Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG051319\
Data File : BG040911.D
Acq On : 13 May 2019 15:00
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 14 01:49:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 2019
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	101	-0.03
2		1,4-Dioxane	0.516	0.558	-8.1	114	-0.02
3		Pyridine	1.496	1.687	-12.8	113	-0.03
4		n-Nitrosodimethylamine	0.830	0.840	-1.2	107	-0.02
5	S	2-Fluorophenol	1.135	1.313	-15.7	120	-0.03
6		Aniline	2.315	2.524	-9.0	112	-0.03
7	S	Phenol-d6	1.747	2.031	-16.3	121	-0.02
8		2-Chlorophenol	1.385	1.603	-15.7	119	-0.03
9		Benzaldehyde	1.083	1.142	-5.4	112	-0.03
10	C	Phenol	1.878	2.130	-13.4	117	-0.03
11		bis(2-Chloroethyl)ether	1.408	1.573	-11.7	116	-0.03
12		1,3-Dichlorobenzene	1.512	1.771	-17.1	121	-0.03
13	C	1,4-Dichlorobenzene	1.533	1.801	-17.5	122	-0.03
14		1,2-Dichlorobenzene	1.478	1.725	-16.7	122	-0.03
15		Benzyl Alcohol	1.818	1.907	-4.9	108	-0.03
16		2,2'-oxybis(1-Chloropropane	2.804	2.499	10.9	93	-0.03
17		2-Methylphenol	1.329	1.495	-12.5	117	-0.03
18		Hexachloroethane	0.629	0.702	-11.6	117	-0.03
19	P	n-Nitroso-di-n-propylamine	1.573	1.594	-1.3	107	-0.03
20		3+4-Methylphenols	1.851	2.101	-13.5	117	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	101	-0.03
22		Acetophenone	0.556	0.649	-16.7	118	-0.03
23	S	Nitrobenzene-d5	0.433	0.520	-20.1	123	-0.03
24		Nitrobenzene	0.462	0.531	-14.9	119	-0.03
25		Isophorone	0.964	1.026	-6.4	109	-0.04
26	C	2-Nitrophenol	0.166	0.239	-44.0#	149	-0.03
27		2,4-Dimethylphenol	0.311	0.353	-13.5	116	-0.03
28		bis(2-Chloroethoxy)methane	0.484	0.531	-9.7	111	-0.03
29	C	2,4-Dichlorophenol	0.353	0.441	-24.9#	125	-0.03
30		1,2,4-Trichlorobenzene	0.392	0.498	-27.0#	129	-0.03
31		Naphthalene	1.063	1.233	-16.0	116	-0.03
32		Benzoic acid	0.213	0.294	-38.0#	145	-0.02
33		4-Chloroaniline	0.471	0.546	-15.9	120	-0.03
34	C	Hexachlorobutadiene	0.289	0.372	-28.7#	131	-0.03
35		Caprolactam	0.139	0.158	-13.7	114	-0.02
36	C	4-Chloro-3-methylphenol	0.445	0.530	-19.1	124	-0.02
37		2-Methylnaphthalene	0.794	0.927	-16.8	118	-0.03
38		1-Methylnaphthalene	0.777	0.905	-16.5	118	-0.03
39	I	Acenaphthene-d10	1.000	1.000	0.0	112	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.745	0.912	-22.4	137	-0.03
41	P	Hexachlorocyclopentadiene	0.440	0.426	3.2	110	-0.03
42	S	2,4,6-Tribromophenol	0.275	0.351	-27.6#	145	-0.03
43	C	2,4,6-Trichlorophenol	0.450	0.547	-21.6#	140	-0.03
44		2,4,5-Trichlorophenol	0.490	0.597	-21.8	139	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.385	1.553	-12.1	125	-0.03
46		1,1'-Biphenyl	1.553	1.691	-8.9	122	-0.03
47		2-Chloronaphthalene	1.215	1.361	-12.0	126	-0.03
48		2-Nitroaniline	0.433	0.512	-18.2	136	-0.03
49		Acenaphthylene	2.062	2.219	-7.6	120	-0.03
50		Dimethylphthalate	1.673	1.871	-11.8	125	-0.03
51		2,6-Dinitrotoluene	0.327	0.416	-27.2#	144	-0.03
52	C	Acenaphthene	1.201	1.272	-5.9	122	-0.03
53		3-Nitroaniline	0.335	0.396	-18.2	133	-0.03
54	P	2,4-Dinitrophenol	0.161	0.243	-50.9#	177#	-0.02
55		Dibenzofuran	1.967	2.188	-11.2	127	-0.03
56	P	4-Nitrophenol	0.290	0.319	-10.0	127	-0.02
57		2,4-Dinitrotoluene	0.444	0.592	-33.3#	153#	-0.02
58		Fluorene	1.590	1.729	-8.7	123	-0.03
59		2,3,4,6-Tetrachlorophenol	0.494	0.588	-19.0	134	-0.03
60		Diethylphthalate	1.766	1.875	-6.2	120	-0.03
61		4-Chlorophenyl-phenylether	0.925	1.078	-16.5	132	-0.03
62		4-Nitroaniline	0.374	0.425	-13.6	129	-0.03
63		Azobenzene	1.818	1.748	3.9	108	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	114	-0.03
65		4,6-Dinitro-2-methylphenol	0.092	0.141	-53.3#	188#	-0.03
66	c	n-Nitrosodiphenylamine	0.588	0.627	-6.6	123	-0.03
67		4-Bromophenyl-phenylether	0.249	0.289	-16.1	135	-0.03
68		Hexachlorobenzene	0.258	0.300	-16.3	136	-0.03
69		Atrazine	0.233	0.200	14.2	96	-0.03
70	C	Pentachlorophenol	0.151	0.181	-19.9	137	-0.02
71		Phenanthrene	1.077	1.169	-8.5	124	-0.03
72		Anthracene	1.083	1.170	-8.0	123	-0.03
73		Carbazole	0.987	1.033	-4.7	119	-0.03
74		Di-n-butylphthalate	1.191	1.197	-0.5	115	-0.04
75	C	Fluoranthene	1.342	1.499	-11.7	128	-0.03
76	I	Chrysene-d12	1.000	1.000	0.0	105	-0.03
77		Benzidine	0.548	0.574	-4.7	104	-0.03
78		Pyrene	1.254	1.504	-19.9	124	-0.03
79	S	Terphenyl-d14	0.903	1.081	-19.7	123	-0.03
80		Butylbenzylphthalate	0.489	0.567	-16.0	122	-0.03
81		Benzo(a)anthracene	1.264	1.519	-20.2	125	-0.03
82		3,3'-Dichlorobenzidine	0.451	0.570	-26.4#	130	-0.04
83		Chrysene	1.202	1.471	-22.4	127	-0.03
84		Bis(2-ethylhexyl)phthalate	0.705	0.781	-10.8	116	-0.05
85	c	Di-n-octyl phthalate	1.188	1.337	-12.5	118	-0.08
86		Indeno(1,2,3-cd)pyrene	1.453	1.807	-24.4	131	-0.11
87	I	Perylene-d12	1.000	1.000	0.0	117	-0.08

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.222	1.397	-14.3	134	-0.06
89	Benzo(k)fluoranthene	1.141	1.263	-10.7	130	-0.06
90 C	Benzo(a)pyrene	1.157	1.304	-12.7	131	-0.07
91	Dibenzo(a,h)anthracene	1.171	1.319	-12.6	131	-0.12
92	Benzo(g,h,i)perylene	1.165	1.300	-11.6	130	-0.12

(#) = Out of Range SPCC's out = 0 CCC's out = 4