Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051315\
Data File : BG017135.D
Acq On : 14 May 2015 7:54
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 14 09:11:15 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 14 06:13:40 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	41.506	-3.8	139	0.00
3		Pyridine	40.000	43.711	-9.3	139	0.00
4		n-Nitrosodimethylamine	40.000	37.706	5.7	119	0.00
5	S	2-Fluorophenol	80.000	81.014	-1.3	128	0.00
6		Aniline	40.000	41.828	-4.6	131	0.00
7	S	Phenol-d6	80.000	84.684	-5.9	133	0.00
8		2-Chlorophenol	40.000	40.668	-1.7	129	0.00
9		Benzaldehyde	40.000	42.438	-6.1	132	0.00
10	C	Phenol	40.000	41.805	-4.5	133	0.00
11		bis(2-Chloroethyl)ether	40.000	44.155	-10.4	137	0.00
12		1,3-Dichlorobenzene	40.000	38.109	4.7	125	0.00
13	C	1,4-Dichlorobenzene	40.000	38.831	2.9	125	0.00
14		1,2-Dichlorobenzene	40.000	38.736	3.2	124	0.00
15		Benzyl Alcohol	40.000	40.643	-1.6	129	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.790	-12.0	145	0.00
17		2-Methylphenol	40.000	41.172	-2.9	134	0.00
18		Hexachloroethane	40.000	38.397	4.0	124	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.035	-7.6	136	0.00
20		3+4-Methylphenols	40.000	41.478	-3.7	130	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	129	0.00
22		Acetophenone	40.000	40.910	-2.3	130	0.00
23	S	Nitrobenzene-d5	80.000	82.291	-2.9	128	0.00
24		Nitrobenzene	40.000	41.462	-3.7	133	0.00
25		Isophorone	40.000	41.479	-3.7	129	0.00
26	C	2-Nitrophenol	40.000	40.315	-0.8	127	0.00
27		2,4-Dimethylphenol	40.000	40.887	-2.2	133	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.044	-5.1	133	0.00
29	C	2,4-Dichlorophenol	40.000	40.619	-1.5	126	0.00
30		1,2,4-Trichlorobenzene	40.000	38.130	4.7	120	0.00
31		Naphthalene	40.000	40.090	-0.2	126	0.00
32		Benzoic acid	40.000	39.306	1.7	123	0.00
33		4-Chloroaniline	40.000	41.479	-3.7	128	0.00
34	C	Hexachlorobutadiene	40.000	36.535	8.7	115	0.00
35		Caprolactam	40.000	42.674	-6.7	141	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.149	-5.4	131	0.00
37		2-Methylnaphthalene	40.000	39.304	1.7	128	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	125	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.775	0.6	121	0.00
40	P	Hexachlorocyclopentadiene	40.000	38.391	4.0	117	0.00
41	S	2,4,6-Tribromophenol	80.000	79.299	0.9	121	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.857	-4.6	124	0.00
43		2,4,5-Trichlorophenol	40.000	41.695	-4.2	126	0.00
44	S	2-Fluorobiphenyl	80.000	79.866	0.2	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051315\
Data File : BG017135.D
Acq On : 14 May 2015 7:54
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 14 09:11:15 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 14 06:13:40 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	40.189	-0.5	122	0.00
46		2-Chloronaphthalene	40.000	40.993	-2.5	125	0.00
47		2-Nitroaniline	40.000	43.895	-9.7	135	0.00
48		Acenaphthylene	40.000	41.427	-3.6	125	0.00
49		Dimethylphthalate	40.000	40.221	-0.6	123	0.00
50		2,6-Dinitrotoluene	40.000	41.633	-4.1	123	0.00
51	C	Acenaphthene	40.000	40.872	-2.2	126	0.00
52		3-Nitroaniline	40.000	42.986	-7.5	130	0.00
53	P	2,4-Dinitrophenol	40.000	39.430	1.4	113	0.00
54		Dibenzofuran	40.000	40.414	-1.0	124	0.00
55	P	4-Nitrophenol	40.000	43.170	-7.9	133	0.00
56		2,4-Dinitrotoluene	40.000	42.545	-6.4	129	0.00
57		Fluorene	40.000	40.863	-2.2	123	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.055	-2.6	122	0.00
59		Diethylphthalate	40.000	39.549	1.1	123	0.00
60		4-Chlorophenyl-phenylether	40.000	40.145	-0.4	121	0.00
61		4-Nitroaniline	40.000	44.009	-10.0	132	0.00
62		Azobenzene	40.000	42.409	-6.0	130	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	126	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.496	1.3	117	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.806	-2.0	125	0.00
66		4-Bromophenyl-phenylether	40.000	38.054	4.9	118	0.00
67		Hexachlorobenzene	40.000	39.630	0.9	122	0.00
68		Atrazine	40.000	39.698	0.8	119	0.00
69	C	Pentachlorophenol	40.000	39.599	1.0	119	0.00
70		Phenanthrene	40.000	41.242	-3.1	125	0.00
71		Anthracene	40.000	40.554	-1.4	125	0.00
72		Carbazole	40.000	41.741	-4.4	128	0.00
73		Di-n-butylphthalate	40.000	41.445	-3.6	128	0.00
74	C	Fluoranthene	40.000	39.757	0.6	122	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	124	0.00
76		Benzidine	40.000	39.418	1.5	116	0.00
77		Pyrene	40.000	41.500	-3.8	125	0.00
78	S	Terphenyl-d14	80.000	80.786	-1.0	122	0.00
79		Butylbenzylphthalate	40.000	42.070	-5.2	127	0.00
80		Benzo(a)anthracene	40.000	40.868	-2.2	122	0.00
81		3,3'-Dichlorobenzidine	40.000	39.819	0.5	120	0.00
82		Chrysene	40.000	40.426	-1.1	122	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.172	-2.9	122	0.00
84	c	Di-n-octyl phthalate	40.000	41.441	-3.6	125	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.257	-0.6	120	0.00
86	I	Perylene-d12	20.000	20.000	0.0	120	0.00
87		Benzo(b)fluoranthene	40.000	40.616	-1.5	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051315\
Data File : BG017135.D
Acq On : 14 May 2015 7:54
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 14 09:11:15 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 14 06:13:40 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	40.995	-2.5	119	0.00
89 C	Benzo(a)pyrene	40.000	41.474	-3.7	123	0.00
90	Dibenzo(a,h)anthracene	40.000	41.215	-3.0	123	-0.01
91	Benzo(g,h,i)perylene	40.000	40.691	-1.7	120	-0.01

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