Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122216\
Data File : BF091877.D
Acq On : 22 Dec 2016 11:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 23 00:56:27 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 16:17:51 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	105	0.00
2		1,4-Dioxane	0.590	0.633	-7.3	111	0.00
3		Pyridine	2.058	1.940	5.7	96	-0.01
4		n-Nitrosodimethylamine	0.776	0.691	11.0	89	-0.01
5	S	2-Fluorophenol	1.198	1.107	7.6	100	0.00
6		Aniline	1.899	1.818	4.3	100	-0.01
7	S	Phenol-d6	1.462	1.490	-1.9	105	0.00
8		2-Chlorophenol	1.362	1.309	3.9	99	-0.01
9		Benzaldehyde	0.904	0.876	3.1	109	0.00
10	C	Phenol	1.666	1.818	-9.1	105	0.00
11		bis(2-Chloroethyl)ether	1.326	1.297	2.2	95	-0.01
12		1,3-Dichlorobenzene	1.455	1.421	2.3	97	-0.01
13	C	1,4-Dichlorobenzene	1.480	1.357	8.3	94	-0.01
14		1,2-Dichlorobenzene	1.285	1.317	-2.5	109	0.00
15		Benzyl Alcohol	1.136	1.026	9.7	93	0.00
16		2,2'-oxybis(1-Chloropropane	1.975	1.892	4.2	100	0.00
17		2-Methylphenol	1.096	1.096	0.0	105	0.00
18		Hexachloroethane	0.549	0.537	2.2	97	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	0.920	5.4	94	-0.01
20		3+4-Methylphenols	1.307	1.337	-2.3	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.493	0.483	2.0	102	0.00
23	S	Nitrobenzene-d5	0.360	0.349	3.1	106	0.00
24		Nitrobenzene	0.383	0.380	0.8	94	0.00
25		Isophorone	0.664	0.659	0.8	105	0.00
26	C	2-Nitrophenol	0.189	0.179	5.3	101	0.00
27		2,4-Dimethylphenol	0.313	0.305	2.6	94	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.374	7.0	95	-0.01
29	C	2,4-Dichlorophenol	0.265	0.272	-2.6	105	0.00
30		1,2,4-Trichlorobenzene	0.291	0.286	1.7	98	0.00
31		Naphthalene	0.915	0.902	1.4	94	-0.01
32		Benzoic acid	0.232	0.239	-3.0	102	-0.01
33		4-Chloroaniline	0.413	0.386	6.5	96	0.00
34	C	Hexachlorobutadiene	0.153	0.149	2.6	100	0.00
35		Caprolactam	0.087	0.083	4.6	97	0.00
36	C	4-Chloro-3-methylphenol	0.305	0.292	4.3	93	-0.01
37		2-Methylnaphthalene	0.628	0.600	4.5	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
39		1,2,4,5-Tetrachlorobenzene	0.505	0.454	10.1	90	0.00
40	P	Hexachlorocyclopentadiene	0.199	0.168	15.6	82	-0.01
41	S	2,4,6-Tribromophenol	0.166	0.157	5.4	109	0.00
42	C	2,4,6-Trichlorophenol	0.353	0.341	3.4	98	0.00
43		2,4,5-Trichlorophenol	0.364	0.341	6.3	101	0.00
44	S	2-Fluorobiphenyl	1.042	0.936	10.2	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122216\
Data File : BF091877.D
Acq On : 22 Dec 2016 11:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 23 00:56:27 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 16:17:51 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	1.369	1.395	-1.9	101	0.00
46		2-Chloronaphthalene	1.084	1.051	3.0	103	0.00
47		2-Nitroaniline	0.386	0.372	3.6	95	-0.01
48		Acenaphthylene	1.668	1.635	2.0	109	0.00
49		Dimethylphthalate	1.279	1.250	2.3	104	0.00
50		2,6-Dinitrotoluene	0.280	0.291	-3.9	115	0.00
51	C	Acenaphthene	1.131	1.035	8.5	103	0.00
52		3-Nitroaniline	0.346	0.375	-8.4	106	0.00
53	P	2,4-Dinitrophenol	0.156	0.136	12.8	85	0.00
54		Dibenzofuran	1.527	1.400	8.3	111	0.00
55	P	4-Nitrophenol	0.235	0.233	0.9	101	0.00
56		2,4-Dinitrotoluene	0.351	0.328	6.6	106	0.00
57		Fluorene	1.111	1.186	-6.8	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.288	0.270	6.2	97	0.00
59		Diethylphthalate	1.242	1.169	5.9	96	-0.01
60		4-Chlorophenyl-phenylether	0.486	0.458	5.8	105	0.00
61		4-Nitroaniline	0.359	0.371	-3.3	103	-0.01
62		Azobenzene	1.368	1.377	-0.7	113	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.132	-9.1	115	0.00
65	c	n-Nitrosodiphenylamine	0.593	0.618	-4.2	119	0.00
66		4-Bromophenyl-phenylether	0.208	0.208	0.0	115	0.00
67		Hexachlorobenzene	0.222	0.216	2.7	108	0.00
68		Atrazine	0.191	0.152	20.4	90	-0.01
69	C	Pentachlorophenol	0.109	0.111	-1.8	103	0.00
70		Phenanthrene	1.084	0.999	7.8	101	-0.01
71		Anthracene	1.086	0.983	9.5	115	0.00
72		Carbazole	1.005	0.970	3.5	120	0.00
73		Di-n-butylphthalate	1.174	1.000	14.8	104	0.00
74	C	Fluoranthene	1.082	0.927	14.3	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	114	0.00
76		Benzidine	0.580	0.461	20.5	95	0.00
77		Pyrene	1.467	1.357	7.5	113	0.00
78	S	Terphenyl-d14	0.825	0.682	17.3	103	0.00
79		Butylbenzylphthalate	0.667	0.679	-1.8	108	0.00
80		Benzo(a)anthracene	1.129	0.975	13.6	99	0.00
81		3,3'-Dichlorobenzidine	0.428	0.378	11.7	107	0.00
82		Chrysene	1.132	1.006	11.1	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.786	0.726	7.6	106	0.00
84	c	Di-n-octyl phthalate	1.456	1.360	6.6	107	0.00
85		Indeno(1,2,3-cd)pyrene	0.981	0.946	3.6	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	117	0.00
87		Benzo(b)fluoranthene	1.245	1.328	-6.7	117	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122216\
Data File : BF091877.D
Acq On : 22 Dec 2016 11:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 23 00:56:27 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 16:17:51 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	1.062	0.857	19.3	97	0.00
89 C	Benzo(a)pyrene	1.105	1.037	6.2	107	0.00
90	Dibenzo(a,h)anthracene	0.908	0.902	0.7	113	-0.01
91	Benzo(g,h,i)perylene	0.945	0.916	3.1	111	-0.01

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