Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049475.D
Acq On : 28 Jun 2022 21:51
Operator : SY/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC050

Quant Time: Jun 29 05:23:20 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 28 10:35:43 2022
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	Pentafluorobenzene	1.000	1.000	0.0	157#	0.00
2	T	Dichlorodifluoromethane	0.803	0.735	8.5	135	0.00
3	P	Chloromethane	1.088	0.814	25.2#	131	0.00
4	C	Vinyl Chloride	0.809	0.778	3.8#	145	0.00
5	T	Bromomethane	0.252	0.288	-14.3	195#	0.00
6	T	Chloroethane	0.377	0.404	-7.2	191#	0.01
7	T	Trichlorofluoromethane	1.031	0.949	8.0	142	0.02
8	T	Diethyl Ether	0.372	0.381	-2.4	167#	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.555	0.576	-3.8	158#	0.02
10	T	Methyl Iodide	0.489	0.771	-57.7#	284#	0.01
11	T	Tert butyl alcohol	0.180	0.214	-18.9	204#	-0.11
12	CM	1,1-Dichloroethene	0.543	0.573	-5.5#	166#	0.02
13	T	Acrolein	0.122	0.081	33.6#	109	0.00
14	T	Allyl chloride	1.098	1.099	-0.1	158#	0.00
15	T	Acrylonitrile	0.433	0.451	-4.2	166#	0.00
16	T	Acetone	0.388	0.366	5.7	163#	0.00
17	T	Carbon Disulfide	1.694	1.614	4.7	156#	0.01
18	T	Methyl Acetate	1.051	1.022	2.8	161#	0.00
19	T	Methyl tert-butyl Ether	2.020	2.134	-5.6	167#	0.00
20	T	Methylene Chloride	0.715	0.713	0.3	173#	0.00
21	T	trans-1,2-Dichloroethene	0.585	0.603	-3.1	162#	0.00
22	T	Diisopropyl ether	2.096	2.177	-3.9	162#	0.00
23	T	Vinyl Acetate	1.837	1.915	-4.2	161#	0.00
24	P	1,1-Dichloroethane	1.204	1.217	-1.1	159#	0.00
25	T	2-Butanone	0.582	0.605	-4.0	172#	0.00
26	T	2,2-Dichloropropane	1.002	0.816	18.6	127	0.00
27	T	cis-1,2-Dichloroethene	0.676	0.722	-6.8	169#	0.00
28	T	Bromochloromethane	0.526	0.512	2.7	149	0.00
29	T	Tetrahydrofuran	0.399	0.401	-0.5	165#	0.00
30	C	Chloroform	1.191	1.183	0.7#	158#	0.00
31	T	Cyclohexane	1.156	1.099	4.9	152#	0.00
32	T	1,1,1-Trichloroethane	1.033	1.032	0.1	152#	0.00
33	S	1,2-Dichloroethane-d4	0.770	0.717	6.9	144	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	166#	0.00
35	S	Dibromofluoromethane	0.350	0.327	6.6	153#	0.00
36	T	1,1-Dichloropropene	0.507	0.489	3.6	158#	0.00
37	T	Ethyl Acetate	0.693	0.650	6.2	143	0.00
38	T	Carbon Tetrachloride	0.532	0.500	6.0	151#	0.00
39	T	Methylcyclohexane	0.608	0.584	3.9	150	0.00
40	TM	Benzene	1.531	1.505	1.7	162#	0.00
41	T	Methacrylonitrile	0.347	0.342	1.4	166#	0.00
42	TM	1,2-Dichloroethane	0.570	0.518	9.1	148	0.00
43	T	Isopropyl Acetate	1.016	0.976	3.9	160#	0.00
44	TM	Trichloroethene	0.364	0.359	1.4	163#	0.00
45	C	1,2-Dichloropropane	0.424	0.407	4.0#	161#	0.00
46	T	Dibromomethane	0.283	0.277	2.1	160#	0.00
47	T	Bromodichloromethane	0.545	0.519	4.8	156#	0.00
48	T	Methyl methacrylate	0.476	0.469	1.5	162#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	T	1,4-Dioxane	0.012	0.013	-8.3	185#	-0.01
50	S	Toluene-d8	1.248	1.188	4.8	156#	0.00
51	T	4-Methyl-2-Pentanone	0.680	0.641	5.7	155#	-0.01
52	CM	Toluene	0.910	0.906	0.4#	164#	0.00
53	T	t-1,3-Dichloropropene	0.584	0.556	4.8	152#	0.00
54	T	cis-1,3-Dichloropropene	0.624	0.615	1.4	157#	0.00
55	T	1,1,2-Trichloroethane	0.376	0.375	0.3	160#	0.00
56	T	Ethyl methacrylate	0.629	0.658	-4.6	164#	0.00
57	T	1,3-Dichloropropane	0.668	0.649	2.8	159#	0.00
58	T	2-Chloroethyl Vinyl ether	0.043	0.075	-74.4#	228#	0.00
59	T	2-Hexanone	0.534	0.501	6.2	157#	-0.01
60	T	Dibromochloromethane	0.391	0.391	0.0	156#	0.00
61	T	1,2-Dibromoethane	0.410	0.405	1.2	164#	0.00
62	S	4-Bromofluorobenzene	0.478	0.441	7.7	149	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	157#	0.00
64	T	Tetrachloroethene	0.321	0.318	0.9	157#	0.00
65	PM	Chlorobenzene	1.022	1.021	0.1	159#	0.00
66	T	1,1,1,2-Tetrachloroethane	0.386	0.400	-3.6	160#	0.00
67	C	Ethyl Benzene	1.877	1.939	-3.3#	161#	0.00
68	T	m/p-Xylenes	0.694	0.718	-3.5	161#	0.00
69	T	o-Xylene	0.689	0.720	-4.5	161#	0.00
70	T	Styrene	1.108	1.185	-6.9	160#	0.00
71	P	Bromoform	0.331	0.341	-3.0	155#	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	147	0.00
73	T	Isopropylbenzene	3.562	4.016	-12.7	157#	0.00
74	T	N-amyl acetate	1.885	2.032	-7.8	156#	0.00
75	P	1,1,2,2-Tetrachloroethane	1.509	1.580	-4.7	154#	0.00
76	T	1,2,3-Trichloropropane	1.695	1.752	-3.4	151#	0.00
77	T	Bromobenzene	0.869	0.938	-7.9	158#	0.00
78	T	n-propylbenzene	4.328	4.613	-6.6	149	0.00
79	T	2-Chlorotoluene	2.666	2.811	-5.4	151#	0.00
80	T	1,3,5-Trimethylbenzene	3.110	3.362	-8.1	152#	0.00
81	T	trans-1,4-Dichloro-2-butene	0.459	0.465	-1.3	145	0.00
82	T	4-Chlorotoluene	2.999	3.180	-6.0	151#	0.00
83	T	tert-Butylbenzene	3.058	3.342	-9.3	152#	0.00
84	T	1,2,4-Trimethylbenzene	3.055	3.375	-10.5	152#	0.00
85	T	sec-Butylbenzene	3.845	4.077	-6.0	146	0.00
86	T	p-Isopropyltoluene	3.110	3.328	-7.0	148	0.00
87	T	1,3-Dichlorobenzene	1.662	1.682	-1.2	150#	0.00
88	T	1,4-Dichlorobenzene	1.671	1.699	-1.7	155#	0.00
89	T	n-Butylbenzene	2.879	2.912	-1.1	149	0.00
90	T	Hexachloroethane	0.619	0.598	3.4	137	0.00
91	T	1,2-Dichlorobenzene	1.670	1.742	-4.3	154#	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.387	0.403	-4.1	155#	0.00
93	T	1,2,4-Trichlorobenzene	1.159	1.113	4.0	155#	0.00
94	T	Hexachlorobutadiene	0.518	0.478	7.7	132	0.00
95	T	Naphthalene	4.457	4.300	3.5	173#	0.00

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Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	1.217	1.166	4.2	154#	0.00

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