Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120215\
Data File : BG019842.D
Acq On : 2 Dec 2015 19:59
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 03 00:10:59 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	109	0.00
2		1,4-Dioxane	40.000	40.072	-0.2	110	0.00
3		Pyridine	40.000	39.524	1.2	107	0.00
4		n-Nitrosodimethylamine	40.000	41.328	-3.3	107	0.00
5	S	2-Fluorophenol	80.000	78.706	1.6	108	0.00
6		Aniline	40.000	39.304	1.7	105	0.00
7	S	Phenol-d6	80.000	78.959	1.3	108	0.00
8		2-Chlorophenol	40.000	39.104	2.2	106	0.00
9		Benzaldehyde	40.000	38.491	3.8	106	0.00
10	C	Phenol	40.000	38.867	2.8	104	0.00
11		bis(2-Chloroethyl)ether	40.000	39.693	0.8	109	0.00
12		1,3-Dichlorobenzene	40.000	38.253	4.4	105	0.00
13	C	1,4-Dichlorobenzene	40.000	38.043	4.9	105	0.00
14		1,2-Dichlorobenzene	40.000	38.952	2.6	107	0.00
15		Benzyl Alcohol	40.000	39.019	2.5	108	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.899	-2.2	112	0.00
17		2-Methylphenol	40.000	39.014	2.5	105	0.00
18		Hexachloroethane	40.000	40.332	-0.8	108	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.159	2.1	107	0.00
20		3+4-Methylphenols	40.000	39.859	0.4	105	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	108	0.00
22		Acetophenone	40.000	39.795	0.5	106	0.00
23	S	Nitrobenzene-d5	80.000	83.317	-4.1	111	0.00
24		Nitrobenzene	40.000	41.782	-4.5	110	0.00
25		Isophorone	40.000	40.287	-0.7	108	0.00
26	C	2-Nitrophenol	40.000	41.545	-3.9	109	0.00
27		2,4-Dimethylphenol	40.000	40.147	-0.4	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.140	-2.9	111	0.00
29	C	2,4-Dichlorophenol	40.000	39.515	1.2	104	0.00
30		1,2,4-Trichlorobenzene	40.000	38.543	3.6	103	0.00
31		Naphthalene	40.000	40.068	-0.2	107	0.00
32		Benzoic acid	40.000	41.717	-4.3	121	0.00
33		4-Chloroaniline	40.000	39.965	0.1	108	0.00
34	C	Hexachlorobutadiene	40.000	38.515	3.7	104	0.00
35		Caprolactam	40.000	40.003	-0.0	111	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.582	1.0	108	0.00
37		2-Methylnaphthalene	40.000	39.609	1.0	107	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.228	4.4	102	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.784	-4.5	108	0.00
41	S	2,4,6-Tribromophenol	80.000	78.086	2.4	104	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.549	3.6	104	0.00
43		2,4,5-Trichlorophenol	40.000	39.025	2.4	106	0.00
44	S	2-Fluorobiphenyl	80.000	79.308	0.9	106	0.00

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 2015
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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	39.608	1.0	105	0.00
46		2-Chloronaphthalene	40.000	39.564	1.1	105	0.00
47		2-Nitroaniline	40.000	43.572	-8.9	113	0.00
48		Acenaphthylene	40.000	39.730	0.7	106	0.00
49		Dimethylphthalate	40.000	39.497	1.3	107	0.00
50		2,6-Dinitrotoluene	40.000	42.956	-7.4	109	0.00
51	C	Acenaphthene	40.000	40.910	-2.3	108	0.00
52		3-Nitroaniline	40.000	43.646	-9.1	113	0.00
53	P	2,4-Dinitrophenol	40.000	39.252	1.9	115	0.00
54		Dibenzofuran	40.000	39.581	1.0	106	0.00
55	P	4-Nitrophenol	40.000	42.690	-6.7	110	0.00
56		2,4-Dinitrotoluene	40.000	44.183	-10.5	111	0.00
57		Fluorene	40.000	39.708	0.7	107	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.395	1.5	102	0.00
59		Diethylphthalate	40.000	39.654	0.9	108	0.00
60		4-Chlorophenyl-phenylether	40.000	39.283	1.8	105	0.00
61		4-Nitroaniline	40.000	42.885	-7.2	111	0.00
62		Azobenzene	40.000	40.938	-2.3	110	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	106	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.995	-2.5	114	0.00
65	c	n-Nitrosodiphenylamine	40.000	39.718	0.7	106	0.00
66		4-Bromophenyl-phenylether	40.000	38.646	3.4	102	0.00
67		Hexachlorobenzene	40.000	38.651	3.4	103	0.00
68		Atrazine	40.000	40.197	-0.5	105	0.00
69	C	Pentachlorophenol	40.000	41.079	-2.7	108	0.00
70		Phenanthrene	40.000	39.595	1.0	105	0.00
71		Anthracene	40.000	39.694	0.8	105	0.00
72		Carbazole	40.000	40.602	-1.5	107	0.00
73		Di-n-butylphthalate	40.000	41.364	-3.4	109	0.00
74	C	Fluoranthene	40.000	40.183	-0.5	106	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	108	0.00
76		Benzidine	40.000	41.944	-4.9	109	0.00
77		Pyrene	40.000	40.492	-1.2	107	0.00
78	S	Terphenyl-d14	80.000	79.477	0.7	105	0.00
79		Butylbenzylphthalate	40.000	41.551	-3.9	110	0.00
80		Benzo(a)anthracene	40.000	39.883	0.3	108	0.00
81		3,3'-Dichlorobenzidine	40.000	39.603	1.0	105	0.00
82		Chrysene	40.000	39.795	0.5	107	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.105	-2.8	111	0.00
84	c	Di-n-octyl phthalate	40.000	42.403	-6.0	112	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.320	-3.3	111	0.00
86	I	Perylene-d12	20.000	20.000	0.0	110	0.00
87		Benzo(b)fluoranthene	40.000	39.909	0.2	111	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.159	2.1	110	0.00
89 C	Benzo(a)pyrene	40.000	40.270	-0.7	112	0.00
90	Dibenzo(a,h)anthracene	40.000	39.806	0.5	112	0.00
91	Benzo(g,h,i)perylene	40.000	40.161	-0.4	112	0.00

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