Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG061115\
Data File : BG017628.D
Acq On : 12 Jun 2015 2:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 12 03:31:54 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 01:16:03 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	113	-0.08
2		1,4-Dioxane	40.000	38.561	3.6	114	-0.09
3		Pyridine	40.000	37.243	6.9	101	-0.09
4		n-Nitrosodimethylamine	40.000	55.882	-39.7#	151	-0.09
5	S	2-Fluorophenol	80.000	76.043	4.9	106	-0.09
6		Aniline	40.000	39.693	0.8	111	-0.08
7	S	Phenol-d6	80.000	76.307	4.6	105	-0.08
8		2-Chlorophenol	40.000	39.292	1.8	111	-0.08
9		Benzaldehyde	40.000	35.027	12.4	94	-0.08
10	C	Phenol	40.000	40.273	-0.7	109	-0.08
11		bis(2-Chloroethyl)ether	40.000	39.465	1.3	109	-0.08
12		1,3-Dichlorobenzene	40.000	39.602	1.0	113	-0.08
13	C	1,4-Dichlorobenzene	40.000	39.685	0.8	112	-0.08
14		1,2-Dichlorobenzene	40.000	39.805	0.5	112	-0.08
15		Benzyl Alcohol	40.000	45.468	-13.7	122	-0.08
16		2,2'-oxybis(1-Chloropropane	40.000	38.108	4.7	106	-0.08
17		2-Methylphenol	40.000	41.721	-4.3	115	-0.08
18		Hexachloroethane	40.000	43.788	-9.5	121	-0.08
19	P	n-Nitroso-di-n-propylamine	40.000	43.890	-9.7	122	-0.08
20		3+4-Methylphenols	40.000	41.011	-2.5	112	-0.08
21	I	Naphthalene-d8	20.000	20.000	0.0	113	-0.08
22		Acetophenone	40.000	42.806	-7.0	118	-0.08
23	S	Nitrobenzene-d5	80.000	91.758	-14.7	125	-0.08
24		Nitrobenzene	40.000	45.435	-13.6	126	-0.08
25		Isophorone	40.000	42.812	-7.0	121	-0.08
26	C	2-Nitrophenol	40.000	43.540	-8.8	118	-0.08
27		2,4-Dimethylphenol	40.000	41.571	-3.9	115	-0.08
28		bis(2-Chloroethoxy)methane	40.000	41.409	-3.5	112	-0.08
29	C	2,4-Dichlorophenol	40.000	45.865	-14.7	125	-0.09
30		1,2,4-Trichlorobenzene	40.000	46.640	-16.6	131	-0.08
31		Naphthalene	40.000	39.995	0.0	111	-0.08
32		Benzoic acid	40.000	43.119	-7.8	116	-0.08
33		4-Chloroaniline	40.000	40.641	-1.6	112	-0.08
34	C	Hexachlorobutadiene	40.000	55.433	-38.6#	154	-0.08
35		Caprolactam	40.000	36.607	8.5	96	-0.09
36	C	4-Chloro-3-methylphenol	40.000	42.413	-6.0	114	-0.09
37		2-Methylnaphthalene	40.000	42.657	-6.6	117	-0.08
38	I	Acenaphthene-d10	20.000	20.000	0.0	128	-0.08
39		1,2,4,5-Tetrachlorobenzene	40.000	48.646	-21.6	158	-0.08
40	P	Hexachlorocyclopentadiene	40.000	29.925	25.2#	90	-0.08
41	S	2,4,6-Tribromophenol	80.000	82.699	-3.4	128	-0.08
42	C	2,4,6-Trichlorophenol	40.000	43.728	-9.3	131	-0.08
43		2,4,5-Trichlorophenol	40.000	43.886	-9.7	140	-0.09
44	S	2-Fluorobiphenyl	80.000	87.162	-9.0	138	-0.08

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Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 12 03:31:54 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 01:16:03 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	40.267	-0.7	126	-0.08
46		2-Chloronaphthalene	40.000	41.602	-4.0	127	-0.08
47		2-Nitroaniline	40.000	41.644	-4.1	127	-0.08
48		Acenaphthylene	40.000	40.409	-1.0	125	-0.08
49		Dimethylphthalate	40.000	39.228	1.9	121	-0.08
50		2,6-Dinitrotoluene	40.000	40.137	-0.3	126	-0.08
51	C	Acenaphthene	40.000	39.687	0.8	126	-0.08
52		3-Nitroaniline	40.000	35.806	10.5	111	-0.08
53	P	2,4-Dinitrophenol	40.000	36.789	8.0	117	-0.07
54		Dibenzofuran	40.000	43.172	-7.9	133	-0.08
55	P	4-Nitrophenol	40.000	30.318	24.2	93	-0.09
56		2,4-Dinitrotoluene	40.000	41.649	-4.1	124	-0.08
57		Fluorene	40.000	43.114	-7.8	134	-0.08
58		2,3,4,6-Tetrachlorophenol	40.000	50.111	-25.3#	149	-0.08
59		Diethylphthalate	40.000	38.969	2.6	121	-0.08
60		4-Chlorophenyl-phenylether	40.000	47.786	-19.5	150	-0.08
61		4-Nitroaniline	40.000	33.789	15.5	99	-0.08
62		Azobenzene	40.000	44.891	-12.2	139	-0.08
63	I	Phenanthrene-d10	20.000	20.000	0.0	132	-0.08
64		4,6-Dinitro-2-methylphenol	40.000	40.477	-1.2	132	-0.07
65	c	n-Nitrosodiphenylamine	40.000	40.268	-0.7	130	-0.08
66		4-Bromophenyl-phenylether	40.000	43.934	-9.8	147	-0.08
67		Hexachlorobenzene	40.000	42.703	-6.8	141	-0.08
68		Atrazine	40.000	41.174	-2.9	131	-0.08
69	C	Pentachlorophenol	40.000	34.193	14.5	111	-0.08
70		Phenanthrene	40.000	39.244	1.9	130	-0.08
71		Anthracene	40.000	38.951	2.6	130	-0.08
72		Carbazole	40.000	36.735	8.2	122	-0.08
73		Di-n-butylphthalate	40.000	36.055	9.9	118	-0.08
74	C	Fluoranthene	40.000	41.146	-2.9	137	-0.08
75	I	Chrysene-d12	20.000	20.000	0.0	133	-0.08
76		Benzidine	40.000	31.985	20.0	99	-0.08
77		Pyrene	40.000	40.943	-2.4	134	-0.08
78	S	Terphenyl-d14	80.000	87.139	-8.9	143	-0.08
79		Butylbenzylphthalate	40.000	37.507	6.2	121	-0.08
80		Benzo(a)anthracene	40.000	43.429	-8.6	142	-0.08
81		3,3'-Dichlorobenzidine	40.000	40.490	-1.2	130	-0.08
82		Chrysene	40.000	44.002	-10.0	144	-0.08
83		Bis(2-ethylhexyl)phthalate	40.000	37.750	5.6	123	-0.08
84	c	Di-n-octyl phthalate	40.000	35.759	10.6	118	-0.09
85		Indeno(1,2,3-cd)pyrene	40.000	39.272	1.8	128	-0.18
86	I	Perylene-d12	20.000	20.000	0.0	124	-0.12
87		Benzo(b)fluoranthene	40.000	42.774	-6.9	129	-0.10

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Instrument :
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LabSampleId :
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Quant Time: Jun 12 03:31:54 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 01:16:03 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)
88	Benzo(k)fluoranthene	40.000	43.215	-8.0	137	-0.11
89 C	Benzo(a)pyrene	40.000	42.506	-6.3	130	-0.12
90	Dibenzo(a,h)anthracene	40.000	41.206	-3.0	126	-0.18
91	Benzo(g,h,i)perylene	40.000	41.478	-3.7	129	-0.20

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