Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050922\
Data File : VX028562.D
Acq On : 09 May 2022 09:50
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: May 09 15:48:18 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
Quant Title : SW846 8260
QLast Update : Tue Apr 19 13:38:04 2022
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	105	0.00
2	T	Dichlorodifluoromethane	0.323	0.380	-17.6	138	0.00
3	P	Chloromethane	0.375	0.364	2.9	117	0.00
4	C	Vinyl Chloride	0.504	0.453	10.1#	105	0.00
5	T	Bromomethane	0.522	0.466	10.7	100	0.00
6	T	Chloroethane	0.461	0.336	27.1#	99	0.00
7	T	Trichlorofluoromethane	1.112	0.887	20.2	93	0.00
8	T	Diethyl Ether	0.382	0.340	11.0	106	0.00
9	T	1,1,2-Trichlorotrifluoroeth	0.405	0.493	-21.7	158#	0.00
10	T	Methyl Iodide	0.509	0.591	-16.1	125	0.00
11	T	Tert butyl alcohol	0.135	0.112	17.0	101	0.00
12	CM	1,1-Dichloroethene	0.384	0.432	-12.5#	145	0.00
13	T	Acrolein	0.080	0.235	-193.7#	284#	0.00
14	T	Allyl chloride	0.645	0.642	0.5	113	0.00
15	T	Acrylonitrile	0.266	0.277	-4.1	113	0.00
16	T	Acetone	0.215	0.230	-7.0	131	0.00
17	T	Carbon Disulfide	1.035	0.877	15.3	99	0.00
18	T	Methyl Acetate	0.580	0.593	-2.2	120	0.00
19	T	Methyl tert-butyl Ether	1.602	1.706	-6.5	112	0.00
20	T	Methylene Chloride	0.484	0.527	-8.9	131	0.00
21	T	trans-1,2-Dichloroethene	0.523	0.484	7.5	102	0.00
22	T	Diisopropyl ether	1.469	1.493	-1.6	111	0.00
23	T	Vinyl Acetate	1.283	1.288	-0.4	106	0.00
24	P	1,1-Dichloroethane	0.867	0.901	-3.9	120	0.00
25	T	2-Butanone	0.386	0.367	4.9	104	0.00
26	T	2,2-Dichloropropane	0.847	0.792	6.5	103	0.00
27	T	cis-1,2-Dichloroethene	0.623	0.624	-0.2	111	0.00
28	T	Bromochloromethane	0.372	0.336	9.7	96	0.00
29	T	Tetrahydrofuran	0.252	0.229	9.1	100	-0.01
30	C	Chloroform	1.041	1.032	0.9#	108	0.00
31	T	Cyclohexane	0.794	0.671	15.5	92	0.00
32	T	1,1,1-Trichloroethane	0.943	0.901	4.5	102	0.00
33	S	1,2-Dichloroethane-d4	0.652	0.622	4.6	103	0.00
34	I	1,4-Difluorobenzene	1.000	1.000	0.0	99	0.00
35	S	Dibromofluoromethane	0.351	0.375	-6.8	111	0.00
36	T	1,1-Dichloropropene	0.457	0.439	3.9	101	0.00
37	T	Ethyl Acetate	0.464	0.456	1.7	102	0.00
38	T	Carbon Tetrachloride	0.516	0.523	-1.4	103	0.00
39	T	Methylcyclohexane	0.577	0.499	13.5	90	0.00
40	TM	Benzene	1.339	1.335	0.3	102	0.00
41	T	Methacrylonitrile	0.239	0.254	-6.3	111	0.00
42	TM	1,2-Dichloroethane	0.497	0.490	1.4	103	-0.01
43	T	Isopropyl Acetate	0.755	0.751	0.5	105	0.00
44	TM	Trichloroethene	0.388	0.403	-3.9	108	0.00
45	C	1,2-Dichloropropane	0.328	0.343	-4.6#	110	0.00
46	T	Dibromomethane	0.257	0.262	-1.9	107	0.00
47	T	Bromodichloromethane	0.494	0.533	-7.9	110	0.00
48	T	Methyl methacrylate	0.347	0.350	-0.9	106	0.00

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LabSampleId :
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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49	T	1,4-Dioxane	0.009	0.008	11.1	89	0.00
50	S	Toluene-d8	1.347	1.294	3.9	98	0.00
51	T	4-Methyl-2-Pentanone	0.485	0.492	-1.4	106	0.00
52	CM	Toluene	0.930	0.912	1.9#	100	0.00
53	T	t-1,3-Dichloropropene	0.525	0.552	-5.1	103	0.00
54	T	cis-1,3-Dichloropropene	0.564	0.588	-4.3	106	0.00
55	T	1,1,2-Trichloroethane	0.378	0.404	-6.9	111	0.00
56	T	Ethyl methacrylate	0.548	0.577	-5.3	105	0.00
57	T	1,3-Dichloropropane	0.610	0.635	-4.1	109	0.00
58	T	2-Chloroethyl Vinyl ether	0.257	0.269	-4.7	101	0.00
59	T	2-Hexanone	0.375	0.380	-1.3	104	0.00
60	T	Dibromochloromethane	0.412	0.453	-10.0	108	0.00
61	T	1,2-Dibromoethane	0.415	0.425	-2.4	102	0.00
62	S	4-Bromofluorobenzene	0.515	0.514	0.2	100	0.00
63	I	Chlorobenzene-d5	1.000	1.000	0.0	95	0.00
64	T	Tetrachloroethene	0.391	0.417	-6.6	109	0.00
65	PM	Chlorobenzene	1.066	1.113	-4.4	103	0.00
66	T	1,1,1,2-Tetrachloroethane	0.397	0.444	-11.8	109	0.00
67	C	Ethyl Benzene	1.788	1.894	-5.9#	101	0.00
68	T	m/p-Xylenes	0.728	0.763	-4.8	99	0.00
69	T	o-Xylene	0.725	0.760	-4.8	100	0.00
70	T	Styrene	1.175	1.300	-10.6	103	0.00
71	P	Bromoform	0.313	0.377	-20.4	113	0.00
72	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
73	T	Isopropylbenzene	3.437	3.577	-4.1	102	0.00
74	T	N-amyl acetate	1.188	1.298	-9.3	107	0.00
75	P	1,1,2,2-Tetrachloroethane	1.115	1.202	-7.8	108	0.00
76	T	1,2,3-Trichloropropane	0.984	1.098	-11.6	108	0.00
77	T	Bromobenzene	0.895	0.935	-4.5	105	0.00
78	T	n-propylbenzene	3.811	4.132	-8.4	103	0.00
79	T	2-Chlorotoluene	2.369	2.463	-4.0	102	0.00
80	T	1,3,5-Trimethylbenzene	2.917	3.048	-4.5	99	0.00
81	T	trans-1,4-Dichloro-2-butene	0.331	0.361	-9.1	103	0.00
82	T	4-Chlorotoluene	2.729	2.858	-4.7	101	0.00
83	T	tert-Butylbenzene	2.979	3.128	-5.0	101	0.00
84	T	1,2,4-Trimethylbenzene	2.900	3.061	-5.6	99	0.00
85	T	sec-Butylbenzene	3.594	3.797	-5.6	100	0.00
86	T	p-Isopropyltoluene	3.100	3.282	-5.9	98	0.00
87	T	1,3-Dichlorobenzene	1.682	1.831	-8.9	104	0.00
88	T	1,4-Dichlorobenzene	1.766	1.846	-4.5	102	0.00
89	T	n-Butylbenzene	2.546	2.744	-7.8	100	0.00
90	T	Hexachloroethane	0.498	0.577	-15.9	103	0.00
91	T	1,2-Dichlorobenzene	1.671	1.822	-9.0	105	0.00
92	T	1,2-Dibromo-3-Chloropropane	0.248	0.255	-2.8	99	0.00
93	T	1,2,4-Trichlorobenzene	0.994	1.156	-16.3	114	0.00
94	T	Hexachlorobutadiene	0.443	0.501	-13.1	114	0.00
95	T	Naphthalene	3.310	3.784	-14.3	106	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.975	1.154	-18.4	114	0.00

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