Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016650.D
Acq On : 23 Apr 2015 11:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 23 15:43:34 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 22 02:10:07 2015
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	132	0.00
2		1,4-Dioxane	0.561	0.520	7.3	118	0.00
3		Pyridine	1.552	1.501	3.3	122	0.00
4		n-Nitrosodimethylamine	0.465	0.444	4.5	123	0.01
5	S	2-Fluorophenol	1.248	1.272	-1.9	131	0.00
6		Aniline	2.408	2.461	-2.2	129	0.00
7	S	Phenol-d6	1.743	1.841	-5.6	132	0.01
8		2-Chlorophenol	1.490	1.562	-4.8	136	0.00
9		Benzaldehyde	0.793	0.709	10.6	132	0.00
10	C	Phenol	1.844	1.916	-3.9	129	0.01
11		bis(2-Chloroethyl)ether	1.444	1.430	1.0	130	0.00
12		1,3-Dichlorobenzene	1.555	1.586	-2.0	135	0.00
13	C	1,4-Dichlorobenzene	1.583	1.619	-2.3	133	0.00
14		1,2-Dichlorobenzene	1.506	1.532	-1.7	133	0.00
15		Benzyl Alcohol	1.109	1.168	-5.3	129	0.00
16		2,2'-oxybis(1-Chloropropane	1.398	1.364	2.4	129	0.00
17		2-Methylphenol	1.243	1.297	-4.3	132	0.00
18		Hexachloroethane	0.569	0.579	-1.8	134	0.00
19	P	n-Nitroso-di-n-propylamine	1.127	1.158	-2.8	132	0.00
20		3+4-Methylphenols	1.712	1.841	-7.5	133	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	135	0.00
22		Acetophenone	0.445	0.448	-0.7	131	0.00
23	S	Nitrobenzene-d5	0.325	0.323	0.6	129	0.00
24		Nitrobenzene	0.338	0.335	0.9	129	0.00
25		Isophorone	0.665	0.680	-2.3	130	0.00
26	C	2-Nitrophenol	0.187	0.207	-10.7	137	0.00
27		2,4-Dimethylphenol	0.317	0.334	-5.4	134	0.00
28		bis(2-Chloroethoxy)methane	0.399	0.404	-1.3	134	0.00
29	C	2,4-Dichlorophenol	0.262	0.296	-13.0	141	0.00
30		1,2,4-Trichlorobenzene	0.279	0.292	-4.7	138	0.00
31		Naphthalene	1.044	1.044	0.0	131	0.00
32		Benzoic acid	0.103	0.191	-85.4#	233#	0.03
33		4-Chloroaniline	0.477	0.503	-5.5	131	0.00
34	C	Hexachlorobutadiene	0.124	0.134	-8.1	149	0.00
35		Caprolactam	0.152	0.158	-3.9	118	0.03
36	C	4-Chloro-3-methylphenol	0.317	0.342	-7.9	131	0.01
37		2-Methylnaphthalene	0.747	0.770	-3.1	133	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	130	0.00
39		1,2,4,5-Tetrachlorobenzene	0.460	0.483	-5.0	140	0.00
40	P	Hexachlorocyclopentadiene	0.143	0.151	-5.6	140	0.00
41	S	2,4,6-Tribromophenol	0.136	0.152	-11.8	127	0.00
42	C	2,4,6-Trichlorophenol	0.326	0.374	-14.7	140	0.00
43		2,4,5-Trichlorophenol	0.347	0.395	-13.8	137	0.00
44	S	2-Fluorobiphenyl	1.208	1.239	-2.6	134	0.00

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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)
45		1,1'-Biphenyl	1.540	1.569	-1.9	132	0.00
46		2-Chloronaphthalene	1.180	1.203	-1.9	131	0.00
47		2-Nitroaniline	0.344	0.352	-2.3	118	0.00
48		Acenaphthylene	2.094	2.112	-0.9	125	0.00
49		Dimethylphthalate	1.479	1.488	-0.6	122	0.00
50		2,6-Dinitrotoluene	0.363	0.377	-3.9	124	0.00
51	C	Acenaphthene	1.250	1.256	-0.5	127	0.00
52		3-Nitroaniline	0.449	0.464	-3.3	117	0.00
53	P	2,4-Dinitrophenol	0.133	0.147	-10.5	123	0.00
54		Dibenzofuran	1.759	1.756	0.2	123	0.00
55	P	4-Nitrophenol	0.340	0.307	9.7	101	0.02
56		2,4-Dinitrotoluene	0.500	0.516	-3.2	114	0.00
57		Fluorene	1.550	1.545	0.3	121	0.00
58		2,3,4,6-Tetrachlorophenol	0.262	0.291	-11.1	127	0.00
59		Diethylphthalate	1.576	1.518	3.7	114	0.00
60		4-Chlorophenyl-phenylether	0.636	0.648	-1.9	128	0.00
61		4-Nitroaniline	0.514	0.488	5.1	103	0.01
62		Azobenzene	1.462	1.413	3.4	115	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
64		4,6-Dinitro-2-methylphenol	0.129	0.137	-6.2	111	0.00
65	c	n-Nitrosodiphenylamine	0.671	0.708	-5.5	118	0.00
66		4-Bromophenyl-phenylether	0.157	0.171	-8.9	123	0.00
67		Hexachlorobenzene	0.164	0.175	-6.7	120	0.00
68		Atrazine	0.187	0.193	-3.2	110	0.00
69	C	Pentachlorophenol	0.073	0.085	-16.4	122	0.00
70		Phenanthrene	1.139	1.146	-0.6	111	0.00
71		Anthracene	1.131	1.150	-1.7	111	0.00
72		Carbazole	1.154	1.132	1.9	105	0.00
73		Di-n-butylphthalate	1.405	1.372	2.3	104	0.00
74	C	Fluoranthene	1.226	1.179	3.8	105	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	106	0.00
76		Benzidine	0.740	0.673	9.1	94	0.00
77		Pyrene	1.343	1.416	-5.4	104	0.00
78	S	Terphenyl-d14	0.830	0.862	-3.9	106	0.00
79		Butylbenzylphthalate	0.703	0.739	-5.1	104	0.00
80		Benzo(a)anthracene	1.206	1.226	-1.7	105	0.00
81		3,3'-Dichlorobenzidine	0.399	0.421	-5.5	108	0.00
82		Chrysene	1.158	1.159	-0.1	103	0.00
83		Bis(2-ethylhexyl)phthalate	0.965	1.012	-4.9	103	0.00
84	c	Di-n-octyl phthalate	1.603	1.687	-5.2	103	0.00
85		Indeno(1,2,3-cd)pyrene	1.271	1.315	-3.5	106	0.01
86	I	Perylene-d12	1.000	1.000	0.0	106	0.00
87		Benzo(b)fluoranthene	1.193	1.237	-3.7	108	0.00

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88	Benzo(k)fluoranthene	1.171	1.172	-0.1	101	0.00
89 C	Benzo(a)pyrene	1.133	1.162	-2.6	105	0.00
90	Dibenzo(a,h)anthracene	1.101	1.134	-3.0	105	0.00
91	Benzo(g,h,i)perylene	1.111	1.142	-2.8	104	0.00

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