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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	132	0.00
2		1,4-Dioxane	40.000	37.079	7.3	118	0.00
3		Pyridine	40.000	38.674	3.3	122	0.00
4		n-Nitrosodimethylamine	40.000	38.242	4.4	123	0.01
5	S	2-Fluorophenol	80.000	81.571	-2.0	131	0.00
6		Aniline	40.000	40.882	-2.2	129	0.00
7	S	Phenol-d6	80.000	84.465	-5.6	132	0.01
8		2-Chlorophenol	40.000	41.945	-4.9	136	0.00
9		Benzaldehyde	40.000	35.751	10.6	132	0.00
10	C	Phenol	40.000	41.553	-3.9	129	0.01
11		bis(2-Chloroethyl)ether	40.000	39.608	1.0	130	0.00
12		1,3-Dichlorobenzene	40.000	40.777	-1.9	135	0.00
13	C	1,4-Dichlorobenzene	40.000	40.904	-2.3	133	0.00
14		1,2-Dichlorobenzene	40.000	40.686	-1.7	133	0.00
15		Benzyl Alcohol	40.000	42.128	-5.3	129	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.022	2.4	129	0.00
17		2-Methylphenol	40.000	41.745	-4.4	132	0.00
18		Hexachloroethane	40.000	40.743	-1.9	134	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.103	-2.8	132	0.00
20		3+4-Methylphenols	40.000	43.020	-7.6	133	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	135	0.00
22		Acetophenone	40.000	40.240	-0.6	131	0.00
23	S	Nitrobenzene-d5	80.000	79.491	0.6	129	0.00
24		Nitrobenzene	40.000	39.685	0.8	129	0.00
25		Isophorone	40.000	40.915	-2.3	130	0.00
26	C	2-Nitrophenol	40.000	44.287	-10.7	137	0.00
27		2,4-Dimethylphenol	40.000	42.111	-5.3	134	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.559	-1.4	134	0.00
29	C	2,4-Dichlorophenol	40.000	45.074	-12.7	141	0.00
30		1,2,4-Trichlorobenzene	40.000	41.895	-4.7	138	0.00
31		Naphthalene	40.000	40.008	-0.0	131	0.00
32		Benzoic acid	40.000	59.444	-48.6#	233	0.03
33		4-Chloroaniline	40.000	42.124	-5.3	131	0.00
34	C	Hexachlorobutadiene	40.000	43.085	-7.7	149	0.00
35		Caprolactam	40.000	41.591	-4.0	118	0.03
36	C	4-Chloro-3-methylphenol	40.000	43.104	-7.8	131	0.01
37		2-Methylnaphthalene	40.000	41.220	-3.0	133	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	130	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	42.035	-5.1	140	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.786	-4.5	140	0.00
41	S	2,4,6-Tribromophenol	80.000	89.601	-12.0	127	0.00
42	C	2,4,6-Trichlorophenol	40.000	45.981	-15.0	140	0.00
43		2,4,5-Trichlorophenol	40.000	45.506	-13.8	137	0.00
44	S	2-Fluorobiphenyl	80.000	82.051	-2.6	134	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	40.753	-1.9	132	0.00
46		2-Chloronaphthalene	40.000	40.797	-2.0	131	0.00
47		2-Nitroaniline	40.000	40.917	-2.3	118	0.00
48		Acenaphthylene	40.000	40.339	-0.8	125	0.00
49		Dimethylphthalate	40.000	40.238	-0.6	122	0.00
50		2,6-Dinitrotoluene	40.000	41.562	-3.9	124	0.00
51	C	Acenaphthene	40.000	40.222	-0.6	127	0.00
52		3-Nitroaniline	40.000	41.385	-3.5	117	0.00
53	P	2,4-Dinitrophenol	40.000	39.694	0.8	123	0.00
54		Dibenzofuran	40.000	39.923	0.2	123	0.00
55	P	4-Nitrophenol	40.000	36.175	9.6	101	0.02
56		2,4-Dinitrotoluene	40.000	41.264	-3.2	114	0.00
57		Fluorene	40.000	39.871	0.3	121	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	44.441	-11.1	127	0.00
59		Diethylphthalate	40.000	38.550	3.6	114	0.00
60		4-Chlorophenyl-phenylether	40.000	40.709	-1.8	128	0.00
61		4-Nitroaniline	40.000	37.995	5.0	103	0.01
62		Azobenzene	40.000	38.645	3.4	115	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	113	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.420	-1.1	111	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.225	-5.6	118	0.00
66		4-Bromophenyl-phenylether	40.000	43.437	-8.6	123	0.00
67		Hexachlorobenzene	40.000	42.661	-6.7	120	0.00
68		Atrazine	40.000	41.266	-3.2	110	0.00
69	C	Pentachlorophenol	40.000	43.248	-8.1	122	0.00
70		Phenanthrene	40.000	40.279	-0.7	111	0.00
71		Anthracene	40.000	40.661	-1.7	111	0.00
72		Carbazole	40.000	39.221	1.9	105	0.00
73		Di-n-butylphthalate	40.000	39.059	2.4	104	0.00
74	C	Fluoranthene	40.000	38.460	3.8	105	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	36.380	9.0	94	0.00
77		Pyrene	40.000	42.167	-5.4	104	0.00
78	S	Terphenyl-d14	80.000	83.080	-3.8	106	0.00
79		Butylbenzylphthalate	40.000	42.049	-5.1	104	0.00
80		Benzo(a)anthracene	40.000	40.670	-1.7	105	0.00
81		3,3'-Dichlorobenzidine	40.000	42.148	-5.4	108	0.00
82		Chrysene	40.000	40.047	-0.1	103	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.922	-4.8	103	0.00
84	c	Di-n-octyl phthalate	40.000	42.107	-5.3	103	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.385	-3.5	106	0.01
86	I	Perylene-d12	20.000	20.000	0.0	106	0.00
87		Benzo(b)fluoranthene	40.000	41.467	-3.7	108	0.00

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88	Benzo(k)fluoranthene	40.000	40.045	-0.1	101	0.00
89 C	Benzo(a)pyrene	40.000	41.037	-2.6	105	0.00
90	Dibenzo(a,h)anthracene	40.000	41.209	-3.0	105	0.00
91	Benzo(g,h,i)perylene	40.000	41.106	-2.8	104	0.00

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