Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG040918\
Data File : BG033691.D
Acq On : 9 Apr 2018 22:55
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 07:24:50 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 05 17:36:07 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	132	0.00
2		1,4-Dioxane	0.542	0.594	-9.6	146	0.00
3		Pyridine	1.497	1.610	-7.5	128	0.00
4		n-Nitrosodimethylamine	0.614	0.659	-7.3	140	0.00
5	S	2-Fluorophenol	1.067	1.144	-7.2	130	0.00
6		Aniline	2.155	2.234	-3.7	126	0.00
7	S	Phenol-d6	1.833	1.891	-3.2	131	0.00
8		2-Chlorophenol	1.219	1.252	-2.7	124	0.00
9		Benzaldehyde	1.165	0.959	17.7	111	0.00
10	C	Phenol	1.809	1.862	-2.9	128	0.00
11		bis(2-Chloroethyl)ether	1.477	1.395	5.6	114	0.00
12		1,3-Dichlorobenzene	1.594	1.597	-0.2	129	0.00
13	C	1,4-Dichlorobenzene	1.597	1.491	6.6	123	0.00
14		1,2-Dichlorobenzene	1.558	1.474	5.4	119	0.00
15		Benzyl Alcohol	1.443	1.489	-3.2	127	0.00
16		2,2'-oxybis(1-Chloropropane	2.408	2.575	-6.9	138	0.00
17		2-Methylphenol	1.233	1.264	-2.5	132	0.00
18		Hexachloroethane	0.616	0.622	-1.0	132	0.00
19	P	n-Nitroso-di-n-propylamine	1.343	1.352	-0.7	121	0.00
20		3+4-Methylphenols	1.755	1.759	-0.2	122	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	120	0.00
22		Acetophenone	0.548	0.563	-2.7	115	0.00
23	S	Nitrobenzene-d5	0.435	0.452	-3.9	121	0.00
24		Nitrobenzene	0.466	0.461	1.1	115	0.00
25		Isophorone	0.845	0.850	-0.6	117	0.00
26	C	2-Nitrophenol	0.176	0.181	-2.8	114	0.00
27		2,4-Dimethylphenol	0.256	0.264	-3.1	126	0.00
28		bis(2-Chloroethoxy)methane	0.422	0.431	-2.1	118	0.00
29	C	2,4-Dichlorophenol	0.315	0.321	-1.9	116	0.00
30		1,2,4-Trichlorobenzene	0.409	0.392	4.2	113	0.00
31		Naphthalene	1.036	1.018	1.7	116	0.00
32		Benzoic acid	0.238	0.209	12.2	118	0.01
33		4-Chloroaniline	0.464	0.440	5.2	109	0.00
34	C	Hexachlorobutadiene	0.310	0.293	5.5	116	0.00
35		Caprolactam	0.123	0.130	-5.7	121	0.00
36	C	4-Chloro-3-methylphenol	0.411	0.412	-0.2	112	0.00
37		2-Methylnaphthalene	0.804	0.758	5.7	114	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	117	0.00
39		1,2,4,5-Tetrachlorobenzene	0.724	0.658	9.1	105	0.00
40	P	Hexachlorocyclopentadiene	0.425	0.312	26.6#	83	0.00
41	S	2,4,6-Tribromophenol	0.299	0.264	11.7	99	0.00
42	C	2,4,6-Trichlorophenol	0.421	0.397	5.7	104	0.00
43		2,4,5-Trichlorophenol	0.498	0.477	4.2	102	0.00
44	S	2-Fluorobiphenyl	1.385	1.318	4.8	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG040918\
Data File : BG033691.D
Acq On : 9 Apr 2018 22:55
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 07:24:50 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 05 17:36:07 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.537	1.476	4.0	110	0.00
46		2-Chloronaphthalene	1.185	1.140	3.8	109	0.00
47		2-Nitroaniline	0.444	0.468	-5.4	117	0.00
48		Acenaphthylene	1.978	1.926	2.6	112	0.00
49		Dimethylphthalate	1.741	1.658	4.8	110	0.00
50		2,6-Dinitrotoluene	0.356	0.347	2.5	107	0.00
51	C	Acenaphthene	1.275	1.225	3.9	109	0.00
52		3-Nitroaniline	0.373	0.366	1.9	111	0.00
53	P	2,4-Dinitrophenol	0.149	0.087	41.6#	62	0.00
54		Dibenzofuran	1.883	1.768	6.1	108	0.00
55	P	4-Nitrophenol	0.262	0.247	5.7	105	0.00
56		2,4-Dinitrotoluene	0.524	0.510	2.7	111	0.00
57		Fluorene	1.604	1.531	4.6	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.468	0.398	15.0	94	0.00
59		Diethylphthalate	1.690	1.687	0.2	115	0.00
60		4-Chlorophenyl-phenylether	0.922	0.841	8.8	107	0.00
61		4-Nitroaniline	0.378	0.377	0.3	113	0.00
62		Azobenzene	1.637	1.636	0.1	114	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
64		4,6-Dinitro-2-methylphenol	0.120	0.105	12.5	95	0.00
65	c	n-Nitrosodiphenylamine	0.571	0.563	1.4	111	0.00
66		4-Bromophenyl-phenylether	0.235	0.222	5.5	105	0.00
67		Hexachlorobenzene	0.248	0.234	5.6	107	0.00
68		Atrazine	0.231	0.232	-0.4	111	0.00
69	C	Pentachlorophenol	0.170	0.137	19.4	90	0.00
70		Phenanthrene	1.042	1.016	2.5	110	0.00
71		Anthracene	1.055	1.052	0.3	112	0.00
72		Carbazole	0.935	0.962	-2.9	115	0.00
73		Di-n-butylphthalate	1.088	1.161	-6.7	119	0.00
74	C	Fluoranthene	1.289	1.274	1.2	111	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	116	0.00
76		Benzidine	0.542	0.575	-6.1	113	0.00
77		Pyrene	1.334	1.298	2.7	112	0.00
78	S	Terphenyl-d14	0.935	0.879	6.0	109	0.00
79		Butylbenzylphthalate	0.492	0.527	-7.1	122	0.00
80		Benzo(a)anthracene	1.274	1.236	3.0	113	0.00
81		3,3'-Dichlorobenzidine	0.453	0.463	-2.2	113	0.00
82		Chrysene	1.233	1.196	3.0	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.679	0.737	-8.5	124	0.00
84	c	Di-n-octyl phthalate	1.039	1.171	-12.7	127	0.00
85		Indeno(1,2,3-cd)pyrene	1.333	1.382	-3.7	118	0.00
86	I	Perylene-d12	1.000	1.000	0.0	120	0.00
87		Benzo(b)fluoranthene	1.155	1.090	5.6	112	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG040918\
Data File : BG033691.D
Acq On : 9 Apr 2018 22:55
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 10 07:24:50 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 05 17:36:07 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.169	1.132	3.2	114	0.00
89 C	Benzo(a)pyrene	1.110	1.085	2.3	115	0.00
90	Dibenzo(a,h)anthracene	1.045	1.057	-1.1	116	0.00
91	Benzo(g,h,i)perylene	1.035	1.043	-0.8	117	-0.02

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