Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG052815\
Data File : BG017337.D
Acq On : 28 May 2015 22:07
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 29 01:15:51 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 29 01:04:23 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	64	-0.05
2		1,4-Dioxane	40.000	45.413	-13.5	73	-0.06
3		Pyridine	40.000	42.144	-5.4	65	-0.06
4		n-Nitrosodimethylamine	40.000	39.703	0.7	61	-0.06
5	S	2-Fluorophenol	80.000	81.833	-2.3	62	-0.06
6		Aniline	40.000	39.651	0.9	60	-0.05
7	S	Phenol-d6	80.000	80.732	-0.9	61	-0.05
8		2-Chlorophenol	40.000	40.260	-0.6	61	-0.05
9		Benzaldehyde	40.000	38.763	3.1	61	-0.05
10	C	Phenol	40.000	41.414	-3.5	64	-0.05
11		bis(2-Chloroethyl)ether	40.000	41.043	-2.6	61	-0.05
12		1,3-Dichlorobenzene	40.000	39.869	0.3	63	-0.06
13	C	1,4-Dichlorobenzene	40.000	39.428	1.4	61	-0.05
14		1,2-Dichlorobenzene	40.000	39.021	2.4	60	-0.06
15		Benzyl Alcohol	40.000	38.413	4.0	59	-0.05
16		2,2'-oxybis(1-Chloropropane	40.000	42.326	-5.8	66	-0.05
17		2-Methylphenol	40.000	39.908	0.2	63	-0.06
18		Hexachloroethane	40.000	39.290	1.8	61	-0.06
19	P	n-Nitroso-di-n-propylamine	40.000	38.703	3.2	59	-0.05
20		3+4-Methylphenols	40.000	39.957	0.1	60	-0.06
21	I	Naphthalene-d8	20.000	20.000	0.0	59	-0.06
22		Acetophenone	40.000	39.894	0.3	59	-0.06
23	S	Nitrobenzene-d5	80.000	81.839	-2.3	59	-0.05
24		Nitrobenzene	40.000	41.270	-3.2	61	-0.05
25		Isophorone	40.000	40.599	-1.5	58	-0.06
26	C	2-Nitrophenol	40.000	39.662	0.8	58	-0.06
27		2,4-Dimethylphenol	40.000	40.272	-0.7	60	-0.06
28		bis(2-Chloroethoxy)methane	40.000	41.218	-3.0	60	-0.06
29	C	2,4-Dichlorophenol	40.000	41.237	-3.1	59	-0.06
30		1,2,4-Trichlorobenzene	40.000	40.037	-0.1	58	-0.06
31		Naphthalene	40.000	40.382	-1.0	59	-0.06
32		Benzoic acid	40.000	30.658	23.4	44	-0.04
33		4-Chloroaniline	40.000	39.013	2.5	55	-0.06
34	C	Hexachlorobutadiene	40.000	39.895	0.3	58	-0.06
35		Caprolactam	40.000	39.721	0.7	60	-0.04
36	C	4-Chloro-3-methylphenol	40.000	41.354	-3.4	60	-0.06
37		2-Methylnaphthalene	40.000	38.955	2.6	59	-0.06
38	I	Acenaphthene-d10	20.000	20.000	0.0	58	-0.05
39		1,2,4,5-Tetrachlorobenzene	40.000	41.024	-2.6	58	-0.05
40	P	Hexachlorocyclopentadiene	40.000	32.708	18.2	46	-0.06
41	S	2,4,6-Tribromophenol	80.000	78.289	2.1	55	-0.04
42	C	2,4,6-Trichlorophenol	40.000	39.938	0.2	55	-0.05
43		2,4,5-Trichlorophenol	40.000	40.184	-0.5	56	-0.06
44	S	2-Fluorobiphenyl	80.000	81.413	-1.8	57	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG052815\
Data File : BG017337.D
Acq On : 28 May 2015 22:07
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 29 01:15:51 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 29 01:04:23 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	40.257	-0.6	57	-0.05
46		2-Chloronaphthalene	40.000	40.796	-2.0	58	-0.05
47		2-Nitroaniline	40.000	41.846	-4.6	60	-0.05
48		Acenaphthylene	40.000	41.216	-3.0	57	-0.06
49		Dimethylphthalate	40.000	39.795	0.5	56	-0.05
50		2,6-Dinitrotoluene	40.000	41.147	-2.9	57	-0.04
51	C	Acenaphthene	40.000	40.343	-0.9	57	-0.05
52		3-Nitroaniline	40.000	40.584	-1.5	57	-0.05
53	P	2,4-Dinitrophenol	40.000	37.458	6.4	50	-0.04
54		Dibenzofuran	40.000	40.335	-0.8	57	-0.05
55	P	4-Nitrophenol	40.000	38.786	3.0	56	-0.05
56		2,4-Dinitrotoluene	40.000	42.252	-5.6	59	-0.04
57		Fluorene	40.000	39.685	0.8	55	-0.05
58		2,3,4,6-Tetrachlorophenol	40.000	38.003	5.0	52	-0.05
59		Diethylphthalate	40.000	38.786	3.0	56	-0.06
60		4-Chlorophenyl-phenylether	40.000	40.057	-0.1	56	-0.05
61		4-Nitroaniline	40.000	42.600	-6.5	59	-0.04
62		Azobenzene	40.000	41.675	-4.2	59	-0.05
63	I	Phenanthrene-d10	20.000	20.000	0.0	59	-0.05
64		4,6-Dinitro-2-methylphenol	40.000	40.384	-1.0	56	-0.04
65	c	n-Nitrosodiphenylamine	40.000	38.915	2.7	56	-0.04
66		4-Bromophenyl-phenylether	40.000	38.686	3.3	56	-0.05
67		Hexachlorobenzene	40.000	39.418	1.5	57	-0.04
68		Atrazine	40.000	38.363	4.1	54	-0.05
69	C	Pentachlorophenol	40.000	35.471	11.3	50	-0.04
70		Phenanthrene	40.000	40.050	-0.1	57	-0.05
71		Anthracene	40.000	40.319	-0.8	58	-0.05
72		Carbazole	40.000	41.474	-3.7	59	-0.05
73		Di-n-butylphthalate	40.000	41.471	-3.7	60	-0.05
74	C	Fluoranthene	40.000	41.583	-4.0	60	-0.05
75	I	Chrysene-d12	20.000	20.000	0.0	71	-0.04
76		Benzidine	40.000	35.280	11.8	59	-0.05
77		Pyrene	40.000	35.844	10.4	61	-0.04
78	S	Terphenyl-d14	80.000	75.749	5.3	65	-0.05
79		Butylbenzylphthalate	40.000	39.563	1.1	68	-0.04
80		Benzo(a)anthracene	40.000	40.514	-1.3	69	-0.05
81		3,3'-Dichlorobenzidine	40.000	40.394	-1.0	70	-0.05
82		Chrysene	40.000	41.141	-2.9	71	-0.04
83		Bis(2-ethylhexyl)phthalate	40.000	40.812	-2.0	69	-0.05
84	c	Di-n-octyl phthalate	40.000	41.388	-3.5	71	-0.06
85		Indeno(1,2,3-cd)pyrene	40.000	41.704	-4.3	71	-0.11
86	I	Perylene-d12	20.000	20.000	0.0	69	-0.07
87		Benzo(b)fluoranthene	40.000	41.447	-3.6	72	-0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG052815\
Data File : BG017337.D
Acq On : 28 May 2015 22:07
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 29 01:15:51 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 29 01:04:23 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	41.186	-3.0	69	-0.06
89 C	Benzo(a)pyrene	40.000	41.465	-3.7	71	-0.07
90	Dibenzo(a,h)anthracene	40.000	41.612	-4.0	72	-0.11
91	Benzo(g,h,i)perylene	40.000	41.562	-3.9	71	-0.12

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