Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
Data File : BG023050.D
Acq On : 13 Jul 2016 00:47
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 13 01:54:39 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	87	0.00
2		1,4-Dioxane	40.000	39.640	0.9	90	0.00
3		Pyridine	40.000	40.260	-0.6	88	0.00
4		n-Nitrosodimethylamine	40.000	41.280	-3.2	89	0.00
5	S	2-Fluorophenol	80.000	80.655	-0.8	90	0.00
6		Aniline	40.000	40.451	-1.1	87	0.00
7	S	Phenol-d6	80.000	85.409	-6.8	91	0.00
8		2-Chlorophenol	40.000	40.861	-2.2	89	0.00
9		Benzaldehyde	40.000	38.801	3.0	86	0.00
10	C	Phenol	40.000	43.154	-7.9	93	0.00
11		bis(2-Chloroethyl)ether	40.000	39.693	0.8	87	0.00
12		1,3-Dichlorobenzene	40.000	38.867	2.8	88	0.00
13	C	1,4-Dichlorobenzene	40.000	39.412	1.5	87	0.00
14		1,2-Dichlorobenzene	40.000	40.369	-0.9	92	0.00
15		Benzyl Alcohol	40.000	42.378	-5.9	90	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.323	-3.3	91	0.00
17		2-Methylphenol	40.000	42.257	-5.6	92	0.00
18		Hexachloroethane	40.000	40.135	-0.3	94	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.762	0.6	84	0.00
20		3+4-Methylphenols	40.000	43.694	-9.2	90	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	90	0.00
22		Acetophenone	40.000	37.476	6.3	85	0.00
23	S	Nitrobenzene-d5	80.000	80.067	-0.1	91	0.00
24		Nitrobenzene	40.000	39.486	1.3	91	0.00
25		Isophorone	40.000	37.720	5.7	83	0.00
26	C	2-Nitrophenol	40.000	39.908	0.2	85	0.00
27		2,4-Dimethylphenol	40.000	38.459	3.9	88	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.132	4.7	85	0.00
29	C	2,4-Dichlorophenol	40.000	39.988	0.0	89	0.00
30		1,2,4-Trichlorobenzene	40.000	38.953	2.6	89	0.00
31		Naphthalene	40.000	39.685	0.8	91	0.00
32		Benzoic acid	40.000	36.020	9.9	77	0.00
33		4-Chloroaniline	40.000	39.574	1.1	89	0.00
34	C	Hexachlorobutadiene	40.000	37.489	6.3	87	0.00
35		Caprolactam	40.000	35.418	11.5	80	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.525	1.2	88	0.00
37		2-Methylnaphthalene	40.000	39.822	0.4	88	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	86	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.739	0.7	87	0.00
40	P	Hexachlorocyclopentadiene	40.000	42.740	-6.9	93	0.00
41	S	2,4,6-Tribromophenol	80.000	64.853	18.9	72	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.286	4.3	82	0.00
43		2,4,5-Trichlorophenol	40.000	38.691	3.3	85	0.00
44	S	2-Fluorobiphenyl	80.000	79.749	0.3	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
Data File : BG023050.D
Acq On : 13 Jul 2016 00:47
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 13 01:54:39 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	40.840	-2.1	89	0.00
46		2-Chloronaphthalene	40.000	41.235	-3.1	89	0.00
47		2-Nitroaniline	40.000	40.726	-1.8	89	0.00
48		Acenaphthylene	40.000	39.912	0.2	87	0.00
49		Dimethylphthalate	40.000	36.910	7.7	82	0.00
50		2,6-Dinitrotoluene	40.000	37.493	6.3	82	0.00
51	C	Acenaphthene	40.000	39.333	1.7	86	0.00
52		3-Nitroaniline	40.000	38.806	3.0	84	0.00
53	P	2,4-Dinitrophenol	40.000	48.598	-21.5	101	0.00
54		Dibenzofuran	40.000	38.908	2.7	86	0.00
55	P	4-Nitrophenol	40.000	33.896	15.3	73	0.01
56		2,4-Dinitrotoluene	40.000	37.367	6.6	82	0.00
57		Fluorene	40.000	38.116	4.7	85	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	35.053	12.4	76	0.00
59		Diethylphthalate	40.000	37.220	7.0	82	0.00
60		4-Chlorophenyl-phenylether	40.000	36.799	8.0	80	0.00
61		4-Nitroaniline	40.000	36.371	9.1	80	0.00
62		Azobenzene	40.000	40.153	-0.4	88	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	79	0.00
64		4,6-Dinitro-2-methylphenol	40.000	35.974	10.1	68	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.051	-5.1	83	0.00
66		4-Bromophenyl-phenylether	40.000	39.001	2.5	77	0.00
67		Hexachlorobenzene	40.000	38.076	4.8	75	0.00
68		Atrazine	40.000	38.572	3.6	77	0.00
69	C	Pentachlorophenol	40.000	36.162	9.6	68	0.00
70		Phenanthrene	40.000	40.839	-2.1	82	0.00
71		Anthracene	40.000	40.566	-1.4	82	0.00
72		Carbazole	40.000	40.427	-1.1	81	0.00
73		Di-n-butylphthalate	40.000	43.102	-7.8	84	0.00
74	C	Fluoranthene	40.000	39.069	2.3	82	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	78	0.02
76		Benzidine	40.000	53.473	-33.7#	102	0.01
77		Pyrene	40.000	39.532	1.2	82	0.00
78	S	Terphenyl-d14	80.000	86.789	-8.5	83	0.00
79		Butylbenzylphthalate	40.000	42.396	-6.0	84	0.02
80		Benzo(a)anthracene	40.000	40.735	-1.8	82	0.02
81		3,3'-Dichlorobenzidine	40.000	39.358	1.6	76	0.02
82		Chrysene	40.000	40.692	-1.7	82	0.02
83		Bis(2-ethylhexyl)phthalate	40.000	41.674	-4.2	84	0.02
84	c	Di-n-octyl phthalate	40.000	42.596	-6.5	84	0.04
85		Indeno(1,2,3-cd)pyrene	40.000	36.653	8.4	74	0.11
86	I	Perylene-d12	20.000	20.000	0.0	77	0.06
87		Benzo(b)fluoranthene	40.000	40.090	-0.2	79	0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
Data File : BG023050.D
Acq On : 13 Jul 2016 00:47
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 13 01:54:39 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	39.824	0.4	76	0.06
89 C	Benzo(a)pyrene	40.000	39.537	1.2	77	0.06
90	Dibenzo(a,h)anthracene	40.000	37.450	6.4	73	0.11
91	Benzo(g,h,i)perylene	40.000	36.133	9.7	70	0.13

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