Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021816\
Data File : BG020952.D
Acq On : 19 Feb 2016 1:43
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 19 02:50:00 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 2016
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	103	0.00
2		1,4-Dioxane	0.432	0.439	-1.6	107	-0.01
3		Pyridine	1.277	1.326	-3.8	103	-0.01
4		n-Nitrosodimethylamine	0.334	0.342	-2.4	103	-0.01
5	S	2-Fluorophenol	1.188	1.206	-1.5	104	0.00
6		Aniline	2.164	2.266	-4.7	106	0.00
7	S	Phenol-d6	1.736	1.828	-5.3	107	0.00
8		2-Chlorophenol	1.484	1.554	-4.7	108	-0.01
9		Benzaldehyde	0.869	0.855	1.6	100	0.00
10	C	Phenol	1.837	1.891	-2.9	106	0.00
11		bis(2-Chloroethyl)ether	1.467	1.528	-4.2	105	-0.01
12		1,3-Dichlorobenzene	1.569	1.618	-3.1	107	-0.01
13	C	1,4-Dichlorobenzene	1.590	1.647	-3.6	106	-0.01
14		1,2-Dichlorobenzene	1.547	1.594	-3.0	105	-0.01
15		Benzyl Alcohol	1.094	1.140	-4.2	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.530	1.522	0.5	102	-0.01
17		2-Methylphenol	1.323	1.395	-5.4	108	0.00
18		Hexachloroethane	0.547	0.564	-3.1	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.120	1.181	-5.4	108	-0.01
20		3+4-Methylphenols	1.844	1.957	-6.1	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	-0.01
22		Acetophenone	0.482	0.494	-2.5	108	0.00
23	S	Nitrobenzene-d5	0.304	0.307	-1.0	107	0.00
24		Nitrobenzene	0.317	0.317	0.0	107	0.00
25		Isophorone	0.641	0.667	-4.1	110	0.00
26	C	2-Nitrophenol	0.167	0.184	-10.2	113	0.00
27		2,4-Dimethylphenol	0.320	0.317	0.9	103	0.00
28		bis(2-Chloroethoxy)methane	0.395	0.407	-3.0	110	-0.01
29	C	2,4-Dichlorophenol	0.287	0.299	-4.2	112	0.00
30		1,2,4-Trichlorobenzene	0.289	0.298	-3.1	110	0.00
31		Naphthalene	1.035	1.052	-1.6	109	-0.01
32		Benzoic acid	0.256	0.248	3.1	105	0.00
33		4-Chloroaniline	0.445	0.467	-4.9	112	0.00
34	C	Hexachlorobutadiene	0.147	0.152	-3.4	110	-0.01
35		Caprolactam	0.110	0.119	-8.2	116	0.00
36	C	4-Chloro-3-methylphenol	0.328	0.342	-4.3	112	0.00
37		2-Methylnaphthalene	0.726	0.738	-1.7	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.587	0.574	2.2	111	0.00
40	P	Hexachlorocyclopentadiene	0.258	0.272	-5.4	116	-0.01
41	S	2,4,6-Tribromophenol	0.197	0.213	-8.1	123	0.00
42	C	2,4,6-Trichlorophenol	0.394	0.411	-4.3	115	-0.01
43		2,4,5-Trichlorophenol	0.414	0.428	-3.4	116	0.00
44	S	2-Fluorobiphenyl	1.424	1.428	-0.3	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.764	1.751	0.7	113	-0.01
46		2-Chloronaphthalene	1.263	1.271	-0.6	115	-0.01
47		2-Nitroaniline	0.303	0.309	-2.0	115	0.00
48		Acenaphthylene	2.196	2.220	-1.1	115	0.00
49		Dimethylphthalate	1.564	1.608	-2.8	116	0.00
50		2,6-Dinitrotoluene	0.335	0.357	-6.6	120	0.00
51	C	Acenaphthene	1.332	1.359	-2.0	116	-0.01
52		3-Nitroaniline	0.389	0.406	-4.4	118	0.00
53	P	2,4-Dinitrophenol	0.116	0.135	-16.4	132	0.00
54		Dibenzofuran	1.762	1.792	-1.7	117	-0.01
55	P	4-Nitrophenol	0.294	0.321	-9.2	121	0.00
56		2,4-Dinitrotoluene	0.433	0.477	-10.2	122	0.00
57		Fluorene	1.622	1.666	-2.7	118	-0.01
58		2,3,4,6-Tetrachlorophenol	0.316	0.336	-6.3	121	-0.01
59		Diethylphthalate	1.608	1.671	-3.9	119	0.00
60		4-Chlorophenyl-phenylether	0.720	0.732	-1.7	119	0.00
61		4-Nitroaniline	0.391	0.413	-5.6	119	0.00
62		Azobenzene	1.264	1.249	1.2	113	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
64		4,6-Dinitro-2-methylphenol	0.094	0.107	-13.8	134	0.00
65	c	n-Nitrosodiphenylamine	0.646	0.656	-1.5	119	-0.01
66		4-Bromophenyl-phenylether	0.211	0.216	-2.4	120	-0.01
67		Hexachlorobenzene	0.225	0.232	-3.1	123	-0.01
68		Atrazine	0.196	0.195	0.5	119	-0.01
69	C	Pentachlorophenol	0.129	0.130	-0.8	120	-0.01
70		Phenanthrene	1.139	1.162	-2.0	122	-0.01
71		Anthracene	1.156	1.166	-0.9	120	-0.01
72		Carbazole	1.037	1.036	0.1	119	-0.01
73		Di-n-butylphthalate	1.330	1.330	0.0	118	-0.01
74	C	Fluoranthene	1.230	1.226	0.3	118	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	119	-0.02
76		Benzidine	0.647	0.592	8.5	106	0.00
77		Pyrene	1.319	1.326	-0.5	120	-0.01
78	S	Terphenyl-d14	0.857	0.861	-0.5	119	-0.01
79		Butylbenzylphthalate	0.632	0.635	-0.5	118	-0.02
80		Benzo(a)anthracene	1.194	1.214	-1.7	122	-0.02
81		3,3'-Dichlorobenzidine	0.443	0.449	-1.4	120	-0.01
82		Chrysene	1.123	1.135	-1.1	121	-0.02
83		Bis(2-ethylhexyl)phthalate	0.936	0.946	-1.1	119	-0.03
84	c	Di-n-octyl phthalate	1.537	1.567	-2.0	120	-0.03
85		Indeno(1,2,3-cd)pyrene	1.291	1.358	-5.2	124	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	121	-0.04
87		Benzo(b)fluoranthene	1.224	1.246	-1.8	122	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.199	1.216	-1.4	123	-0.03
89 C	Benzo(a)pyrene	1.166	1.183	-1.5	122	-0.04
90	Dibenzo(a,h)anthracene	1.134	1.153	-1.7	123	-0.06
91	Benzo(g,h,i)perylene	1.126	1.152	-2.3	123	-0.07

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