Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032222\
Data File : BM034312.D
Acq On : 22 Mar 2022 16:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 22 17:30:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 16 02:38:26 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	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2		1,4-Dioxane	0.504	0.474	6.0	98	0.00
3		Pyridine	1.365	1.339	1.9	101	-0.01
4		n-Nitrosodimethylamine	0.648	0.634	2.2	103	0.00
5	S	2-Fluorophenol	1.115	1.111	0.4	101	-0.01
6		Aniline	1.824	1.850	-1.4	103	-0.01
7	S	Phenol-d6	1.596	1.605	-0.6	102	0.00
8		2-Chlorophenol	1.304	1.298	0.5	103	-0.01
9		Benzaldehyde	0.860	0.819	4.8	102	0.00
10	C	Phenol	1.588	1.600	-0.8	103	-0.01
11		bis(2-Chloroethyl)ether	1.303	1.293	0.8	103	0.00
12		1,3-Dichlorobenzene	1.479	1.444	2.4	102	-0.01
13	C	1,4-Dichlorobenzene	1.514	1.474	2.6	103	0.00
14		1,2-Dichlorobenzene	1.476	1.423	3.6	101	0.00
15		Benzyl Alcohol	1.270	1.306	-2.8	103	-0.01
16		2,2'-oxybis(1-Chloropropane	1.788	1.807	-1.1	107	-0.01
17		2-Methylphenol	1.112	1.116	-0.4	103	0.00
18		Hexachloroethane	0.558	0.560	-0.4	104	-0.01
19	P	n-Nitroso-di-n-propylamine	1.155	1.164	-0.8	104	-0.01
20		3+4-Methylphenols	1.538	1.566	-1.8	104	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	103	-0.01
22		Acetophenone	0.513	0.506	1.4	105	-0.01
23	S	Nitrobenzene-d5	0.411	0.411	0.0	105	0.00
24		Nitrobenzene	0.384	0.380	1.0	105	0.00
25		Isophorone	0.687	0.694	-1.0	105	-0.01
26	C	2-Nitrophenol	0.170	0.174	-2.4	104	0.00
27		2,4-Dimethylphenol	0.266	0.268	-0.8	106	0.00
28		bis(2-Chloroethoxy)methane	0.440	0.443	-0.7	105	-0.01
29	C	2,4-Dichlorophenol	0.309	0.312	-1.0	104	0.00
30		1,2,4-Trichlorobenzene	0.352	0.339	3.7	103	-0.01
31		Naphthalene	1.042	1.022	1.9	105	0.00
32		Benzoic acid	0.220	0.219	0.5	102	0.00
33		4-Chloroaniline	0.411	0.426	-3.6	107	-0.01
34	C	Hexachlorobutadiene	0.223	0.213	4.5	103	-0.01
35		Caprolactam	0.108	0.120	-11.1	116	0.00
36	C	4-Chloro-3-methylphenol	0.355	0.372	-4.8	108	0.00
37		2-Methylnaphthalene	0.743	0.741	0.3	107	-0.01
38		1-Methylnaphthalene	0.727	0.734	-1.0	108	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
40		1,2,4,5-Tetrachlorobenzene	0.604	0.563	6.8	106	-0.01
41	P	Hexachlorocyclopentadiene	0.349	0.325	6.9	102	0.00
42	S	2,4,6-Tribromophenol	0.270	0.280	-3.7	118	0.00
43	C	2,4,6-Trichlorophenol	0.417	0.408	2.2	108	0.00
44		2,4,5-Trichlorophenol	0.447	0.441	1.3	108	0.00
45	S	2-Fluorobiphenyl	1.464	1.391	5.0	107	0.00
46		1,1'-Biphenyl	1.525	1.445	5.2	108	-0.01
47		2-Chloronaphthalene	1.171	1.123	4.1	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.358	0.386	-7.8	117	0.00
49		Acenaphthylene	1.824	1.799	1.4	111	0.00
50		Dimethylphthalate	1.522	1.507	1.0	112	0.00
51		2,6-Dinitrotoluene	0.312	0.322	-3.2	115	0.00
52	C	Acenaphthene	1.229	1.216	1.1	111	0.00
53		3-Nitroaniline	0.313	0.337	-7.7	116	0.00
54	P	2,4-Dinitrophenol	0.175	0.188	-7.4	113	0.00
55		Dibenzofuran	1.831	1.798	1.8	112	-0.01
56	P	4-Nitrophenol	0.254	0.281	-10.6	121	0.00
57		2,4-Dinitrotoluene	0.422	0.450	-6.6	117	0.00
58		Fluorene	1.487	1.494	-0.5	114	0.00
59		2,3,4,6-Tetrachlorophenol	0.403	0.408	-1.2	113	0.00
60		Diethylphthalate	1.562	1.605	-2.8	115	0.00
61		4-Chlorophenyl-phenylether	0.759	0.751	1.1	113	-0.01
62		4-Nitroaniline	0.301	0.351	-16.6	123	0.00
63		Azobenzene	1.506	1.568	-4.1	117	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	119	-0.01
65		4,6-Dinitro-2-methylphenol	0.130	0.133	-2.3	119	0.00
66	c	n-Nitrosodiphenylamine	0.635	0.606	4.6	116	0.00
67		4-Bromophenyl-phenylether	0.230	0.219	4.8	117	0.00
68		Hexachlorobenzene	0.256	0.242	5.5	116	0.00
69		Atrazine	0.209	0.217	-3.8	121	0.00
70	C	Pentachlorophenol	0.164	0.166	-1.2	119	0.00
71		Phenanthrene	1.120	1.106	1.3	122	0.00
72		Anthracene	1.090	1.086	0.4	121	-0.01
73		Carbazole	1.017	1.066	-4.8	127	0.00
74		Di-n-butylphthalate	1.264	1.322	-4.6	125	0.00
75	C	Fluoranthene	1.256	1.316	-4.8	130	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	150	0.00
77		Benzidine	0.293	0.266	9.2	128	0.00
78		Pyrene	1.409	1.292	8.3	133	0.00
79	S	Terphenyl-d14	1.151	1.061	7.8	131	0.00
80		Butylbenzylphthalate	0.596	0.599	-0.5	144	0.00
81		Benzo(a)anthracene	1.262	1.251	0.9	153#	0.00
82		3,3'-Dichlorobenzidine	0.422	0.434	-2.8	158#	0.00
83		Chrysene	1.289	1.231	4.5	148	0.00
84		Bis(2-ethylhexyl)phthalate	0.880	0.890	-1.1	150	0.00
85	c	Di-n-octyl phthalate	1.397	1.486	-6.4	166#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	166#	0.00
87		Indeno(1,2,3-cd)pyrene	1.287	1.204	6.4	160#	0.00
88		Benzo(b)fluoranthene	1.196	1.189	0.6	167#	0.00
89		Benzo(k)fluoranthene	1.267	1.323	-4.4	171#	0.00
90	C	Benzo(a)pyrene	1.014	1.022	-0.8	170#	0.00
91		Dibenzo(a,h)anthracene	1.043	0.989	5.2	165#	0.00
92		Benzo(g,h,i)perylene	1.080	0.970	10.2	155#	0.00

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