Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062118\
Data File : BG035314.D
Acq On : 22 Jun 2018 1:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 22 05:35:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 2018
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	156	-0.02
2		1,4-Dioxane	40.000	38.754	3.1	152	-0.01
3		Pyridine	40.000	39.275	1.8	143	0.00
4		n-Nitrosodimethylamine	40.000	35.341	11.6	130	-0.01
5	S	2-Fluorophenol	80.000	76.865	3.9	146	-0.02
6		Aniline	40.000	36.267	9.3	139	-0.02
7	S	Phenol-d6	80.000	76.273	4.7	144	-0.01
8		2-Chlorophenol	40.000	36.823	7.9	139	-0.02
9		Benzaldehyde	40.000	31.158	22.1	115	-0.02
10	C	Phenol	40.000	36.844	7.9	141	-0.02
11		bis(2-Chloroethyl)ether	40.000	36.736	8.2	141	-0.02
12		1,3-Dichlorobenzene	40.000	36.837	7.9	145	-0.02
13	C	1,4-Dichlorobenzene	40.000	37.983	5.0	146	-0.02
14		1,2-Dichlorobenzene	40.000	38.562	3.6	147	-0.02
15		Benzyl Alcohol	40.000	35.446	11.4	134	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	33.815	15.5	131	-0.02
17		2-Methylphenol	40.000	38.103	4.7	141	-0.02
18		Hexachloroethane	40.000	37.022	7.4	139	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	34.659	13.4	131	-0.02
20		3+4-Methylphenols	40.000	37.745	5.6	143	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	152	-0.02
22		Acetophenone	40.000	36.359	9.1	137	-0.02
23	S	Nitrobenzene-d5	80.000	82.662	-3.3	148	-0.02
24		Nitrobenzene	40.000	40.064	-0.2	145	-0.02
25		Isophorone	40.000	35.074	12.3	132	-0.02
26	C	2-Nitrophenol	40.000	44.434	-11.1	167	-0.02
27		2,4-Dimethylphenol	40.000	37.201	7.0	141	-0.02
28		bis(2-Chloroethoxy)methane	40.000	37.267	6.8	139	-0.02
29	C	2,4-Dichlorophenol	40.000	39.652	0.9	146	-0.02
30		1,2,4-Trichlorobenzene	40.000	40.255	-0.6	147	-0.02
31		Naphthalene	40.000	37.903	5.2	143	-0.02
32		Benzoic acid	40.000	40.626	-1.6	153	0.00
33		4-Chloroaniline	40.000	37.638	5.9	139	-0.02
34	C	Hexachlorobutadiene	40.000	40.506	-1.3	152	-0.03
35		Caprolactam	40.000	37.898	5.3	143	-0.02
36	C	4-Chloro-3-methylphenol	40.000	39.149	2.1	145	-0.01
37		2-Methylnaphthalene	40.000	38.036	4.9	140	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	159	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	39.221	1.9	151	-0.02
40	P	Hexachlorocyclopentadiene	40.000	36.776	8.1	138	-0.03
41	S	2,4,6-Tribromophenol	80.000	84.044	-5.1	164	-0.02
42	C	2,4,6-Trichlorophenol	40.000	40.037	-0.1	152	-0.02
43		2,4,5-Trichlorophenol	40.000	40.663	-1.7	160	-0.02
44	S	2-Fluorobiphenyl	80.000	73.522	8.1	145	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062118\
Data File : BG035314.D
Acq On : 22 Jun 2018 1:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 22 05:35:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 2018
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	37.010	7.5	144	-0.02
46		2-Chloronaphthalene	40.000	36.670	8.3	146	-0.02
47		2-Nitroaniline	40.000	40.499	-1.2	159	-0.02
48		Acenaphthylene	40.000	36.517	8.7	143	-0.02
49		Dimethylphthalate	40.000	36.391	9.0	144	-0.03
50		2,6-Dinitrotoluene	40.000	42.761	-6.9	164	-0.02
51	C	Acenaphthene	40.000	37.950	5.1	148	-0.02
52		3-Nitroaniline	40.000	43.613	-9.0	162	-0.01
53	P	2,4-Dinitrophenol	40.000	47.788	-19.5	182	-0.02
54		Dibenzofuran	40.000	36.609	8.5	143	-0.02
55	P	4-Nitrophenol	40.000	37.175	7.1	137	0.00
56		2,4-Dinitrotoluene	40.000	45.116	-12.8	172	-0.02
57		Fluorene	40.000	36.178	9.6	141	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	41.434	-3.6	160	-0.02
59		Diethylphthalate	40.000	34.848	12.9	138	-0.03
60		4-Chlorophenyl-phenylether	40.000	37.958	5.1	150	-0.03
61		4-Nitroaniline	40.000	41.544	-3.9	155	-0.02
62		Azobenzene	40.000	33.406	16.5	132	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	160	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	48.464	-21.2	190	-0.02
65	c	n-Nitrosodiphenylamine	40.000	36.763	8.1	142	-0.02
66		4-Bromophenyl-phenylether	40.000	38.861	2.8	154	-0.03
67		Hexachlorobenzene	40.000	39.435	1.4	157	-0.02
68		Atrazine	40.000	36.076	9.8	140	-0.03
69	C	Pentachlorophenol	40.000	39.790	0.5	152	-0.02
70		Phenanthrene	40.000	36.977	7.6	144	-0.02
71		Anthracene	40.000	37.404	6.5	145	-0.02
72		Carbazole	40.000	36.458	8.9	141	-0.02
73		Di-n-butylphthalate	40.000	35.480	11.3	138	-0.03
74	C	Fluoranthene	40.000	36.845	7.9	143	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	165	-0.03
76		Benzidine	40.000	23.303	41.7#	94	-0.02
77		Pyrene	40.000	35.572	11.1	145	-0.02
78	S	Terphenyl-d14	80.000	71.618	10.5	145	-0.02
79		Butylbenzylphthalate	40.000	35.182	12.0	138	-0.03
80		Benzo(a)anthracene	40.000	37.084	7.3	152	-0.02
81		3,3'-Dichlorobenzidine	40.000	36.420	8.9	144	-0.03
82		Chrysene	40.000	37.265	6.8	152	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	34.045	14.9	136	-0.05
84	c	Di-n-octyl phthalate	40.000	34.267	14.3	136	-0.06
85		Indeno(1,2,3-cd)pyrene	40.000	37.898	5.3	153	-0.07
86	I	Perylene-d12	20.000	20.000	0.0	165	-0.05
87		Benzo(b)fluoranthene	40.000	36.982	7.5	149	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062118\
Data File : BG035314.D
Acq On : 22 Jun 2018 1:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 22 05:35:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 08 17:16:28 2018
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	36.506	8.7	154	-0.04
89 C	Benzo(a)pyrene	40.000	36.714	8.2	150	-0.05
90	Dibenzo(a,h)anthracene	40.000	37.132	7.2	153	-0.09
91	Benzo(g,h,i)perylene	40.000	37.633	5.9	154	-0.07

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