Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
Data File : BF092464.D
Acq On : 25 Jan 2017 9:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 25 10:32:59 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 24 13:48:07 2017
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	106	0.00
2		1,4-Dioxane	0.684	0.673	1.6	106	-0.01
3		Pyridine	1.946	1.870	3.9	103	0.00
4		n-Nitrosodimethylamine	0.955	0.909	4.8	101	0.00
5	S	2-Fluorophenol	1.285	1.304	-1.5	110	0.01
6		Aniline	2.442	2.471	-1.2	106	0.00
7	S	Phenol-d6	1.637	1.581	3.4	104	0.00
8		2-Chlorophenol	1.350	1.385	-2.6	109	0.00
9		Benzaldehyde	1.162	1.079	7.1	102	0.00
10	C	Phenol	1.976	1.849	6.4	96	0.00
11		bis(2-Chloroethyl)ether	1.479	1.557	-5.3	107	0.00
12		1,3-Dichlorobenzene	1.597	1.601	-0.3	106	0.00
13	C	1,4-Dichlorobenzene	1.607	1.665	-3.6	105	0.00
14		1,2-Dichlorobenzene	1.525	1.440	5.6	107	0.00
15		Benzyl Alcohol	1.234	1.203	2.5	102	0.00
16		2,2'-oxybis(1-Chloropropane	2.925	2.843	2.8	104	0.00
17		2-Methylphenol	1.154	1.143	1.0	104	0.00
18		Hexachloroethane	0.554	0.566	-2.2	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.115	1.046	6.2	102	0.00
20		3+4-Methylphenols	1.400	1.492	-6.6	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	-0.01
22		Acetophenone	0.476	0.463	2.7	104	0.00
23	S	Nitrobenzene-d5	0.366	0.361	1.4	106	0.00
24		Nitrobenzene	0.384	0.401	-4.4	104	0.00
25		Isophorone	0.721	0.688	4.6	103	0.00
26	C	2-Nitrophenol	0.161	0.170	-5.6	116	0.00
27		2,4-Dimethylphenol	0.334	0.329	1.5	108	0.00
28		bis(2-Chloroethoxy)methane	0.474	0.428	9.7	104	0.00
29	C	2,4-Dichlorophenol	0.290	0.280	3.4	105	0.00
30		1,2,4-Trichlorobenzene	0.313	0.302	3.5	110	0.00
31		Naphthalene	1.024	0.952	7.0	106	0.00
32		Benzoic acid	0.231	0.223	3.5	95	0.00
33		4-Chloroaniline	0.433	0.413	4.6	109	0.00
34	C	Hexachlorobutadiene	0.171	0.173	-1.2	109	0.00
35		Caprolactam	0.097	0.093	4.1	102	0.00
36	C	4-Chloro-3-methylphenol	0.313	0.336	-7.3	111	0.00
37		2-Methylnaphthalene	0.648	0.601	7.3	108	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.504	0.511	-1.4	107	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.269	-6.3	108	0.00
41	S	2,4,6-Tribromophenol	0.136	0.134	1.5	107	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.380	1.8	109	0.00
43		2,4,5-Trichlorophenol	0.354	0.368	-4.0	107	0.00
44	S	2-Fluorobiphenyl	1.082	1.095	-1.2	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.446	1.281	11.4	105	0.00
46		2-Chloronaphthalene	1.181	1.095	7.3	103	0.00
47		2-Nitroaniline	0.398	0.393	1.3	108	0.00
48		Acenaphthylene	1.732	1.738	-0.3	108	0.00
49		Dimethylphthalate	1.272	1.334	-4.9	110	0.00
50		2,6-Dinitrotoluene	0.260	0.292	-12.3	109	0.00
51	C	Acenaphthene	1.076	1.026	4.6	111	0.00
52		3-Nitroaniline	0.326	0.389	-19.3	109	0.00
53	P	2,4-Dinitrophenol	0.093	0.127	-36.6#	132	-0.01
54		Dibenzofuran	1.461	1.392	4.7	109	0.00
55	P	4-Nitrophenol	0.259	0.299	-15.4	109	0.00
56		2,4-Dinitrotoluene	0.345	0.352	-2.0	113	0.00
57		Fluorene	1.084	1.031	4.9	111	0.00
58		2,3,4,6-Tetrachlorophenol	0.265	0.278	-4.9	115	0.00
59		Diethylphthalate	1.220	1.171	4.0	108	0.00
60		4-Chlorophenyl-phenylether	0.526	0.491	6.7	100	-0.01
61		4-Nitroaniline	0.326	0.374	-14.7	120	0.00
62		Azobenzene	1.388	1.316	5.2	98	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.01
64		4,6-Dinitro-2-methylphenol	0.099	0.141	-42.4#	133	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.676	-8.2	121	0.00
66		4-Bromophenyl-phenylether	0.202	0.196	3.0	103	-0.01
67		Hexachlorobenzene	0.204	0.204	0.0	111	0.00
68		Atrazine	0.214	0.196	8.4	92	-0.01
69	C	Pentachlorophenol	0.131	0.144	-9.9	109	0.00
70		Phenanthrene	1.101	1.023	7.1	100	-0.01
71		Anthracene	1.076	1.146	-6.5	115	0.00
72		Carbazole	1.028	0.916	10.9	92	-0.01
73		Di-n-butylphthalate	1.175	1.199	-2.0	110	-0.01
74	C	Fluoranthene	1.077	1.030	4.4	106	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	106	-0.01
76		Benzidine	0.742	0.814	-9.7	110	-0.01
77		Pyrene	1.449	1.312	9.5	108	-0.01
78	S	Terphenyl-d14	0.751	0.669	10.9	107	-0.01
79		Butylbenzylphthalate	0.587	0.624	-6.3	107	-0.02
80		Benzo(a)anthracene	1.075	1.040	3.3	108	-0.01
81		3,3'-Dichlorobenzidine	0.408	0.422	-3.4	107	-0.01
82		Chrysene	1.148	1.099	4.3	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.739	0.755	-2.2	110	-0.01
84	c	Di-n-octyl phthalate	1.342	1.467	-9.3	109	-0.02
85		Indeno(1,2,3-cd)pyrene	0.849	0.876	-3.2	113	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	111	-0.01
87		Benzo(b)fluoranthene	1.284	1.356	-5.6	112	-0.01

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88	Benzo(k)fluoranthene	1.068	0.943	11.7	109	-0.01
89 C	Benzo(a)pyrene	1.086	1.085	0.1	111	-0.01
90	Dibenzo(a,h)anthracene	0.878	0.865	1.5	113	-0.01
91	Benzo(g,h,i)perylene	0.888	0.845	4.8	113	-0.02

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