Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022226.D
Acq On : 8 Jun 2016 7:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 07:59:43 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 06 16:12:45 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	109	0.00
2		1,4-Dioxane	0.496	0.481	3.0	114	0.00
3		Pyridine	1.340	1.359	-1.4	109	0.00
4		n-Nitrosodimethylamine	0.300	0.330	-10.0	115	0.00
5	S	2-Fluorophenol	1.034	1.109	-7.3	112	0.00
6		Aniline	2.066	2.231	-8.0	110	0.00
7	S	Phenol-d6	1.626	1.762	-8.4	112	0.00
8		2-Chlorophenol	1.430	1.478	-3.4	109	0.00
9		Benzaldehyde	0.895	0.937	-4.7	107	0.00
10	C	Phenol	1.677	1.859	-10.9	116	0.00
11		bis(2-Chloroethyl)ether	1.337	1.388	-3.8	109	0.00
12		1,3-Dichlorobenzene	1.533	1.536	-0.2	110	0.00
13	C	1,4-Dichlorobenzene	1.527	1.535	-0.5	109	0.00
14		1,2-Dichlorobenzene	1.539	1.568	-1.9	112	0.00
15		Benzyl Alcohol	1.051	1.088	-3.5	111	0.00
16		2,2'-oxybis(1-Chloropropane	1.387	1.429	-3.0	113	0.00
17		2-Methylphenol	1.251	1.346	-7.6	117	0.00
18		Hexachloroethane	0.558	0.571	-2.3	113	0.00
19	P	n-Nitroso-di-n-propylamine	1.101	1.154	-4.8	108	0.00
20		3+4-Methylphenols	1.795	1.857	-3.5	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.431	0.441	-2.3	110	0.00
23	S	Nitrobenzene-d5	0.287	0.299	-4.2	111	0.00
24		Nitrobenzene	0.288	0.304	-5.6	114	0.00
25		Isophorone	0.671	0.646	3.7	107	0.00
26	C	2-Nitrophenol	0.156	0.153	1.9	102	0.00
27		2,4-Dimethylphenol	0.283	0.288	-1.8	110	0.00
28		bis(2-Chloroethoxy)methane	0.385	0.381	1.0	108	0.00
29	C	2,4-Dichlorophenol	0.262	0.275	-5.0	109	0.00
30		1,2,4-Trichlorobenzene	0.293	0.288	1.7	107	0.00
31		Naphthalene	0.998	0.978	2.0	111	0.00
32		Benzoic acid	0.087	0.086	1.1	103	0.01
33		4-Chloroaniline	0.415	0.413	0.5	107	0.00
34	C	Hexachlorobutadiene	0.157	0.152	3.2	111	0.00
35		Caprolactam	0.144	0.125	13.2	96	0.00
36	C	4-Chloro-3-methylphenol	0.311	0.322	-3.5	111	0.00
37		2-Methylnaphthalene	0.737	0.731	0.8	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.515	0.528	-2.5	108	0.00
40	P	Hexachlorocyclopentadiene	0.233	0.229	1.7	102	0.00
41	S	2,4,6-Tribromophenol	0.226	0.215	4.9	99	0.00
42	C	2,4,6-Trichlorophenol	0.345	0.360	-4.3	106	0.00
43		2,4,5-Trichlorophenol	0.382	0.390	-2.1	108	0.00
44	S	2-Fluorobiphenyl	1.312	1.343	-2.4	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.505	1.560	-3.7	107	0.00
46		2-Chloronaphthalene	1.154	1.185	-2.7	106	0.00
47		2-Nitroaniline	0.275	0.269	2.2	100	0.00
48		Acenaphthylene	2.024	2.034	-0.5	107	0.00
49		Dimethylphthalate	1.658	1.576	4.9	103	0.00
50		2,6-Dinitrotoluene	0.360	0.364	-1.1	104	0.00
51	C	Acenaphthene	1.213	1.214	-0.1	107	0.00
52		3-Nitroaniline	0.400	0.392	2.0	110	0.00
53	P	2,4-Dinitrophenol	0.108	0.118	-9.3	104	0.00
54		Dibenzofuran	1.893	1.874	1.0	105	0.00
55	P	4-Nitrophenol	0.276	0.263	4.7	93	0.00
56		2,4-Dinitrotoluene	0.484	0.486	-0.4	100	0.00
57		Fluorene	1.631	1.616	0.9	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.342	0.335	2.0	106	0.00
59		Diethylphthalate	1.771	1.663	6.1	102	0.00
60		4-Chlorophenyl-phenylether	0.738	0.741	-0.4	105	0.00
61		4-Nitroaniline	0.424	0.403	5.0	100	0.00
62		Azobenzene	1.341	1.378	-2.8	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	0.092	0.092	0.0	94	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.608	-5.6	104	0.00
66		4-Bromophenyl-phenylether	0.187	0.193	-3.2	102	0.00
67		Hexachlorobenzene	0.218	0.218	0.0	101	0.00
68		Atrazine	0.213	0.204	4.2	97	0.00
69	C	Pentachlorophenol	0.084	0.091	-8.3	109	0.00
70		Phenanthrene	1.086	1.104	-1.7	104	0.00
71		Anthracene	1.077	1.104	-2.5	103	0.00
72		Carbazole	1.020	1.001	1.9	99	0.00
73		Di-n-butylphthalate	1.334	1.285	3.7	99	0.00
74	C	Fluoranthene	1.249	1.200	3.9	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.523	0.489	6.5	82	0.00
77		Pyrene	1.243	1.282	-3.1	100	0.00
78	S	Terphenyl-d14	0.842	0.856	-1.7	98	0.00
79		Butylbenzylphthalate	0.577	0.578	-0.2	97	0.00
80		Benzo(a)anthracene	1.121	1.147	-2.3	101	0.00
81		3,3'-Dichlorobenzidine	0.315	0.318	-1.0	95	0.00
82		Chrysene	1.091	1.108	-1.6	102	0.00
83		Bis(2-ethylhexyl)phthalate	0.848	0.857	-1.1	99	0.00
84	c	Di-n-octyl phthalate	1.408	1.453	-3.2	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.224	1.298	-6.0	102	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	100	0.00
87		Benzo(b)fluoranthene	1.166	1.180	-1.2	100	0.00

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88	Benzo(k)fluoranthene	1.148	1.204	-4.9	104	0.00
89 C	Benzo(a)pyrene	1.115	1.158	-3.9	102	0.00
90	Dibenzo(a,h)anthracene	1.050	1.107	-5.4	102	-0.01
91	Benzo(g,h,i)perylene	1.093	1.146	-4.8	104	0.00

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