Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
Data File : BF092483.D
Acq On : 25 Jan 2017 18:17
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 10:06:46 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 24 13:48:07 2017
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	107	0.00
2		1,4-Dioxane	0.684	0.648	5.3	103	-0.01
3		Pyridine	1.946	1.824	6.3	102	0.00
4		n-Nitrosodimethylamine	0.955	0.874	8.5	98	0.00
5	S	2-Fluorophenol	1.285	1.241	3.4	106	0.00
6		Aniline	2.442	2.393	2.0	103	0.00
7	S	Phenol-d6	1.637	1.549	5.4	103	0.00
8		2-Chlorophenol	1.350	1.354	-0.3	107	0.00
9		Benzaldehyde	1.162	1.051	9.6	100	0.00
10	C	Phenol	1.976	1.972	0.2	103	0.00
11		bis(2-Chloroethyl)ether	1.479	1.538	-4.0	106	0.00
12		1,3-Dichlorobenzene	1.597	1.633	-2.3	109	0.00
13	C	1,4-Dichlorobenzene	1.607	1.664	-3.5	106	0.00
14		1,2-Dichlorobenzene	1.525	1.410	7.5	106	0.00
15		Benzyl Alcohol	1.234	1.147	7.1	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.925	2.811	3.9	104	0.00
17		2-Methylphenol	1.154	1.126	2.4	103	0.00
18		Hexachloroethane	0.554	0.566	-2.2	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.115	1.068	4.2	105	0.00
20		3+4-Methylphenols	1.400	1.445	-3.2	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.476	0.464	2.5	103	-0.01
23	S	Nitrobenzene-d5	0.366	0.375	-2.5	108	0.00
24		Nitrobenzene	0.384	0.394	-2.6	100	0.00
25		Isophorone	0.721	0.709	1.7	104	0.00
26	C	2-Nitrophenol	0.161	0.172	-6.8	115	0.00
27		2,4-Dimethylphenol	0.334	0.316	5.4	101	-0.01
28		bis(2-Chloroethoxy)methane	0.474	0.418	11.8	100	0.00
29	C	2,4-Dichlorophenol	0.290	0.280	3.4	103	0.00
30		1,2,4-Trichlorobenzene	0.313	0.311	0.6	111	0.00
31		Naphthalene	1.024	0.956	6.6	104	0.00
32		Benzoic acid	0.231	0.228	1.3	95	0.00
33		4-Chloroaniline	0.433	0.391	9.7	101	0.00
34	C	Hexachlorobutadiene	0.171	0.180	-5.3	111	0.00
35		Caprolactam	0.097	0.093	4.1	100	0.00
36	C	4-Chloro-3-methylphenol	0.313	0.317	-1.3	103	-0.01
37		2-Methylnaphthalene	0.648	0.619	4.5	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.504	0.536	-6.3	112	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.267	-5.5	108	0.00
41	S	2,4,6-Tribromophenol	0.136	0.145	-6.6	116	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.383	1.0	110	-0.01
43		2,4,5-Trichlorophenol	0.354	0.366	-3.4	107	0.00
44	S	2-Fluorobiphenyl	1.082	1.021	5.6	100	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
Data File : BF092483.D
Acq On : 25 Jan 2017 18:17
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 10:06:46 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 24 13:48:07 2017
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	1.446	1.254	13.3	103	0.00
46		2-Chloronaphthalene	1.181	1.052	10.9	99	0.00
47		2-Nitroaniline	0.398	0.359	9.8	98	0.00
48		Acenaphthylene	1.732	1.809	-4.4	113	0.00
49		Dimethylphthalate	1.272	1.304	-2.5	108	0.00
50		2,6-Dinitrotoluene	0.260	0.299	-15.0	111	0.00
51	C	Acenaphthene	1.076	1.069	0.7	117	0.00
52		3-Nitroaniline	0.326	0.358	-9.8	101	0.00
53	P	2,4-Dinitrophenol	0.093	0.124	-33.3#	130	-0.01
54		Dibenzofuran	1.461	1.476	-1.0	116	0.00
55	P	4-Nitrophenol	0.259	0.261	-0.8	95	0.00
56		2,4-Dinitrotoluene	0.345	0.383	-11.0	124	0.00
57		Fluorene	1.084	1.141	-5.3	123	0.00
58		2,3,4,6-Tetrachlorophenol	0.265	0.283	-6.8	117	0.00
59		Diethylphthalate	1.220	1.230	-0.8	113	0.00
60		4-Chlorophenyl-phenylether	0.526	0.481	8.6	99	-0.01
61		4-Nitroaniline	0.326	0.324	0.6	104	-0.01
62		Azobenzene	1.388	1.276	8.1	95	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	-0.01
64		4,6-Dinitro-2-methylphenol	0.099	0.130	-31.3#	124	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.636	-1.8	114	0.00
66		4-Bromophenyl-phenylether	0.202	0.194	4.0	103	-0.01
67		Hexachlorobenzene	0.204	0.212	-3.9	116	0.00
68		Atrazine	0.214	0.229	-7.0	109	-0.01
69	C	Pentachlorophenol	0.131	0.131	0.0	100	-0.01
70		Phenanthrene	1.101	1.031	6.4	102	-0.01
71		Anthracene	1.076	1.088	-1.1	110	0.00
72		Carbazole	1.028	1.009	1.8	102	-0.01
73		Di-n-butylphthalate	1.175	1.103	6.1	102	-0.01
74	C	Fluoranthene	1.077	1.095	-1.7	113	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	107	-0.01
76		Benzidine	0.742	0.743	-0.1	101	-0.02
77		Pyrene	1.449	1.303	10.1	108	-0.01
78	S	Terphenyl-d14	0.751	0.724	3.6	117	-0.01
79		Butylbenzylphthalate	0.587	0.555	5.5	96	-0.02
80		Benzo(a)anthracene	1.075	1.016	5.5	107	-0.01
81		3,3'-Dichlorobenzidine	0.408	0.417	-2.2	107	-0.01
82		Chrysene	1.148	0.973	15.2	100	-0.01
83		Bis(2-ethylhexyl)phthalate	0.739	0.719	2.7	105	-0.02
84	c	Di-n-octyl phthalate	1.342	1.394	-3.9	104	-0.02
85		Indeno(1,2,3-cd)pyrene	0.849	0.815	4.0	106	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
87		Benzo(b)fluoranthene	1.284	1.185	7.7	88	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
Data File : BF092483.D
Acq On : 25 Jan 2017 18:17
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 10:06:46 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 24 13:48:07 2017
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	1.068	1.187	-11.1	123	-0.01
89 C	Benzo(a)pyrene	1.086	1.095	-0.8	100	-0.02
90	Dibenzo(a,h)anthracene	0.878	0.905	-3.1	105	-0.02
91	Benzo(g,h,i)perylene	0.888	0.892	-0.5	107	-0.02

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