T	n-propylbenzene	4.834	5.349	-10.7	143	0.00
79	T	2-Chlorotoluene	2.999	3.283	-9.5	148	0.00
80	T	1,3,5-Trimethylbenzene	3.304	3.764	-13.9	150	0.00
81	T	trans-1,4-Dichloro-2-butene	0.472	0.474	-0.4	129	0.00
82	T	4-Chlorotoluene	2.934	3.242	-10.5	143	0.00
83	T	tert-Butylbenzene	2.776	3.005	-8.2	140	0.00
84	T	1,2,4-Trimethylbenzene	3.271	3.701	-13.1	146	0.00
85	T	sec-Butylbenzene	3.820	4.294	-12.4	148	0.00
86	T	p-Isopropyltoluene	3.052	3.521	-15.4	148	0.00
87	T	1,3-Dichlorobenzene	1.724	1.889	-9.6	141	0.00
88	T	1,4-Dichlorobenzene	1.726	1.834	-6.3	139	0.00
89	T	n-Butylbenzene	2.512	2.976	-18.5	148	0.00
90	T	Hexachloroethane	0.612	0.662	-8.2	151#	0.00
91	T	1,2-Dichlorobenzene	1.715	1.859	-8.4	144	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.284	0.241	15.1	119	0.00
93	T	1,2,4-Trichlorobenzene	0.608	0.745	-22.5	151#	0.00
94	T	Hexachlorobutadiene	0.329	0.425	-29.2#	180#	0.00
95	T	Naphthalene	2.054	2.096	-2.0	120	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.666	0.775	-16.4	147	0.00

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