Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080516\
Data File : BF089624.D
Acq On : 5 Aug 2016 11:27
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 23:11:21 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	104	0.00
2		1,4-Dioxane	0.684	0.601	12.1	108	0.01
3		Pyridine	1.257	1.299	-3.3	111	0.00
4		n-Nitrosodimethylamine	0.492	0.513	-4.3	101	0.00
5	S	2-Fluorophenol	1.193	1.167	2.2	104	0.01
6		Aniline	2.098	2.180	-3.9	107	0.00
7	S	Phenol-d6	1.619	1.602	1.1	105	0.01
8		2-Chlorophenol	1.465	1.441	1.6	104	0.00
9		Benzaldehyde	0.587	0.626	-6.6	104	0.00
10	C	Phenol	1.662	1.709	-2.8	105	0.00
11		bis(2-Chloroethyl)ether	1.428	1.388	2.8	106	0.00
12		1,3-Dichlorobenzene	1.591	1.514	4.8	104	0.00
13	C	1,4-Dichlorobenzene	1.587	1.514	4.6	107	0.00
14		1,2-Dichlorobenzene	1.673	1.547	7.5	105	0.00
15		Benzyl Alcohol	1.061	1.036	2.4	99	0.00
16		2,2'-oxybis(1-Chloropropane	1.818	1.703	6.3	104	0.00
17		2-Methylphenol	1.058	1.029	2.7	104	0.00
18		Hexachloroethane	0.669	0.622	7.0	104	0.00
19	P	n-Nitroso-di-n-propylamine	1.169	1.105	5.5	100	0.00
20		3+4-Methylphenols	1.336	1.379	-3.2	106	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.477	0.464	2.7	94	0.00
23	S	Nitrobenzene-d5	0.362	0.364	-0.6	103	0.00
24		Nitrobenzene	0.343	0.356	-3.8	104	0.00
25		Isophorone	0.692	0.663	4.2	104	0.00
26	C	2-Nitrophenol	0.158	0.161	-1.9	105	0.00
27		2,4-Dimethylphenol	0.283	0.297	-4.9	107	0.01
28		bis(2-Chloroethoxy)methane	0.419	0.429	-2.4	105	0.00
29	C	2,4-Dichlorophenol	0.268	0.277	-3.4	108	0.00
30		1,2,4-Trichlorobenzene	0.306	0.301	1.6	103	0.00
31		Naphthalene	1.063	1.018	4.2	104	0.00
32		Benzoic acid	0.178	0.169	5.1	97	0.01
33		4-Chloroaniline	0.399	0.416	-4.3	103	0.00
34	C	Hexachlorobutadiene	0.176	0.168	4.5	103	0.00
35		Caprolactam	0.078	0.081	-3.8	105	0.01
36	C	4-Chloro-3-methylphenol	0.312	0.315	-1.0	101	0.00
37		2-Methylnaphthalene	0.654	0.641	2.0	105	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	0.752	0.748	0.5	104	0.00
40	P	Hexachlorocyclopentadiene	0.204	0.232	-13.7	101	0.00
41	S	2,4,6-Tribromophenol	0.165	0.168	-1.8	101	0.00
42	C	2,4,6-Trichlorophenol	0.371	0.369	0.5	101	0.00
43		2,4,5-Trichlorophenol	0.396	0.419	-5.8	110	0.00
44	S	2-Fluorobiphenyl	1.539	1.443	6.2	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080516\
Data File : BF089624.D
Acq On : 5 Aug 2016 11:27
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 23:11:21 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.804	1.765	2.2	104	0.00
46		2-Chloronaphthalene	1.352	1.312	3.0	104	0.00
47		2-Nitroaniline	0.436	0.440	-0.9	106	0.01
48		Acenaphthylene	2.228	2.137	4.1	104	0.00
49		Dimethylphthalate	1.568	1.510	3.7	103	0.00
50		2,6-Dinitrotoluene	0.312	0.322	-3.2	101	0.00
51	C	Acenaphthene	1.287	1.226	4.7	102	0.00
52		3-Nitroaniline	0.373	0.373	0.0	102	0.00
53	P	2,4-Dinitrophenol	0.105	0.080	23.8	80	0.01
54		Dibenzofuran	1.769	1.729	2.3	104	0.00
55	P	4-Nitrophenol	0.192	0.173	9.9	83	0.02
56		2,4-Dinitrotoluene	0.438	0.435	0.7	102	0.00
57		Fluorene	1.512	1.445	4.4	103	0.00
58		2,3,4,6-Tetrachlorophenol	0.301	0.289	4.0	100	0.00
59		Diethylphthalate	1.621	1.591	1.9	104	0.00
60		4-Chlorophenyl-phenylether	0.693	0.686	1.0	103	0.00
61		4-Nitroaniline	0.335	0.331	1.2	98	0.00
62		Azobenzene	1.568	1.593	-1.6	103	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	0.112	0.104	7.1	91	0.01
65	c	n-Nitrosodiphenylamine	0.675	0.656	2.8	102	0.00
66		4-Bromophenyl-phenylether	0.193	0.194	-0.5	101	0.00
67		Hexachlorobenzene	0.213	0.201	5.6	103	0.00
68		Atrazine	0.189	0.194	-2.6	98	0.00
69	C	Pentachlorophenol	0.081	0.086	-6.2	90	0.00
70		Phenanthrene	1.088	1.053	3.2	103	0.00
71		Anthracene	1.166	1.091	6.4	101	0.00
72		Carbazole	0.972	0.954	1.9	100	0.00
73		Di-n-butylphthalate	1.314	1.276	2.9	100	0.00
74	C	Fluoranthene	1.118	1.030	7.9	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.606	0.626	-3.3	95	0.00
77		Pyrene	1.768	1.655	6.4	99	0.00
78	S	Terphenyl-d14	1.013	0.926	8.6	98	0.00
79		Butylbenzylphthalate	0.752	0.750	0.3	96	0.00
80		Benzo(a)anthracene	1.149	1.082	5.8	96	0.00
81		3,3'-Dichlorobenzidine	0.362	0.354	2.2	93	0.00
82		Chrysene	1.206	1.110	8.0	93	0.00
83		Bis(2-ethylhexyl)phthalate	1.042	0.991	4.9	95	0.00
84	c	Di-n-octyl phthalate	1.499	1.501	-0.1	95	-0.01
85		Indeno(1,2,3-cd)pyrene	0.915	0.879	3.9	104	0.00
86	I	Perylene-d12	1.000	1.000	0.0	98	0.00
87		Benzo(b)fluoranthene	1.230	1.165	5.3	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080516\
Data File : BF089624.D
Acq On : 5 Aug 2016 11:27
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 05 23:11:21 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 01:39:23 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.241	1.228	1.0	100	0.00
89 C	Benzo(a)pyrene	1.152	1.085	5.8	97	0.00
90	Dibenzo(a,h)anthracene	1.076	1.064	1.1	104	0.00
91	Benzo(g,h,i)perylene	1.101	1.118	-1.5	107	0.01

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