Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021226.D
Acq On : 3 Mar 2016 00:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 01:09:26 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	98	-0.02
2		1,4-Dioxane	40.000	33.143	17.1	86	0.00
3		Pyridine	40.000	35.181	12.0	81	0.00
4		n-Nitrosodimethylamine	40.000	30.320	24.2	70	0.00
5	S	2-Fluorophenol	80.000	75.930	5.1	92	0.00
6		Aniline	40.000	35.531	11.2	85	0.00
7	S	Phenol-d6	80.000	74.271	7.2	87	0.00
8		2-Chlorophenol	40.000	39.287	1.8	93	0.00
9		Benzaldehyde	40.000	34.079	14.8	88	-0.02
10	C	Phenol	40.000	36.972	7.6	89	0.00
11		bis(2-Chloroethyl)ether	40.000	35.988	10.0	84	-0.02
12		1,3-Dichlorobenzene	40.000	39.727	0.7	96	-0.02
13	C	1,4-Dichlorobenzene	40.000	40.080	-0.2	97	-0.02
14		1,2-Dichlorobenzene	40.000	40.541	-1.4	96	0.00
15		Benzyl Alcohol	40.000	35.309	11.7	82	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	32.262	19.3	78	0.00
17		2-Methylphenol	40.000	36.761	8.1	88	0.00
18		Hexachloroethane	40.000	39.050	2.4	94	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	32.166	19.6	77	-0.02
20		3+4-Methylphenols	40.000	36.101	9.7	87	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	91	0.00
22		Acetophenone	40.000	36.848	7.9	83	0.00
23	S	Nitrobenzene-d5	80.000	75.506	5.6	83	0.00
24		Nitrobenzene	40.000	37.940	5.2	85	0.00
25		Isophorone	40.000	33.713	15.7	76	0.00
26	C	2-Nitrophenol	40.000	39.889	0.3	89	0.00
27		2,4-Dimethylphenol	40.000	36.749	8.1	82	0.00
28		bis(2-Chloroethoxy)methane	40.000	35.755	10.6	82	0.00
29	C	2,4-Dichlorophenol	40.000	39.122	2.2	89	0.00
30		1,2,4-Trichlorobenzene	40.000	42.673	-6.7	96	0.00
31		Naphthalene	40.000	39.599	1.0	90	0.00
32		Benzoic acid	40.000	32.564	18.6	71	0.00
33		4-Chloroaniline	40.000	36.335	9.2	82	0.00
34	C	Hexachlorobutadiene	40.000	45.650	-14.1	102	-0.02
35		Caprolactam	40.000	30.814	23.0	70	0.00
36	C	4-Chloro-3-methylphenol	40.000	34.599	13.5	79	0.00
37		2-Methylnaphthalene	40.000	37.794	5.5	86	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	82	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	45.865	-14.7	91	0.00
40	P	Hexachlorocyclopentadiene	40.000	25.296	36.8#	48	0.00
41	S	2,4,6-Tribromophenol	80.000	77.224	3.5	78	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.860	-2.1	84	0.00
43		2,4,5-Trichlorophenol	40.000	40.292	-0.7	82	0.00
44	S	2-Fluorobiphenyl	80.000	84.841	-6.1	85	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021226.D
Acq On : 3 Mar 2016 00:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 01:09:26 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	41.539	-3.8	84	0.00
46		2-Chloronaphthalene	40.000	41.603	-4.0	83	0.00
47		2-Nitroaniline	40.000	34.682	13.3	70	0.00
48		Acenaphthylene	40.000	39.247	1.9	79	0.00
49		Dimethylphthalate	40.000	36.479	8.8	75	0.00
50		2,6-Dinitrotoluene	40.000	37.546	6.1	76	0.00
51	C	Acenaphthene	40.000	36.665	8.3	72	0.00
52		3-Nitroaniline	40.000	34.953	12.6	71	0.00
53	P	2,4-Dinitrophenol	40.000	26.606	33.5#	45	0.00
54		Dibenzofuran	40.000	38.322	4.2	79	0.00
55	P	4-Nitrophenol	40.000	35.922	10.2	71	0.00
56		2,4-Dinitrotoluene	40.000	36.430	8.9	73	0.00
57		Fluorene	40.000	37.068	7.3	75	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	36.594	8.5	75	0.00
59		Diethylphthalate	40.000	34.395	14.0	71	0.00
60		4-Chlorophenyl-phenylether	40.000	38.406	4.0	78	0.00
61		4-Nitroaniline	40.000	35.787	10.5	72	0.00
62		Azobenzene	40.000	33.977	15.1	69	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	75	0.00
64		4,6-Dinitro-2-methylphenol	40.000	30.620	23.4	53	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.646	-1.6	74	0.00
66		4-Bromophenyl-phenylether	40.000	41.737	-4.3	75	0.00
67		Hexachlorobenzene	40.000	42.652	-6.6	77	0.00
68		Atrazine	40.000	41.493	-3.7	76	0.00
69	C	Pentachlorophenol	40.000	38.902	2.7	68	0.00
70		Phenanthrene	40.000	40.134	-0.3	74	0.00
71		Anthracene	40.000	40.997	-2.5	75	0.00
72		Carbazole	40.000	40.618	-1.5	75	0.00
73		Di-n-butylphthalate	40.000	39.406	1.5	71	0.00
74	C	Fluoranthene	40.000	40.344	-0.9	74	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	74	0.00
76		Benzidine	40.000	40.580	-1.4	72	0.00
77		Pyrene	40.000	41.257	-3.1	75	0.00
78	S	Terphenyl-d14	80.000	82.427	-3.0	74	0.00
79		Butylbenzylphthalate	40.000	39.043	2.4	70	0.00
80		Benzo(a)anthracene	40.000	40.265	-0.7	73	0.00
81		3,3'-Dichlorobenzidine	40.000	40.584	-1.5	72	0.00
82		Chrysene	40.000	40.491	-1.2	73	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	38.645	3.4	69	0.00
84	c	Di-n-octyl phthalate	40.000	39.182	2.0	70	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	38.551	3.6	72	0.00
86	I	Perylene-d12	20.000	20.000	0.0	76	0.00
87		Benzo(b)fluoranthene	40.000	39.895	0.3	74	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021226.D
Acq On : 3 Mar 2016 00:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 01:09:26 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	40.578	-1.4	76	0.00
89 C	Benzo(a)pyrene	40.000	40.909	-2.3	76	0.00
90	Dibenzo(a,h)anthracene	40.000	36.001	10.0	68	0.00
91	Benzo(g,h,i)perylene	40.000	34.690	13.3	67	-0.02

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