Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100665.D
Acq On : 17 Nov 2017 9:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 17 15:30:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	51	0.00
2		1,4-Dioxane	0.608	0.594	2.3	50	0.00
3		Pyridine	1.659	1.562	5.8	47#	0.00
4		n-Nitrosodimethylamine	0.805	0.727	9.7	45#	0.00
5	S	2-Fluorophenol	1.248	1.265	-1.4	53	0.00
6		Aniline	2.174	2.075	4.6	49#	0.00
7	S	Phenol-d6	1.539	1.526	0.8	51	0.00
8		2-Chlorophenol	1.320	1.362	-3.2	53	0.00
9		Benzaldehyde	1.099	0.997	9.3	47#	0.00
10	C	Phenol	1.615	1.604	0.7	52	0.00
11		bis(2-Chloroethyl)ether	1.357	1.328	2.1	51	0.00
12		1,3-Dichlorobenzene	1.522	1.584	-4.1	54	0.00
13	C	1,4-Dichlorobenzene	1.540	1.617	-5.0	54	0.00
14		1,2-Dichlorobenzene	1.424	1.494	-4.9	54	0.00
15		Benzyl Alcohol	1.201	1.205	-0.3	50	0.00
16		2,2'-oxybis(1-Chloropropane	2.170	2.032	6.4	48#	0.00
17		2-Methylphenol	1.128	1.154	-2.3	53	0.00
18		Hexachloroethane	0.508	0.535	-5.3	53	0.00
19	P	n-Nitroso-di-n-propylamine	1.067	1.011	5.2	49#	0.00
20		3+4-Methylphenols	1.427	1.437	-0.7	51	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	50	0.00
22		Acetophenone	0.488	0.482	1.2	51	0.00
23	S	Nitrobenzene-d5	0.303	0.343	-13.2	55	0.00
24		Nitrobenzene	0.311	0.346	-11.3	52	0.00
25		Isophorone	0.636	0.600	5.7	48#	0.00
26	C	2-Nitrophenol	0.132	0.186	-40.9#	66	0.00
27		2,4-Dimethylphenol	0.285	0.289	-1.4	50	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.409	1.7	50#	0.00
29	C	2,4-Dichlorophenol	0.272	0.292	-7.4	52	0.00
30		1,2,4-Trichlorobenzene	0.294	0.311	-5.8	54	0.00
31		Naphthalene	0.963	0.988	-2.6	52	0.00
32		Benzoic acid	0.186	0.236	-26.9#	61	0.00
33		4-Chloroaniline	0.401	0.409	-2.0	51	0.00
34	C	Hexachlorobutadiene	0.175	0.183	-4.6	53	0.00
35		Caprolactam	0.093	0.090	3.2	48#	0.00
36	C	4-Chloro-3-methylphenol	0.292	0.293	-0.3	49#	0.00
37		2-Methylnaphthalene	0.640	0.655	-2.3	52	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	48#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.663	0.743	-12.1	54	0.00
40	P	Hexachlorocyclopentadiene	0.312	0.346	-10.9	50#	0.00
41	S	2,4,6-Tribromophenol	0.204	0.234	-14.7	53	0.00
42	C	2,4,6-Trichlorophenol	0.420	0.461	-9.8	50	0.00
43		2,4,5-Trichlorophenol	0.436	0.502	-15.1	54	0.00
44	S	2-Fluorobiphenyl	1.313	1.421	-8.2	54	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100665.D
Acq On : 17 Nov 2017 9:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 17 15:30:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	1.730	1.838	-6.2	53	0.00
46		2-Chloronaphthalene	1.255	1.371	-9.2	53	0.00
47		2-Nitroaniline	0.350	0.412	-17.7	53	0.00
48		Acenaphthylene	2.080	2.205	-6.0	52	0.00
49		Dimethylphthalate	1.527	1.586	-3.9	51	0.00
50		2,6-Dinitrotoluene	0.302	0.352	-16.6	54	0.00
51	C	Acenaphthene	1.265	1.363	-7.7	53	0.00
52		3-Nitroaniline	0.357	0.404	-13.2	53	0.00
53	P	2,4-Dinitrophenol	0.091	0.170	-86.8#	90	0.00
54		Dibenzofuran	1.773	1.846	-4.1	51	0.00
55	P	4-Nitrophenol	0.290	0.318	-9.7	51	0.00
56		2,4-Dinitrotoluene	0.381	0.464	-21.8	55	0.00
57		Fluorene	1.299	1.433	-10.3	53	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.382	-7.0	50#	0.00
59		Diethylphthalate	1.528	1.534	-0.4	49#	0.00
60		4-Chlorophenyl-phenylether	0.637	0.709	-11.3	54	0.00
61		4-Nitroaniline	0.338	0.396	-17.2	53	0.00
62		Azobenzene	1.439	1.375	4.4	47#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	49#	0.00
64		4,6-Dinitro-2-methylphenol	0.087	0.140	-60.9#	73	0.00
65	c	n-Nitrosodiphenylamine	0.695	0.705	-1.4	50	0.00
66		4-Bromophenyl-phenylether	0.214	0.226	-5.6	52	0.00
67		Hexachlorobenzene	0.224	0.235	-4.9	51	0.00
68		Atrazine	0.211	0.215	-1.9	49#	0.00
69	C	Pentachlorophenol	0.161	0.167	-3.7	50#	0.00
70		Phenanthrene	1.053	1.080	-2.6	52	0.00
71		Anthracene	1.068	1.101	-3.1	52	0.00
72		Carbazole	0.986	1.008	-2.2	52	0.00
73		Di-n-butylphthalate	1.154	1.217	-5.5	52	0.00
74	C	Fluoranthene	1.116	1.146	-2.7	52	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	50#	0.00
76		Benzidine	0.801	0.794	0.9	46#	0.00
77		Pyrene	1.426	1.459	-2.3	51	0.00
78	S	Terphenyl-d14	0.870	0.887	-2.0	53	0.00
79		Butylbenzylphthalate	0.581	0.628	-8.1	52	0.00
80		Benzo(a)anthracene	1.204	1.265	-5.1	52	0.00
81		3,3'-Dichlorobenzidine	0.432	0.448	-3.7	49#	0.00
82		Chrysene	1.166	1.168	-0.2	50	0.00
83		Bis(2-ethylhexyl)phthalate	0.765	0.861	-12.5	53	0.00
84	c	Di-n-octyl phthalate	1.314	1.373	-4.5	49#	0.00
85		Indeno(1,2,3-cd)pyrene	0.943	0.921	2.3	50#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	46#	0.00
87		Benzo(b)fluoranthene	1.185	1.206	-1.8	47#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100665.D
Acq On : 17 Nov 2017 9:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 17 15:30:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 17 15:22:08 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	1.120	1.240	-10.7	48#	0.00
89 C	Benzo(a)pyrene	1.050	1.104	-5.1	47#	0.00
90	Dibenzo(a,h)anthracene	0.859	0.919	-7.0	50#	0.00
91	Benzo(g,h,i)perylene	0.841	0.908	-8.0	51	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1