Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051515\
Data File : BG017189.D
Acq On : 16 May 2015 00:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 16 01:09:08 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 16 00:01:34 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	38.595	3.5	104	0.00
3		Pyridine	40.000	41.122	-2.8	106	0.00
4		n-Nitrosodimethylamine	40.000	38.285	4.3	98	0.00
5	S	2-Fluorophenol	80.000	79.304	0.9	101	0.00
6		Aniline	40.000	41.653	-4.1	106	0.00
7	S	Phenol-d6	80.000	84.395	-5.5	107	0.00
8		2-Chlorophenol	40.000	40.298	-0.7	103	0.00
9		Benzaldehyde	40.000	39.780	0.5	102	0.00
10	C	Phenol	40.000	44.035	-10.1	114	0.00
11		bis(2-Chloroethyl)ether	40.000	41.398	-3.5	104	0.00
12		1,3-Dichlorobenzene	40.000	37.651	5.9	100	0.00
13	C	1,4-Dichlorobenzene	40.000	38.414	4.0	100	0.00
14		1,2-Dichlorobenzene	40.000	38.491	3.8	100	0.00
15		Benzyl Alcohol	40.000	39.743	0.6	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.886	-4.7	110	0.00
17		2-Methylphenol	40.000	40.868	-2.2	108	0.00
18		Hexachloroethane	40.000	39.876	0.3	104	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.559	-3.9	106	0.00
20		3+4-Methylphenols	40.000	42.066	-5.2	107	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	105	0.00
22		Acetophenone	40.000	38.651	3.4	100	0.00
23	S	Nitrobenzene-d5	80.000	82.219	-2.8	104	0.00
24		Nitrobenzene	40.000	40.094	-0.2	105	0.00
25		Isophorone	40.000	40.341	-0.9	102	0.00
26	C	2-Nitrophenol	40.000	39.395	1.5	101	0.00
27		2,4-Dimethylphenol	40.000	38.626	3.4	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.907	0.2	103	0.00
29	C	2,4-Dichlorophenol	40.000	42.005	-5.0	106	0.00
30		1,2,4-Trichlorobenzene	40.000	38.383	4.0	99	0.00
31		Naphthalene	40.000	39.669	0.8	102	0.00
32		Benzoic acid	40.000	36.911	7.7	94	0.00
33		4-Chloroaniline	40.000	41.854	-4.6	105	0.00
34	C	Hexachlorobutadiene	40.000	38.052	4.9	97	-0.01
35		Caprolactam	40.000	42.120	-5.3	113	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.974	-7.4	109	0.00
37		2-Methylnaphthalene	40.000	39.626	0.9	105	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.422	3.9	98	0.00
40	P	Hexachlorocyclopentadiene	40.000	33.140	17.1	85	0.00
41	S	2,4,6-Tribromophenol	80.000	78.872	1.4	101	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.856	0.4	99	0.00
43		2,4,5-Trichlorophenol	40.000	42.057	-5.1	107	0.00
44	S	2-Fluorobiphenyl	80.000	77.900	2.6	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051515\
Data File : BG017189.D
Acq On : 16 May 2015 00:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 16 01:09:08 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 16 00:01:34 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	39.464	1.3	101	0.00
46		2-Chloronaphthalene	40.000	39.872	0.3	103	0.00
47		2-Nitroaniline	40.000	42.142	-5.4	109	0.00
48		Acenaphthylene	40.000	40.434	-1.1	102	0.00
49		Dimethylphthalate	40.000	39.106	2.2	101	0.00
50		2,6-Dinitrotoluene	40.000	42.797	-7.0	107	0.00
51	C	Acenaphthene	40.000	40.625	-1.6	105	0.00
52		3-Nitroaniline	40.000	42.816	-7.0	109	0.00
53	P	2,4-Dinitrophenol	40.000	31.911	20.2	77	0.00
54		Dibenzofuran	40.000	39.883	0.3	103	0.00
55	P	4-Nitrophenol	40.000	36.862	7.8	96	0.00
56		2,4-Dinitrotoluene	40.000	42.651	-6.6	109	0.00
57		Fluorene	40.000	40.456	-1.1	102	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.764	0.6	100	0.00
59		Diethylphthalate	40.000	38.885	2.8	102	0.00
60		4-Chlorophenyl-phenylether	40.000	39.525	1.2	100	0.00
61		4-Nitroaniline	40.000	41.982	-5.0	106	0.00
62		Azobenzene	40.000	41.337	-3.3	106	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
64		4,6-Dinitro-2-methylphenol	40.000	36.854	7.9	90	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.574	-3.9	105	0.00
66		4-Bromophenyl-phenylether	40.000	40.560	-1.4	103	0.00
67		Hexachlorobenzene	40.000	39.849	0.4	101	0.00
68		Atrazine	40.000	39.213	2.0	97	0.00
69	C	Pentachlorophenol	40.000	34.154	14.6	84	0.00
70		Phenanthrene	40.000	41.510	-3.8	104	0.00
71		Anthracene	40.000	41.623	-4.1	105	0.00
72		Carbazole	40.000	40.636	-1.6	103	0.00
73		Di-n-butylphthalate	40.000	41.957	-4.9	107	0.00
74	C	Fluoranthene	40.000	39.462	1.3	100	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	101	0.00
76		Benzidine	40.000	34.544	13.6	83	0.00
77		Pyrene	40.000	41.480	-3.7	102	0.00
78	S	Terphenyl-d14	80.000	83.377	-4.2	103	0.00
79		Butylbenzylphthalate	40.000	41.821	-4.6	103	0.00
80		Benzo(a)anthracene	40.000	41.006	-2.5	100	0.00
81		3,3'-Dichlorobenzidine	40.000	38.698	3.3	96	0.00
82		Chrysene	40.000	40.625	-1.6	100	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.820	-4.6	101	0.00
84	c	Di-n-octyl phthalate	40.000	41.784	-4.5	103	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.947	-2.4	100	0.00
86	I	Perylene-d12	20.000	20.000	0.0	101	0.00
87		Benzo(b)fluoranthene	40.000	39.542	1.1	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG051515\
Data File : BG017189.D
Acq On : 16 May 2015 00:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 16 01:09:08 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 16 00:01:34 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	42.117	-5.3	103	0.00
89 C	Benzo(a)pyrene	40.000	40.951	-2.4	102	0.00
90	Dibenzo(a,h)anthracene	40.000	40.364	-0.9	101	-0.01
91	Benzo(g,h,i)perylene	40.000	39.633	0.9	99	0.00

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