Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG110316\
Data File : BG024646.D
Acq On : 3 Nov 2016 10:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 03 14:21:08 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 26 11:14:13 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	111	0.00
2		1,4-Dioxane	0.499	0.505	-1.2	117	0.00
3		Pyridine	1.493	1.531	-2.5	118	0.00
4		n-Nitrosodimethylamine	0.773	0.766	0.9	112	0.00
5	S	2-Fluorophenol	1.220	1.210	0.8	116	0.00
6		Aniline	2.121	2.057	3.0	110	0.00
7	S	Phenol-d6	1.674	1.670	0.2	114	0.00
8		2-Chlorophenol	1.241	1.200	3.3	114	0.00
9		Benzaldehyde	1.034	1.001	3.2	111	0.00
10	C	Phenol	1.725	1.759	-2.0	117	0.00
11		bis(2-Chloroethyl)ether	1.296	1.278	1.4	116	0.00
12		1,3-Dichlorobenzene	1.536	1.483	3.5	115	0.00
13	C	1,4-Dichlorobenzene	1.517	1.531	-0.9	117	0.00
14		1,2-Dichlorobenzene	1.459	1.439	1.4	115	0.00
15		Benzyl Alcohol	1.557	1.505	3.3	111	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.320	3.5	111	0.00
17		2-Methylphenol	1.143	1.102	3.6	112	0.00
18		Hexachloroethane	0.583	0.593	-1.7	123	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.164	5.7	105	0.00
20		3+4-Methylphenols	1.535	1.557	-1.4	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.514	0.506	1.6	110	0.00
23	S	Nitrobenzene-d5	0.439	0.447	-1.8	115	0.00
24		Nitrobenzene	0.489	0.493	-0.8	112	0.00
25		Isophorone	0.817	0.796	2.6	107	0.00
26	C	2-Nitrophenol	0.181	0.191	-5.5	118	0.00
27		2,4-Dimethylphenol	0.318	0.324	-1.9	111	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.420	0.7	113	0.00
29	C	2,4-Dichlorophenol	0.333	0.340	-2.1	112	0.00
30		1,2,4-Trichlorobenzene	0.395	0.399	-1.0	114	0.00
31		Naphthalene	0.998	1.009	-1.1	112	0.00
32		Benzoic acid	0.264	0.159	39.8#	67	0.01
33		4-Chloroaniline	0.433	0.433	0.0	106	0.00
34	C	Hexachlorobutadiene	0.309	0.314	-1.6	118	0.00
35		Caprolactam	0.122	0.128	-4.9	114	0.00
36	C	4-Chloro-3-methylphenol	0.402	0.416	-3.5	111	0.00
37		2-Methylnaphthalene	0.728	0.722	0.8	111	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.696	-3.1	115	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.361	5.0	102	0.00
41	S	2,4,6-Tribromophenol	0.299	0.314	-5.0	107	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.413	-2.0	111	0.00
43		2,4,5-Trichlorophenol	0.438	0.448	-2.3	109	0.00
44	S	2-Fluorobiphenyl	1.203	1.205	-0.2	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.383	0.4	109	0.00
46		2-Chloronaphthalene	1.057	1.073	-1.5	112	0.00
47		2-Nitroaniline	0.433	0.451	-4.2	104	0.00
48		Acenaphthylene	1.621	1.623	-0.1	106	0.00
49		Dimethylphthalate	1.420	1.415	0.4	104	0.00
50		2,6-Dinitrotoluene	0.303	0.316	-4.3	105	0.00
51	C	Acenaphthene	1.208	1.097	9.2	97	0.00
52		3-Nitroaniline	0.314	0.333	-6.1	109	0.00
53	P	2,4-Dinitrophenol	0.220	0.209	5.0	94	0.00
54		Dibenzofuran	1.670	1.678	-0.5	107	0.00
55	P	4-Nitrophenol	0.273	0.264	3.3	100	0.00
56		2,4-Dinitrotoluene	0.440	0.476	-8.2	106	0.00
57		Fluorene	1.385	1.412	-1.9	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.460	-1.3	104	0.00
59		Diethylphthalate	1.489	1.461	1.9	102	0.00
60		4-Chlorophenyl-phenylether	0.777	0.800	-3.0	110	0.00
61		4-Nitroaniline	0.343	0.361	-5.2	108	0.00
62		Azobenzene	1.485	1.444	2.8	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.134	2.9	101	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.538	1.6	106	0.00
66		4-Bromophenyl-phenylether	0.243	0.245	-0.8	109	0.00
67		Hexachlorobenzene	0.268	0.273	-1.9	108	0.00
68		Atrazine	0.235	0.229	2.6	103	0.00
69	C	Pentachlorophenol	0.177	0.186	-5.1	107	0.00
70		Phenanthrene	1.019	1.016	0.3	108	0.00
71		Anthracene	1.028	1.019	0.9	106	0.00
72		Carbazole	0.938	0.936	0.2	109	0.00
73		Di-n-butylphthalate	1.081	1.061	1.9	105	0.00
74	C	Fluoranthene	1.289	1.311	-1.7	107	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76		Benzidine	0.471	0.535	-13.6	121	0.00
77		Pyrene	0.978	1.001	-2.4	106	0.00
78	S	Terphenyl-d14	0.669	0.655	2.1	105	0.00
79		Butylbenzylphthalate	0.408	0.406	0.5	106	0.00
80		Benzo(a)anthracene	1.099	1.103	-0.4	110	0.00
81		3,3'-Dichlorobenzidine	0.443	0.447	-0.9	108	0.00
82		Chrysene	1.046	1.055	-0.9	108	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.571	2.1	105	-0.01
84	c	Di-n-octyl phthalate	0.968	0.973	-0.5	107	-0.01
85		Indeno(1,2,3-cd)pyrene	1.444	1.396	3.3	108	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	100	-0.02
87		Benzo(b)fluoranthene	1.081	1.104	-2.1	110	0.00

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88	Benzo(k)fluoranthene	1.044	1.040	0.4	106	-0.01
89 C	Benzo(a)pyrene	1.056	1.065	-0.9	108	-0.01
90	Dibenzo(a,h)anthracene	1.159	1.138	1.8	109	-0.03
91	Benzo(g,h,i)perylene	1.158	1.118	3.5	109	-0.03

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