Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG052015\
Data File : BG017228.D
Acq On : 21 May 2015 00:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 21 03:39:45 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 21 02:15:39 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	122	-0.01
2		1,4-Dioxane	40.000	38.798	3.0	119	-0.02
3		Pyridine	40.000	39.164	2.1	114	-0.01
4		n-Nitrosodimethylamine	40.000	36.947	7.6	108	-0.02
5	S	2-Fluorophenol	80.000	77.900	2.6	113	-0.02
6		Aniline	40.000	38.184	4.5	110	-0.01
7	S	Phenol-d6	80.000	77.711	2.9	112	-0.01
8		2-Chlorophenol	40.000	37.747	5.6	110	-0.01
9		Benzaldehyde	40.000	37.027	7.4	111	-0.02
10	C	Phenol	40.000	39.544	1.1	116	-0.01
11		bis(2-Chloroethyl)ether	40.000	38.861	2.8	111	-0.01
12		1,3-Dichlorobenzene	40.000	36.855	7.9	111	-0.01
13	C	1,4-Dichlorobenzene	40.000	36.735	8.2	109	-0.01
14		1,2-Dichlorobenzene	40.000	36.531	8.7	108	-0.01
15		Benzyl Alcohol	40.000	36.895	7.8	108	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	39.572	1.1	118	-0.01
17		2-Methylphenol	40.000	37.637	5.9	113	-0.02
18		Hexachloroethane	40.000	38.130	4.7	113	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	36.258	9.4	106	-0.01
20		3+4-Methylphenols	40.000	36.620	8.5	106	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	113	-0.01
22		Acetophenone	40.000	37.527	6.2	105	-0.02
23	S	Nitrobenzene-d5	80.000	77.206	3.5	106	-0.01
24		Nitrobenzene	40.000	38.789	3.0	109	-0.01
25		Isophorone	40.000	37.547	6.1	103	-0.02
26	C	2-Nitrophenol	40.000	37.202	7.0	103	-0.02
27		2,4-Dimethylphenol	40.000	38.000	5.0	108	-0.02
28		bis(2-Chloroethoxy)methane	40.000	38.304	4.2	106	-0.01
29	C	2,4-Dichlorophenol	40.000	39.264	1.8	107	-0.01
30		1,2,4-Trichlorobenzene	40.000	36.192	9.5	100	-0.02
31		Naphthalene	40.000	38.086	4.8	105	-0.02
32		Benzoic acid	40.000	35.313	11.7	97	0.00
33		4-Chloroaniline	40.000	37.600	6.0	102	-0.02
34	C	Hexachlorobutadiene	40.000	37.057	7.4	102	-0.02
35		Caprolactam	40.000	39.304	1.7	114	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.748	0.6	109	-0.01
37		2-Methylnaphthalene	40.000	35.896	10.3	103	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	106	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	39.748	0.6	102	-0.01
40	P	Hexachlorocyclopentadiene	40.000	34.450	13.9	89	-0.01
41	S	2,4,6-Tribromophenol	80.000	80.562	-0.7	104	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.142	-0.4	101	-0.01
43		2,4,5-Trichlorophenol	40.000	42.144	-5.4	109	-0.01
44	S	2-Fluorobiphenyl	80.000	76.808	4.0	99	-0.01

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
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	38.135	4.7	99	-0.01
46		2-Chloronaphthalene	40.000	38.447	3.9	100	-0.01
47		2-Nitroaniline	40.000	42.114	-5.3	110	-0.01
48		Acenaphthylene	40.000	39.637	0.9	101	-0.01
49		Dimethylphthalate	40.000	37.856	5.4	98	-0.01
50		2,6-Dinitrotoluene	40.000	40.346	-0.9	102	-0.01
51	C	Acenaphthene	40.000	39.317	1.7	103	-0.01
52		3-Nitroaniline	40.000	42.026	-5.1	108	-0.01
53	P	2,4-Dinitrophenol	40.000	31.548	21.1	77	0.00
54		Dibenzofuran	40.000	38.424	3.9	100	-0.01
55	P	4-Nitrophenol	40.000	39.467	1.3	104	-0.01
56		2,4-Dinitrotoluene	40.000	42.426	-6.1	109	0.00
57		Fluorene	40.000	38.803	3.0	99	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	42.106	-5.3	107	-0.01
59		Diethylphthalate	40.000	38.180	4.6	101	-0.01
60		4-Chlorophenyl-phenylether	40.000	38.231	4.4	98	-0.01
61		4-Nitroaniline	40.000	44.224	-10.6	113	0.00
62		Azobenzene	40.000	40.004	-0.0	104	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	114	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	37.350	6.6	100	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.281	6.8	103	0.00
66		4-Bromophenyl-phenylether	40.000	36.407	9.0	102	0.00
67		Hexachlorobenzene	40.000	37.553	6.1	104	0.00
68		Atrazine	40.000	37.588	6.0	102	-0.01
69	C	Pentachlorophenol	40.000	30.422	23.9#	82	0.00
70		Phenanthrene	40.000	39.053	2.4	107	-0.01
71		Anthracene	40.000	38.836	2.9	108	0.00
72		Carbazole	40.000	40.610	-1.5	112	0.00
73		Di-n-butylphthalate	40.000	40.722	-1.8	113	0.00
74	C	Fluoranthene	40.000	40.648	-1.6	113	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	137	0.00
76		Benzidine	40.000	29.694	25.8#	97	-0.01
77		Pyrene	40.000	34.780	13.0	115	0.00
78	S	Terphenyl-d14	80.000	71.144	11.1	118	-0.01
79		Butylbenzylphthalate	40.000	37.722	5.7	125	0.00
80		Benzo(a)anthracene	40.000	37.511	6.2	124	-0.01
81		3,3'-Dichlorobenzidine	40.000	36.706	8.2	122	0.00
82		Chrysene	40.000	37.366	6.6	124	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.693	5.8	123	-0.01
84	c	Di-n-octyl phthalate	40.000	38.640	3.4	128	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.771	3.1	128	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	135	-0.01
87		Benzo(b)fluoranthene	40.000	36.289	9.3	123	0.00

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88	Benzo(k)fluoranthene	40.000	38.923	2.7	127	0.00
89 C	Benzo(a)pyrene	40.000	38.122	4.7	127	0.00
90	Dibenzo(a,h)anthracene	40.000	38.174	4.6	128	-0.02
91	Benzo(g,h,i)perylene	40.000	38.097	4.8	127	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 1