Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100706.D
Acq On : 18 Nov 2017 6:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 00:10:04 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	41#	0.00
2		1,4-Dioxane	0.608	0.586	3.6	39#	-0.02
3		Pyridine	1.659	1.618	2.5	39#	-0.01
4		n-Nitrosodimethylamine	0.805	0.711	11.7	35#	-0.02
5	S	2-Fluorophenol	1.248	1.273	-2.0	42#	0.00
6		Aniline	2.174	2.052	5.6	38#	0.00
7	S	Phenol-d6	1.539	1.533	0.4	41#	0.00
8		2-Chlorophenol	1.320	1.373	-4.0	43#	0.00
9		Benzaldehyde	1.099	1.008	8.3	38#	0.00
10	C	Phenol	1.615	1.564	3.2	40#	0.00
11		bis(2-Chloroethyl)ether	1.357	1.316	3.0	40#	0.00
12		1,3-Dichlorobenzene	1.522	1.614	-6.0	43#	0.00
13	C	1,4-Dichlorobenzene	1.540	1.637	-6.3	44#	0.00
14		1,2-Dichlorobenzene	1.424	1.522	-6.9	44#	0.00
15		Benzyl Alcohol	1.201	1.142	4.9	38#	0.00
16		2,2'-oxybis(1-Chloropropane	2.170	1.713	21.1	32#	0.00
17		2-Methylphenol	1.128	1.110	1.6	40#	0.00
18		Hexachloroethane	0.508	0.465	8.5	37#	0.00
19	P	n-Nitroso-di-n-propylamine	1.067	0.987	7.5	38#	0.00
20		3+4-Methylphenols	1.427	1.393	2.4	40#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	38#	0.00
22		Acetophenone	0.488	0.504	-3.3	40#	0.00
23	S	Nitrobenzene-d5	0.303	0.350	-15.5	42#	0.00
24		Nitrobenzene	0.311	0.350	-12.5	39#	0.00
25		Isophorone	0.636	0.599	5.8	35#	0.00
26	C	2-Nitrophenol	0.132	0.177	-34.1#	47#	0.00
27		2,4-Dimethylphenol	0.285	0.277	2.8	36#	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.420	-1.0	38#	0.00
29	C	2,4-Dichlorophenol	0.272	0.285	-4.8	38#	0.00
30		1,2,4-Trichlorobenzene	0.294	0.313	-6.5	40#	0.00
31		Naphthalene	0.963	0.992	-3.0	39#	0.00
32		Benzoic acid	0.186	0.159	14.5	31#	-0.03
33		4-Chloroaniline	0.401	0.400	0.2	37#	0.00
34	C	Hexachlorobutadiene	0.175	0.196	-12.0	42#	0.00
35		Caprolactam	0.093	0.076	18.3	31#	-0.01
36	C	4-Chloro-3-methylphenol	0.292	0.261	10.6	33#	0.00
37		2-Methylnaphthalene	0.640	0.633	1.1	38#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	29#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.663	0.865	-30.5#	38#	0.00
40	P	Hexachlorocyclopentadiene	0.312	0.066	78.8#	6#	0.00
41	S	2,4,6-Tribromophenol	0.204	0.200	2.0	28#	0.00
42	C	2,4,6-Trichlorophenol	0.420	0.496	-18.1	33#	0.00
43		2,4,5-Trichlorophenol	0.436	0.499	-14.4	32#	0.00
44	S	2-Fluorobiphenyl	1.313	1.705	-29.9#	39#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100706.D
Acq On : 18 Nov 2017 6:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 00:10:04 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	2.119	-22.5	37#	0.00
46		2-Chloronaphthalene	1.255	1.473	-17.4	34#	0.00
47		2-Nitroaniline	0.350	0.416	-18.9	33#	0.00
48		Acenaphthylene	2.080	2.259	-8.6	32#	0.00
49		Dimethylphthalate	1.527	1.769	-15.8	35#	0.00
50		2,6-Dinitrotoluene	0.302	0.349	-15.6	32#	0.00
51	C	Acenaphthene	1.265	1.323	-4.6	31#	0.00
52		3-Nitroaniline	0.357	0.384	-7.6	30#	0.00
53	P	2,4-Dinitrophenol	0.091	0.029#	68.1#	9#	0.00
54		Dibenzofuran	1.773	1.818	-2.5	30#	0.00
55	P	4-Nitrophenol	0.290	0.242	16.6	23#	0.00
56		2,4-Dinitrotoluene	0.381	0.402	-5.5	29#	0.00
57		Fluorene	1.299	1.360	-4.7	30#	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.337	5.6	26#	0.00
59		Diethylphthalate	1.528	1.602	-4.8	31#	0.00
60		4-Chlorophenyl-phenylether	0.637	0.709	-11.3	33#	0.00
61		4-Nitroaniline	0.338	0.340	-0.6	28#	0.00
62		Azobenzene	1.439	1.241	13.8	26#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	24#	0.00
64		4,6-Dinitro-2-methylphenol	0.087	0.045	48.3#	12#	0.00
65	c	n-Nitrosodiphenylamine	0.695	0.826	-18.8	29#	0.00
66		4-Bromophenyl-phenylether	0.214	0.251	-17.3	28#	0.00
67		Hexachlorobenzene	0.224	0.250	-11.6	27#	0.00
68		Atrazine	0.211	0.229	-8.5	26#	0.00
69	C	Pentachlorophenol	0.161	0.169	-5.0	25#	0.00
70		Phenanthrene	1.053	1.133	-7.6	27#	0.00
71		Anthracene	1.068	1.161	-8.7	27#	0.00
72		Carbazole	0.986	1.051	-6.6	26#	0.00
73		Di-n-butylphthalate	1.154	1.323	-14.6	28#	0.00
74	C	Fluoranthene	1.116	1.305	-16.9	29#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	33#	0.00
76		Benzidine	0.801	0.688	14.1	27#	0.00
77		Pyrene	1.426	1.281	10.2	30#	0.00
78	S	Terphenyl-d14	0.870	0.811	6.8	33#	0.00
79		Butylbenzylphthalate	0.581	0.569	2.1	32#	0.00
80		Benzo(a)anthracene	1.204	1.276	-6.0	35#	0.00
81		3,3'-Dichlorobenzidine	0.432	0.481	-11.3	35#	0.00
82		Chrysene	1.166	1.217	-4.4	35#	0.00
83		Bis(2-ethylhexyl)phthalate	0.765	0.842	-10.1	35#	0.00
84	c	Di-n-octyl phthalate	1.314	1.496	-13.9	36#	0.00
85		Indeno(1,2,3-cd)pyrene	0.943	0.818	13.3	30#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	31#	0.00
87		Benzo(b)fluoranthene	1.185	1.275	-7.6	34#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF111717\
Data File : BF100706.D
Acq On : 18 Nov 2017 6:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 20 00:10:04 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	1.282	-14.5	34#	0.00
89 C	Benzo(a)pyrene	1.050	1.120	-6.7	32#	0.00
90	Dibenzo(a,h)anthracene	0.859	0.854	0.6	31#	0.00
91	Benzo(g,h,i)perylene	0.841	0.771	8.3	29#	0.00

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