Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016667.D
Acq On : 23 Apr 2015 23:46
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SSTDCCC040EC

Quant Time: Apr 24 02:27:41 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 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	102	0.00
2		1,4-Dioxane	40.000	44.005	-10.0	105	0.00
3		Pyridine	40.000	43.705	-9.3	104	0.00
4		n-Nitrosodimethylamine	40.000	41.571	-3.9	103	0.00
5	S	2-Fluorophenol	80.000	85.169	-6.5	103	0.00
6		Aniline	40.000	42.243	-5.6	100	0.00
7	S	Phenol-d6	80.000	84.715	-5.9	100	0.00
8		2-Chlorophenol	40.000	41.797	-4.5	101	0.00
9		Benzaldehyde	40.000	37.132	7.2	101	0.00
10	C	Phenol	40.000	41.986	-5.0	101	0.00
11		bis(2-Chloroethyl)ether	40.000	41.169	-2.9	101	0.00
12		1,3-Dichlorobenzene	40.000	41.243	-3.1	101	0.00
13	C	1,4-Dichlorobenzene	40.000	41.939	-4.8	103	0.00
14		1,2-Dichlorobenzene	40.000	41.334	-3.3	101	0.00
15		Benzyl Alcohol	40.000	42.820	-7.1	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.906	-2.3	100	0.00
17		2-Methylphenol	40.000	42.767	-6.9	100	0.00
18		Hexachloroethane	40.000	42.622	-6.6	105	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.176	-5.4	98	0.00
20		3+4-Methylphenols	40.000	42.313	-5.8	101	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	43.209	-8.0	100	0.00
23	S	Nitrobenzene-d5	80.000	86.967	-8.7	100	0.00
24		Nitrobenzene	40.000	43.337	-8.3	100	0.00
25		Isophorone	40.000	42.907	-7.3	97	0.00
26	C	2-Nitrophenol	40.000	44.164	-10.4	100	0.00
27		2,4-Dimethylphenol	40.000	42.864	-7.2	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.521	-6.3	99	0.00
29	C	2,4-Dichlorophenol	40.000	43.040	-7.6	97	0.00
30		1,2,4-Trichlorobenzene	40.000	42.043	-5.1	99	0.00
31		Naphthalene	40.000	41.941	-4.9	99	0.00
32		Benzoic acid	40.000	34.032	14.9	79	0.00
33		4-Chloroaniline	40.000	42.075	-5.2	97	0.00
34	C	Hexachlorobutadiene	40.000	42.363	-5.9	101	0.00
35		Caprolactam	40.000	42.549	-6.4	91	-0.02
36	C	4-Chloro-3-methylphenol	40.000	42.819	-7.0	96	0.00
37		2-Methylnaphthalene	40.000	41.517	-3.8	97	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	95	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	42.280	-5.7	95	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.609	-1.5	97	0.00
41	S	2,4,6-Tribromophenol	80.000	83.824	-4.8	91	0.00
42	C	2,4,6-Trichlorophenol	40.000	42.754	-6.9	91	0.00
43		2,4,5-Trichlorophenol	40.000	42.785	-7.0	93	0.00
44	S	2-Fluorobiphenyl	80.000	85.332	-6.7	96	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016667.D
Acq On : 23 Apr 2015 23:46
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SSTDCCC040EC

Quant Time: Apr 24 02:27:41 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 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	42.334	-5.8	95	0.00
46		2-Chloronaphthalene	40.000	42.309	-5.8	96	0.00
47		2-Nitroaniline	40.000	43.986	-10.0	93	0.00
48		Acenaphthylene	40.000	42.556	-6.4	94	0.00
49		Dimethylphthalate	40.000	41.859	-4.6	93	0.00
50		2,6-Dinitrotoluene	40.000	41.932	-4.8	91	0.00
51	C	Acenaphthene	40.000	41.646	-4.1	94	0.00
52		3-Nitroaniline	40.000	42.435	-6.1	89	0.00
53	P	2,4-Dinitrophenol	40.000	37.172	7.1	83	0.00
54		Dibenzofuran	40.000	41.208	-3.0	93	0.00
55	P	4-Nitrophenol	40.000	42.337	-5.8	92	0.00
56		2,4-Dinitrotoluene	40.000	43.192	-8.0	91	0.00
57		Fluorene	40.000	41.511	-3.8	93	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.346	-3.4	90	0.00
59		Diethylphthalate	40.000	41.761	-4.4	92	0.00
60		4-Chlorophenyl-phenylether	40.000	40.550	-1.4	91	0.00
61		4-Nitroaniline	40.000	45.358	-13.4	94	0.00
62		Azobenzene	40.000	41.628	-4.1	92	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
64		4,6-Dinitro-2-methylphenol	40.000	42.384	-6.0	89	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.544	-6.4	92	0.00
66		4-Bromophenyl-phenylether	40.000	41.511	-3.8	90	0.00
67		Hexachlorobenzene	40.000	42.143	-5.4	93	0.00
68		Atrazine	40.000	42.138	-5.3	90	0.00
69	C	Pentachlorophenol	40.000	40.340	-0.9	84	0.00
70		Phenanthrene	40.000	42.224	-5.6	92	0.00
71		Anthracene	40.000	42.774	-6.9	92	0.00
72		Carbazole	40.000	43.614	-9.0	94	0.00
73		Di-n-butylphthalate	40.000	44.406	-11.0	93	0.00
74	C	Fluoranthene	40.000	43.170	-7.9	94	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
76		Benzidine	40.000	40.880	-2.2	98	0.00
77		Pyrene	40.000	40.925	-2.3	94	0.00
78	S	Terphenyl-d14	80.000	82.119	-2.6	94	0.00
79		Butylbenzylphthalate	40.000	42.892	-7.2	94	0.00
80		Benzo(a)anthracene	40.000	41.377	-3.4	95	0.00
81		3,3'-Dichlorobenzidine	40.000	42.554	-6.4	96	0.00
82		Chrysene	40.000	41.250	-3.1	95	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	43.618	-9.0	96	0.00
84	c	Di-n-octyl phthalate	40.000	43.792	-9.5	95	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.576	-3.9	96	0.00
86	I	Perylene-d12	20.000	20.000	0.0	95	0.00
87		Benzo(b)fluoranthene	40.000	41.608	-4.0	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016667.D
Acq On : 23 Apr 2015 23:46
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SSTDCCC040EC

Quant Time: Apr 24 02:27:41 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 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	43.152	-7.9	97	0.00
89 C	Benzo(a)pyrene	40.000	42.140	-5.4	96	0.00
90	Dibenzo(a,h)anthracene	40.000	42.108	-5.3	95	0.00
91	Benzo(g,h,i)perylene	40.000	42.032	-5.1	95	0.00

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