Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG040918\
Data File : BG033673.D
Acq On : 9 Apr 2018 10:22
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 06:11:07 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 05 17:36:07 2018
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	117	0.00
2		1,4-Dioxane	0.542	0.585	-7.9	128	0.00
3		Pyridine	1.497	1.670	-11.6	118	0.00
4		n-Nitrosodimethylamine	0.614	0.693	-12.9	131	0.00
5	S	2-Fluorophenol	1.067	1.183	-10.9	120	0.00
6		Aniline	2.155	2.269	-5.3	114	0.00
7	S	Phenol-d6	1.833	1.889	-3.1	117	0.00
8		2-Chlorophenol	1.219	1.270	-4.2	113	0.00
9		Benzaldehyde	1.165	0.945	18.9	97	0.00
10	C	Phenol	1.809	1.889	-4.4	116	0.00
11		bis(2-Chloroethyl)ether	1.477	1.550	-4.9	113	0.00
12		1,3-Dichlorobenzene	1.594	1.590	0.3	115	0.00
13	C	1,4-Dichlorobenzene	1.597	1.601	-0.3	118	0.00
14		1,2-Dichlorobenzene	1.558	1.555	0.2	112	0.00
15		Benzyl Alcohol	1.443	1.583	-9.7	120	0.00
16		2,2'-oxybis(1-Chloropropane	2.408	2.764	-14.8	132	0.00
17		2-Methylphenol	1.233	1.338	-8.5	124	0.00
18		Hexachloroethane	0.616	0.625	-1.5	118	0.00
19	P	n-Nitroso-di-n-propylamine	1.343	1.471	-9.5	118	0.00
20		3+4-Methylphenols	1.755	1.839	-4.8	114	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
22		Acetophenone	0.548	0.557	-1.6	109	0.00
23	S	Nitrobenzene-d5	0.435	0.429	1.4	111	0.00
24		Nitrobenzene	0.466	0.466	0.0	112	0.00
25		Isophorone	0.845	0.865	-2.4	115	0.00
26	C	2-Nitrophenol	0.176	0.184	-4.5	111	0.00
27		2,4-Dimethylphenol	0.256	0.274	-7.0	126	0.00
28		bis(2-Chloroethoxy)methane	0.422	0.408	3.3	108	0.00
29	C	2,4-Dichlorophenol	0.315	0.318	-1.0	111	0.00
30		1,2,4-Trichlorobenzene	0.409	0.381	6.8	106	0.00
31		Naphthalene	1.036	1.017	1.8	112	0.00
32		Benzoic acid	0.238	0.208	12.6	113	0.00
33		4-Chloroaniline	0.464	0.417	10.1	99	0.00
34	C	Hexachlorobutadiene	0.310	0.284	8.4	108	0.00
35		Caprolactam	0.123	0.130	-5.7	117	0.00
36	C	4-Chloro-3-methylphenol	0.411	0.404	1.7	106	0.00
37		2-Methylnaphthalene	0.804	0.758	5.7	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.724	0.682	5.8	102	0.00
40	P	Hexachlorocyclopentadiene	0.425	0.350	17.6	87	0.00
41	S	2,4,6-Tribromophenol	0.299	0.250	16.4	88	0.00
42	C	2,4,6-Trichlorophenol	0.421	0.406	3.6	99	0.00
43		2,4,5-Trichlorophenol	0.498	0.461	7.4	92	0.00
44	S	2-Fluorobiphenyl	1.385	1.349	2.6	104	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.537	1.532	0.3	107	0.00
46		2-Chloronaphthalene	1.185	1.167	1.5	105	0.00
47		2-Nitroaniline	0.444	0.472	-6.3	110	0.00
48		Acenaphthylene	1.978	1.950	1.4	106	0.00
49		Dimethylphthalate	1.741	1.729	0.7	107	0.00
50		2,6-Dinitrotoluene	0.356	0.359	-0.8	103	0.00
51	C	Acenaphthene	1.275	1.271	0.3	105	0.00
52		3-Nitroaniline	0.373	0.358	4.0	101	0.00
53	P	2,4-Dinitrophenol	0.149	0.141	5.4	94	0.00
54		Dibenzofuran	1.883	1.787	5.1	102	0.00
55	P	4-Nitrophenol	0.262	0.251	4.2	100	0.00
56		2,4-Dinitrotoluene	0.524	0.502	4.2	102	0.00
57		Fluorene	1.604	1.528	4.7	104	0.00
58		2,3,4,6-Tetrachlorophenol	0.468	0.412	12.0	91	0.00
59		Diethylphthalate	1.690	1.702	-0.7	109	0.00
60		4-Chlorophenyl-phenylether	0.922	0.854	7.4	102	0.00
61		4-Nitroaniline	0.378	0.387	-2.4	109	0.00
62		Azobenzene	1.637	1.615	1.3	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	0.120	0.124	-3.3	102	0.00
65	c	n-Nitrosodiphenylamine	0.571	0.572	-0.2	103	0.00
66		4-Bromophenyl-phenylether	0.235	0.222	5.5	97	0.00
67		Hexachlorobenzene	0.248	0.236	4.8	99	0.00
68		Atrazine	0.231	0.231	0.0	102	0.00
69	C	Pentachlorophenol	0.170	0.139	18.2	84	0.00
70		Phenanthrene	1.042	1.011	3.0	101	0.00
71		Anthracene	1.055	1.037	1.7	102	0.00
72		Carbazole	0.935	0.951	-1.7	104	0.00
73		Di-n-butylphthalate	1.088	1.162	-6.8	109	0.00
74	C	Fluoranthene	1.289	1.238	4.0	99	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76		Benzidine	0.542	0.594	-9.6	101	0.00
77		Pyrene	1.334	1.331	0.2	100	0.00
78	S	Terphenyl-d14	0.935	0.910	2.7	98	0.00
79		Butylbenzylphthalate	0.492	0.538	-9.3	108	0.00
80		Benzo(a)anthracene	1.274	1.228	3.6	97	0.00
81		3,3'-Dichlorobenzidine	0.453	0.468	-3.3	99	0.00
82		Chrysene	1.233	1.199	2.8	98	0.00
83		Bis(2-ethylhexyl)phthalate	0.679	0.755	-11.2	110	0.00
84	c	Di-n-octyl phthalate	1.039	1.202	-15.7	113	0.00
85		Indeno(1,2,3-cd)pyrene	1.333	1.359	-2.0	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Benzo(b)fluoranthene	1.155	1.098	4.9	97	0.00

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88	Benzo(k)fluoranthene	1.169	1.142	2.3	100	0.00
89 C	Benzo(a)pyrene	1.110	1.090	1.8	100	0.00
90	Dibenzo(a,h)anthracene	1.045	1.057	-1.1	100	0.00
91	Benzo(g,h,i)perylene	1.035	1.041	-0.6	101	-0.02

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