Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	107	0.00
2		1,4-Dioxane	40.000	38.036	4.9	108	0.00
3		Pyridine	40.000	37.625	5.9	103	0.00
4		n-Nitrosodimethylamine	40.000	38.989	2.5	112	0.00
5	S	2-Fluorophenol	80.000	77.264	3.4	108	0.00
6		Aniline	40.000	37.802	5.5	106	0.00
7	S	Phenol-d6	80.000	77.920	2.6	109	0.00
8		2-Chlorophenol	40.000	39.123	2.2	111	0.00
9		Benzaldehyde	40.000	38.833	2.9	106	0.00
10	C	Phenol	40.000	39.069	2.3	110	0.00
11		bis(2-Chloroethyl)ether	40.000	38.322	4.2	106	0.00
12		1,3-Dichlorobenzene	40.000	36.854	7.9	105	0.00
13	C	1,4-Dichlorobenzene	40.000	36.756	8.1	103	0.00
14		1,2-Dichlorobenzene	40.000	36.401	9.0	102	0.00
15		Benzyl Alcohol	40.000	38.238	4.4	107	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.544	6.1	107	0.00
17		2-Methylphenol	40.000	37.210	7.0	104	0.00
18		Hexachloroethane	40.000	37.106	7.2	104	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	37.498	6.3	107	0.00
20		3+4-Methylphenols	40.000	38.172	4.6	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	37.153	7.1	106	0.00
23	S	Nitrobenzene-d5	80.000	77.782	2.8	113	0.00
24		Nitrobenzene	40.000	38.741	3.1	111	0.00
25		Isophorone	40.000	36.503	8.7	106	0.00
26	C	2-Nitrophenol	40.000	39.042	2.4	112	0.00
27		2,4-Dimethylphenol	40.000	36.796	8.0	104	0.00
28		bis(2-Chloroethoxy)methane	40.000	36.148	9.6	107	0.00
29	C	2,4-Dichlorophenol	40.000	35.823	10.4	103	0.00
30		1,2,4-Trichlorobenzene	40.000	36.004	10.0	103	0.00
31		Naphthalene	40.000	36.103	9.7	105	0.00
32		Benzoic acid	40.000	40.993	-2.5	117	0.00
33		4-Chloroaniline	40.000	37.062	7.3	105	0.00
34	C	Hexachlorobutadiene	40.000	34.924	12.7	102	0.00
35		Caprolactam	40.000	36.559	8.6	105	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.064	7.3	107	0.00
37		2-Methylnaphthalene	40.000	34.739	13.2	102	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	37.216	7.0	102	0.00
40	P	Hexachlorocyclopentadiene	40.000	34.513	13.7	95	0.00
41	S	2,4,6-Tribromophenol	80.000	77.489	3.1	106	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.380	4.0	100	0.00
43		2,4,5-Trichlorophenol	40.000	38.784	3.0	104	0.00
44	S	2-Fluorobiphenyl	80.000	74.574	6.8	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	37.286	6.8	102	0.00
46		2-Chloronaphthalene	40.000	37.878	5.3	103	0.00
47		2-Nitroaniline	40.000	43.831	-9.6	117	0.00
48		Acenaphthylene	40.000	38.034	4.9	103	-0.01
49		Dimethylphthalate	40.000	37.228	6.9	102	0.00
50		2,6-Dinitrotoluene	40.000	40.111	-0.3	105	0.00
51	C	Acenaphthene	40.000	37.159	7.1	102	0.00
52		3-Nitroaniline	40.000	42.072	-5.2	109	0.00
53	P	2,4-Dinitrophenol	40.000	37.338	6.7	107	0.00
54		Dibenzofuran	40.000	37.836	5.4	102	0.00
55	P	4-Nitrophenol	40.000	39.527	1.2	106	0.00
56		2,4-Dinitrotoluene	40.000	43.560	-8.9	115	0.00
57		Fluorene	40.000	37.064	7.3	101	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.648	3.4	101	0.00
59		Diethylphthalate	40.000	37.496	6.3	103	0.00
60		4-Chlorophenyl-phenylether	40.000	36.353	9.1	101	0.00
61		4-Nitroaniline	40.000	39.747	0.6	107	0.00
62		Azobenzene	40.000	38.712	3.2	106	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.981	-2.5	113	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.109	7.2	102	0.00
66		4-Bromophenyl-phenylether	40.000	36.628	8.4	101	0.00
67		Hexachlorobenzene	40.000	37.881	5.3	105	0.00
68		Atrazine	40.000	36.258	9.4	97	0.00
69	C	Pentachlorophenol	40.000	37.739	5.7	100	0.00
70		Phenanthrene	40.000	36.918	7.7	101	0.00
71		Anthracene	40.000	36.486	8.8	100	0.00
72		Carbazole	40.000	36.921	7.7	99	0.00
73		Di-n-butylphthalate	40.000	37.380	6.5	101	-0.01
74	C	Fluoranthene	40.000	35.946	10.1	96	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	97	0.00
76		Benzidine	40.000	34.007	15.0	83	0.00
77		Pyrene	40.000	37.700	5.7	98	0.00
78	S	Terphenyl-d14	80.000	77.208	3.5	98	0.00
79		Butylbenzylphthalate	40.000	38.948	2.6	100	-0.01
80		Benzo(a)anthracene	40.000	38.118	4.7	98	0.00
81		3,3'-Dichlorobenzidine	40.000	37.239	6.9	95	-0.01
82		Chrysene	40.000	37.893	5.3	97	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.892	2.8	98	-0.01
84	c	Di-n-octyl phthalate	40.000	39.226	1.9	99	-0.01
85		Indeno(1,2,3-cd)pyrene	40.000	38.230	4.4	101	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	99	-0.01
87		Benzo(b)fluoranthene	40.000	37.392	6.5	98	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	37.532	6.2	99	-0.01
89 C	Benzo(a)pyrene	40.000	37.379	6.6	98	-0.01
90	Dibenzo(a,h)anthracene	40.000	37.656	5.9	101	-0.03
91	Benzo(g,h,i)perylene	40.000	37.485	6.3	102	-0.01

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