Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041316\
Data File : BG021628.D
Acq On : 13 Apr 2016 19:33
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 14 02:22:06 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 13 15:48:13 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	100	0.00
2		1,4-Dioxane	0.535	0.537	-0.4	99	0.00
3		Pyridine	1.604	1.544	3.7	99	0.00
4		n-Nitrosodimethylamine	0.795	0.764	3.9	101	0.00
5	S	2-Fluorophenol	1.201	1.190	0.9	100	0.00
6		Aniline	2.373	2.308	2.7	101	0.00
7	S	Phenol-d6	1.794	1.781	0.7	101	0.00
8		2-Chlorophenol	1.440	1.428	0.8	101	0.00
9		Benzaldehyde	1.037	0.985	5.0	101	0.00
10	C	Phenol	1.930	1.904	1.3	100	0.00
11		bis(2-Chloroethyl)ether	1.544	1.520	1.6	102	0.00
12		1,3-Dichlorobenzene	1.608	1.582	1.6	99	0.00
13	C	1,4-Dichlorobenzene	1.637	1.610	1.6	99	0.00
14		1,2-Dichlorobenzene	1.570	1.537	2.1	100	0.00
15		Benzyl Alcohol	1.371	1.326	3.3	100	0.00
16		2,2'-oxybis(1-Chloropropane	3.308	3.203	3.2	99	0.00
17		2-Methylphenol	1.334	1.316	1.3	102	0.00
18		Hexachloroethane	0.596	0.596	0.0	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.394	1.330	4.6	101	0.00
20		3+4-Methylphenols	1.826	1.792	1.9	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.557	0.542	2.7	100	0.00
23	S	Nitrobenzene-d5	0.389	0.387	0.5	101	0.00
24		Nitrobenzene	0.417	0.410	1.7	100	0.00
25		Isophorone	0.852	0.789	7.4	99	0.00
26	C	2-Nitrophenol	0.159	0.156	1.9	101	0.00
27		2,4-Dimethylphenol	0.361	0.340	5.8	100	0.00
28		bis(2-Chloroethoxy)methane	0.453	0.435	4.0	99	0.00
29	C	2,4-Dichlorophenol	0.308	0.306	0.6	101	0.00
30		1,2,4-Trichlorobenzene	0.331	0.330	0.3	101	0.00
31		Naphthalene	1.070	1.046	2.2	100	0.00
32		Benzoic acid	0.192	0.220	-14.6	123	0.00
33		4-Chloroaniline	0.463	0.447	3.5	102	0.00
34	C	Hexachlorobutadiene	0.214	0.212	0.9	101	0.00
35		Caprolactam	0.141	0.134	5.0	99	0.00
36	C	4-Chloro-3-methylphenol	0.387	0.371	4.1	101	0.00
37		2-Methylnaphthalene	0.735	0.722	1.8	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	0.625	0.629	-0.6	102	0.00
40	P	Hexachlorocyclopentadiene	0.347	0.358	-3.2	101	0.00
41	S	2,4,6-Tribromophenol	0.216	0.213	1.4	103	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.407	-2.0	103	0.00
43		2,4,5-Trichlorophenol	0.430	0.421	2.1	98	0.00
44	S	2-Fluorobiphenyl	1.365	1.352	1.0	101	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.685	1.657	1.7	100	0.00
46		2-Chloronaphthalene	1.246	1.238	0.6	102	0.00
47		2-Nitroaniline	0.427	0.424	0.7	103	0.00
48		Acenaphthylene	2.060	2.023	1.8	101	0.00
49		Dimethylphthalate	1.729	1.656	4.2	102	0.00
50		2,6-Dinitrotoluene	0.330	0.330	0.0	102	0.00
51	C	Acenaphthene	1.317	1.308	0.7	102	0.00
52		3-Nitroaniline	0.366	0.367	-0.3	104	0.00
53	P	2,4-Dinitrophenol	0.106	0.112	-5.7	112	0.00
54		Dibenzofuran	1.798	1.758	2.2	102	0.00
55	P	4-Nitrophenol	0.313	0.331	-5.8	105	0.00
56		2,4-Dinitrotoluene	0.446	0.470	-5.4	108	0.00
57		Fluorene	1.606	1.558	3.0	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.355	0.6	104	0.00
59		Diethylphthalate	1.805	1.732	4.0	102	0.00
60		4-Chlorophenyl-phenylether	0.791	0.766	3.2	102	0.00
61		4-Nitroaniline	0.402	0.421	-4.7	104	0.00
62		Azobenzene	1.547	1.485	4.0	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	0.077	0.082	-6.5	105	0.00
65	c	n-Nitrosodiphenylamine	0.600	0.595	0.8	101	0.00
66		4-Bromophenyl-phenylether	0.214	0.216	-0.9	102	0.00
67		Hexachlorobenzene	0.226	0.224	0.9	104	0.00
68		Atrazine	0.237	0.249	-5.1	101	0.00
69	C	Pentachlorophenol	0.129	0.137	-6.2	105	0.00
70		Phenanthrene	1.102	1.096	0.5	103	0.00
71		Anthracene	1.119	1.131	-1.1	103	0.00
72		Carbazole	1.044	1.068	-2.3	103	0.00
73		Di-n-butylphthalate	1.329	1.383	-4.1	101	0.00
74	C	Fluoranthene	1.316	1.364	-3.6	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.623	0.648	-4.0	105	0.00
77		Pyrene	1.153	1.144	0.8	103	0.00
78	S	Terphenyl-d14	0.802	0.791	1.4	103	0.00
79		Butylbenzylphthalate	0.535	0.531	0.7	100	0.00
80		Benzo(a)anthracene	1.151	1.130	1.8	100	0.00
81		3,3'-Dichlorobenzidine	0.911	0.902	1.0	101	0.00
82		Chrysene	1.106	1.085	1.9	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.783	0.770	1.7	101	0.00
84	c	Di-n-octyl phthalate	1.312	1.309	0.2	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.315	1.306	0.7	101	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	101	0.00
87		Benzo(b)fluoranthene	1.160	1.161	-0.1	99	0.00

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88	Benzo(k)fluoranthene	1.135	1.154	-1.7	104	0.00
89 C	Benzo(a)pyrene	1.124	1.124	0.0	102	0.00
90	Dibenzo(a,h)anthracene	1.127	1.131	-0.4	101	-0.01
91	Benzo(g,h,i)perylene	1.106	1.108	-0.2	101	0.00

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