Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060622\
Data File : VU048876.D
Acq On : 06 Jun 2022 12:14
Operator : SY/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC050

Quant Time: Jun 07 05:13:29 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260
QLast Update : Fri May 27 06:42:31 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	88	0.00
2	T	Dichlorodifluoromethane	0.572	0.502	12.2	84	0.00
3	P	Chloromethane	0.737	0.667	9.5	88	0.00
4	C	Vinyl Chloride	0.682	0.647	5.1#	87	0.00
5	T	Bromomethane	0.426	0.335	21.4#	76	-0.02
6	T	Chloroethane	0.425	0.367	13.6	83	-0.02
7	T	Trichlorofluoromethane	0.975	0.897	8.0	85	0.00
8	T	Diethyl Ether	0.360	0.357	0.8	91	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.534	0.504	5.6	90	0.00
10	T	Methyl Iodide	0.704	0.722	-2.6	79	0.00
11	T	Tert butyl alcohol	0.190	0.216	-13.7	114	0.00
12	CM	1,1-Dichloroethene	0.509	0.480	5.7#	88	0.00
13	T	Acrolein	0.073	0.055	24.7#	79	0.00
14	T	Allyl chloride	0.986	1.004	-1.8	97	0.00
15	T	Acrylonitrile	0.413	0.450	-9.0	99	0.00
16	T	Acetone	0.383	0.435	-13.6	112	0.00
17	T	Carbon Disulfide	1.486	1.174	21.0#	78	0.00
18	T	Methyl Acetate	0.961	1.115	-16.0	111	0.00
19	T	Methyl tert-butyl Ether	1.931	2.043	-5.8	96	0.00
20	T	Methylene Chloride	0.651	0.618	5.1	92	0.00
21	T	trans-1,2-Dichloroethene	0.582	0.522	10.3	85	0.00
22	T	Diisopropyl ether	1.818	2.040	-12.2	99	0.00
23	T	Vinyl Acetate	1.575	1.772	-12.5	96	0.00
24	P	1,1-Dichloroethane	1.142	1.165	-2.0	94	0.00
25	T	2-Butanone	0.550	0.636	-15.6	105	0.00
26	T	2,2-Dichloropropane	1.024	1.009	1.5	93	0.00
27	T	cis-1,2-Dichloroethene	0.674	0.664	1.5	93	0.00
28	T	Bromochloromethane	0.448	0.487	-8.7	94	0.00
29	T	Tetrahydrofuran	0.346	0.388	-12.1	100	0.00
30	C	Chloroform	1.143	1.180	-3.2#	94	0.00
31	T	Cyclohexane	0.946	0.900	4.9	89	0.00
32	T	1,1,1-Trichloroethane	1.037	1.011	2.5	90	0.00
33	S	1,2-Dichloroethane-d4	0.751	0.743	1.1	86	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	90	0.00
35	S	Dibromofluoromethane	0.357	0.349	2.2	88	0.00
36	T	1,1-Dichloropropene	0.521	0.489	6.1	90	0.00
37	T	Ethyl Acetate	0.658	0.693	-5.3	99	0.00
38	T	Carbon Tetrachloride	0.597	0.544	8.9	86	0.00
39	T	Methylcyclohexane	0.546	0.553	-1.3	95	0.00
40	TM	Benzene	1.537	1.478	3.8	93	0.00
41	T	Methacrylonitrile	0.333	0.381	-14.4	102	0.00
42	TM	1,2-Dichloroethane	0.602	0.587	2.5	92	0.00
43	T	Isopropyl Acetate	0.955	1.023	-7.1	100	0.00
44	TM	Trichloroethene	0.396	0.360	9.1	87	0.00
45	C	1,2-Dichloropropane	0.387	0.425	-9.8#	97	0.00
46	T	Dibromomethane	0.289	0.292	-1.0	93	0.00
47	T	Bromodichloromethane	0.585	0.580	0.9	95	0.00
48	T	Methyl methacrylate	0.429	0.489	-14.0	102	0.00

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

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	T	1,4-Dioxane	0.010	0.013	-30.0#	131	0.00
50	S	Toluene-d8	1.279	1.224	4.3	86	0.00
51	T	4-Methyl-2-Pentanone	0.620	0.764	-23.2#	110	0.00
52	CM	Toluene	0.919	0.923	-0.4#	91	0.00
53	T	t-1,3-Dichloropropene	0.631	0.646	-2.4	95	0.00
54	T	cis-1,3-Dichloropropene	0.653	0.682	-4.4	95	0.00
55	T	1,1,2-Trichloroethane	0.406	0.410	-1.0	97	0.00
56	T	Ethyl methacrylate	0.566	0.685	-21.0#	106	0.00
57	T	1,3-Dichloropropane	0.683	0.696	-1.9	96	0.00
58	T	2-Chloroethyl Vinyl ether	0.111	0.122	-9.9	98	0.00
59	T	2-Hexanone	0.509	0.622	-22.2#	114	0.00
60	T	Dibromochloromethane	0.455	0.463	-1.8	95	0.00
61	T	1,2-Dibromoethane	0.447	0.428	4.3	92	0.00
62	S	4-Bromofluorobenzene	0.484	0.503	-3.9	92	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	87	0.00
64	T	Tetrachloroethene	0.384	0.336	12.5	84	0.00
65	PM	Chlorobenzene	1.058	1.077	-1.8	92	0.00
66	T	1,1,1,2-Tetrachloroethane	0.401	0.431	-7.5	95	0.00
67	C	Ethyl Benzene	1.784	1.919	-7.6#	92	0.00
68	T	m/p-Xylenes	0.698	0.760	-8.9	92	0.00
69	T	o-Xylene	0.685	0.761	-11.1	96	0.00
70	T	Styrene	1.127	1.312	-16.4	96	0.00
71	P	Bromoform	0.371	0.428	-15.4	100	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	93	0.00
73	T	Isopropylbenzene	3.228	3.465	-7.3	97	0.00
74	T	N-amyl acetate	1.396	1.813	-29.9#	113	0.00
75	P	1,1,2,2-Tetrachloroethane	1.317	1.533	-16.4	114	0.00
76	T	1,2,3-Trichloropropane	1.612	1.747	-8.4	107	0.00
77	T	Bromobenzene	0.867	0.876	-1.0	98	0.00
78	T	n-propylbenzene	3.846	4.127	-7.3	96	0.00
79	T	2-Chlorotoluene	2.392	2.556	-6.9	100	0.00
80	T	1,3,5-Trimethylbenzene	2.807	3.097	-10.3	99	0.00
81	T	trans-1,4-Dichloro-2-butene	0.444	0.475	-7.0	101	0.00
82	T	4-Chlorotoluene	2.760	2.942	-6.6	97	0.00
83	T	tert-Butylbenzene	2.694	3.132	-16.3	106	0.00
84	T	1,2,4-Trimethylbenzene	2.732	3.174	-16.2	101	0.00
85	T	sec-Butylbenzene	3.371	3.924	-16.4	104	0.00
86	T	p-Isopropyltoluene	2.894	3.373	-16.6	104	0.00
87	T	1,3-Dichlorobenzene	1.700	1.731	-1.8	99	0.00
88	T	1,4-Dichlorobenzene	1.756	1.737	1.1	98	0.00
89	T	n-Butylbenzene	2.544	2.927	-15.1	101	0.00
90	T	Hexachloroethane	0.591	0.628	-6.3	108	0.00
91	T	1,2-Dichlorobenzene	1.717	1.806	-5.2	104	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.379	0.387	-2.1	98	0.00
93	T	1,2,4-Trichlorobenzene	1.127	1.063	5.7	89	0.00
94	T	Hexachlorobutadiene	0.540	0.568	-5.2	110	0.00
95	T	Naphthalene	3.495	3.619	-3.5	93	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	1.116	1.095	1.9	91	0.00

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