Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060415\
Data File : BG017456.D
Acq On : 4 Jun 2015 23:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 05 04:00:56 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 03:51:43 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	91	0.00
2		1,4-Dioxane	40.000	39.567	1.1	94	0.00
3		Pyridine	40.000	41.105	-2.8	89	0.00
4		n-Nitrosodimethylamine	40.000	45.005	-12.5	97	0.00
5	S	2-Fluorophenol	80.000	84.386	-5.5	95	0.00
6		Aniline	40.000	41.063	-2.7	92	0.00
7	S	Phenol-d6	80.000	79.645	0.4	88	-0.01
8		2-Chlorophenol	40.000	39.792	0.5	90	-0.01
9		Benzaldehyde	40.000	38.610	3.5	83	0.00
10	C	Phenol	40.000	42.803	-7.0	93	0.00
11		bis(2-Chloroethyl)ether	40.000	41.800	-4.5	93	0.00
12		1,3-Dichlorobenzene	40.000	39.843	0.4	91	0.00
13	C	1,4-Dichlorobenzene	40.000	39.027	2.4	89	0.00
14		1,2-Dichlorobenzene	40.000	39.397	1.5	89	0.00
15		Benzyl Alcohol	40.000	40.735	-1.8	88	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.839	-4.6	93	0.00
17		2-Methylphenol	40.000	40.605	-1.5	90	0.00
18		Hexachloroethane	40.000	40.241	-0.6	89	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.166	-0.4	90	-0.01
20		3+4-Methylphenols	40.000	39.640	0.9	87	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	-0.01
22		Acetophenone	40.000	39.493	1.3	89	0.00
23	S	Nitrobenzene-d5	80.000	82.200	-2.8	91	0.00
24		Nitrobenzene	40.000	40.667	-1.7	91	0.00
25		Isophorone	40.000	36.980	7.6	84	0.00
26	C	2-Nitrophenol	40.000	39.228	1.9	86	0.00
27		2,4-Dimethylphenol	40.000	39.478	1.3	89	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.799	3.0	85	0.00
29	C	2,4-Dichlorophenol	40.000	40.876	-2.2	90	0.00
30		1,2,4-Trichlorobenzene	40.000	38.639	3.4	88	0.00
31		Naphthalene	40.000	39.016	2.5	88	0.00
32		Benzoic acid	40.000	34.724	13.2	76	0.00
33		4-Chloroaniline	40.000	37.947	5.1	85	-0.01
34	C	Hexachlorobutadiene	40.000	40.159	-0.4	91	0.00
35		Caprolactam	40.000	34.982	12.5	74	-0.01
36	C	4-Chloro-3-methylphenol	40.000	38.851	2.9	85	0.00
37		2-Methylnaphthalene	40.000	39.502	1.2	88	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.645	3.4	88	0.00
40	P	Hexachlorocyclopentadiene	40.000	36.673	8.3	81	0.00
41	S	2,4,6-Tribromophenol	80.000	81.035	-1.3	88	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.774	3.1	81	-0.01
43		2,4,5-Trichlorophenol	40.000	39.573	1.1	88	-0.01
44	S	2-Fluorobiphenyl	80.000	80.016	-0.0	88	0.00

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Quant Time: Jun 05 04:00:56 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 03:51:43 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)
45		1,1'-Biphenyl	40.000	39.893	0.3	87	0.00
46		2-Chloronaphthalene	40.000	39.852	0.4	84	0.00
47		2-Nitroaniline	40.000	40.540	-1.3	86	0.00
48		Acenaphthylene	40.000	40.505	-1.3	88	0.00
49		Dimethylphthalate	40.000	37.351	6.6	81	0.00
50		2,6-Dinitrotoluene	40.000	38.944	2.6	85	0.00
51	C	Acenaphthene	40.000	39.862	0.3	88	0.00
52		3-Nitroaniline	40.000	38.657	3.4	84	0.00
53	P	2,4-Dinitrophenol	40.000	36.410	9.0	81	0.00
54		Dibenzofuran	40.000	40.502	-1.3	87	0.00
55	P	4-Nitrophenol	40.000	37.948	5.1	81	0.00
56		2,4-Dinitrotoluene	40.000	40.679	-1.7	85	0.00
57		Fluorene	40.000	39.784	0.5	86	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.979	0.1	83	0.00
59		Diethylphthalate	40.000	37.732	5.7	82	-0.01
60		4-Chlorophenyl-phenylether	40.000	39.514	1.2	86	0.00
61		4-Nitroaniline	40.000	42.174	-5.4	86	0.00
62		Azobenzene	40.000	42.095	-5.2	91	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	84	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.638	0.9	82	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.031	-2.6	84	0.00
66		4-Bromophenyl-phenylether	40.000	41.227	-3.1	88	0.00
67		Hexachlorobenzene	40.000	41.013	-2.5	86	0.00
68		Atrazine	40.000	40.274	-0.7	82	0.00
69	C	Pentachlorophenol	40.000	39.489	1.3	83	0.00
70		Phenanthrene	40.000	41.704	-4.3	88	0.00
71		Anthracene	40.000	41.748	-4.4	88	0.00
72		Carbazole	40.000	40.981	-2.5	86	0.00
73		Di-n-butylphthalate	40.000	40.252	-0.6	84	0.00
74	C	Fluoranthene	40.000	41.006	-2.5	87	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	85	0.00
76		Benzidine	40.000	41.539	-3.8	82	0.00
77		Pyrene	40.000	41.426	-3.6	86	0.00
78	S	Terphenyl-d14	80.000	85.108	-6.4	89	0.00
79		Butylbenzylphthalate	40.000	41.428	-3.6	85	0.00
80		Benzo(a)anthracene	40.000	41.594	-4.0	87	0.00
81		3,3'-Dichlorobenzidine	40.000	41.822	-4.6	86	0.00
82		Chrysene	40.000	41.612	-4.0	87	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.382	-6.0	88	0.00
84	c	Di-n-octyl phthalate	40.000	42.408	-6.0	90	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	42.131	-5.3	88	0.00
86	I	Perylene-d12	20.000	20.000	0.0	88	0.00
87		Benzo(b)fluoranthene	40.000	40.017	-0.0	85	0.00

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Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	40.832	-2.1	92	0.00
89 C	Benzo(a)pyrene	40.000	40.379	-0.9	88	0.00
90	Dibenzo(a,h)anthracene	40.000	40.423	-1.1	87	0.00
91	Benzo(g,h,i)perylene	40.000	40.363	-0.9	89	0.00

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