Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF013118\
Data File : BF102659.D
Acq On : 31 Jan 2018 17:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 18:45:20 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 18:38:00 2018
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	114	0.00
2		1,4-Dioxane	0.627	0.660	-5.3	118	0.00
3		Pyridine	1.638	1.711	-4.5	114	0.00
4		n-Nitrosodimethylamine	0.642	0.659	-2.6	113	0.00
5	S	2-Fluorophenol	1.319	1.380	-4.6	117	0.00
6		Aniline	2.103	2.085	0.9	112	0.00
7	S	Phenol-d6	1.590	1.620	-1.9	116	0.00
8		2-Chlorophenol	1.383	1.427	-3.2	115	0.00
9		Benzaldehyde	0.892	0.870	2.5	111	0.00
10	C	Phenol	1.736	1.712	1.4	111	0.00
11		bis(2-Chloroethyl)ether	1.320	1.390	-5.3	117	0.00
12		1,3-Dichlorobenzene	1.644	1.654	-0.6	114	0.00
13	C	1,4-Dichlorobenzene	1.647	1.660	-0.8	114	0.00
14		1,2-Dichlorobenzene	1.507	1.507	0.0	112	0.00
15		Benzyl Alcohol	1.122	1.084	3.4	108	0.00
16		2,2'-oxybis(1-Chloropropane	2.215	2.175	1.8	109	0.00
17		2-Methylphenol	1.201	1.211	-0.8	112	0.00
18		Hexachloroethane	0.574	0.591	-3.0	115	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	0.992	-2.0	115	0.00
20		3+4-Methylphenols	1.412	1.543	-9.3	119	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	118	0.00
22		Acetophenone	0.477	0.466	2.3	118	0.00
23	S	Nitrobenzene-d5	0.348	0.344	1.1	116	0.00
24		Nitrobenzene	0.356	0.357	-0.3	116	0.00
25		Isophorone	0.620	0.614	1.0	116	0.00
26	C	2-Nitrophenol	0.187	0.191	-2.1	115	0.00
27		2,4-Dimethylphenol	0.272	0.268	1.5	114	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.383	0.0	116	0.00
29	C	2,4-Dichlorophenol	0.277	0.276	0.4	116	0.00
30		1,2,4-Trichlorobenzene	0.297	0.281	5.4	112	0.00
31		Naphthalene	0.999	0.956	4.3	114	0.00
32		Benzoic acid	0.228	0.206	9.6	105	0.00
33		4-Chloroaniline	0.420	0.410	2.4	114	0.00
34	C	Hexachlorobutadiene	0.159	0.149	6.3	109	0.00
35		Caprolactam	0.092	0.089	3.3	112	0.00
36	C	4-Chloro-3-methylphenol	0.292	0.292	0.0	116	0.00
37		2-Methylnaphthalene	0.665	0.636	4.4	113	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
39		1,2,4,5-Tetrachlorobenzene	0.602	0.583	3.2	112	0.00
40	P	Hexachlorocyclopentadiene	0.269	0.254	5.6	102	0.00
41	S	2,4,6-Tribromophenol	0.198	0.168	15.2	99	0.00
42	C	2,4,6-Trichlorophenol	0.403	0.370	8.2	106	0.00
43		2,4,5-Trichlorophenol	0.454	0.426	6.2	107	0.00
44	S	2-Fluorobiphenyl	1.360	1.282	5.7	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF013118\
Data File : BF102659.D
Acq On : 31 Jan 2018 17:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 18:45:20 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 18:38:00 2018
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	1.788	1.707	4.5	112	0.00
46		2-Chloronaphthalene	1.390	1.329	4.4	111	0.00
47		2-Nitroaniline	0.433	0.443	-2.3	118	0.00
48		Acenaphthylene	2.166	2.009	7.2	108	0.00
49		Dimethylphthalate	1.653	1.558	5.7	111	0.00
50		2,6-Dinitrotoluene	0.356	0.342	3.9	108	0.00
51	C	Acenaphthene	1.265	1.197	5.4	112	0.00
52		3-Nitroaniline	0.422	0.422	0.0	115	0.00
53	P	2,4-Dinitrophenol	0.145	0.161	-11.0	119	0.00
54		Dibenzofuran	1.712	1.635	4.5	108	0.00
55	P	4-Nitrophenol	0.278	0.246	11.5	96	0.00
56		2,4-Dinitrotoluene	0.438	0.403	8.0	105	0.00
57		Fluorene	1.356	1.346	0.7	115	0.00
58		2,3,4,6-Tetrachlorophenol	0.324	0.293	9.6	106	0.00
59		Diethylphthalate	1.606	1.489	7.3	109	0.00
60		4-Chlorophenyl-phenylether	0.588	0.584	0.7	113	0.00
61		4-Nitroaniline	0.421	0.407	3.3	111	0.00
62		Azobenzene	1.471	1.420	3.5	112	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
64		4,6-Dinitro-2-methylphenol	0.133	0.120	9.8	97	0.00
65	c	n-Nitrosodiphenylamine	0.734	0.713	2.9	113	0.00
66		4-Bromophenyl-phenylether	0.215	0.210	2.3	111	0.00
67		Hexachlorobenzene	0.241	0.215	10.8	102	0.00
68		Atrazine	0.217	0.203	6.5	108	0.00
69	C	Pentachlorophenol	0.126	0.109	13.5	91	0.00
70		Phenanthrene	1.165	1.101	5.5	110	0.00
71		Anthracene	1.190	1.140	4.2	111	0.00
72		Carbazole	1.161	1.122	3.4	112	0.00
73		Di-n-butylphthalate	1.451	1.357	6.5	107	0.00
74	C	Fluoranthene	1.188	1.104	7.1	108	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.672	0.656	2.4	105	0.00
77		Pyrene	1.546	1.634	-5.7	106	0.00
78	S	Terphenyl-d14	0.887	0.890	-0.3	103	0.00
79		Butylbenzylphthalate	0.770	0.795	-3.2	101	0.00
80		Benzo(a)anthracene	1.231	1.268	-3.0	105	0.00
81		3,3'-Dichlorobenzidine	0.452	0.430	4.9	94	0.00
82		Chrysene	1.250	1.239	0.9	100	0.00
83		Bis(2-ethylhexyl)phthalate	1.034	1.039	-0.5	99	0.00
84	c	Di-n-octyl phthalate	1.682	1.542	8.3	91	0.00
85		Indeno(1,2,3-cd)pyrene	1.033	0.901	12.8	94	0.00
86	I	Perylene-d12	1.000	1.000	0.0	92	0.00
87		Benzo(b)fluoranthene	1.296	1.243	4.1	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF013118\
Data File : BF102659.D
Acq On : 31 Jan 2018 17:16
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 31 18:45:20 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 31 18:38:00 2018
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	1.225	1.278	-4.3	99	0.00
89 C	Benzo(a)pyrene	1.169	1.159	0.9	91	0.00
90	Dibenzo(a,h)anthracene	0.947	0.904	4.5	92	0.00
91	Benzo(g,h,i)perylene	0.935	0.952	-1.8	99	0.00

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