Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032818\
Data File : BF104008.D
Acq On : 28 Mar 2018 14:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 16:55:01 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 28 13:41:47 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	112	0.00
2		1,4-Dioxane	0.541	0.518	4.3	107	0.00
3		Pyridine	1.369	1.356	0.9	107	0.00
4		n-Nitrosodimethylamine	0.407	0.393	3.4	110	0.00
5	S	2-Fluorophenol	1.254	1.216	3.0	106	0.00
6		Aniline	1.847	1.842	0.3	105	0.00
7	S	Phenol-d6	1.543	1.453	5.8	105	0.00
8		2-Chlorophenol	1.376	1.336	2.9	106	0.00
9		Benzaldehyde	0.819	0.830	-1.3	114	0.00
10	C	Phenol	1.710	1.586	7.3	105	0.00
11		bis(2-Chloroethyl)ether	1.473	1.471	0.1	111	0.00
12		1,3-Dichlorobenzene	1.546	1.492	3.5	107	0.00
13	C	1,4-Dichlorobenzene	1.657	1.592	3.9	107	0.00
14		1,2-Dichlorobenzene	1.489	1.430	4.0	107	0.00
15		Benzyl Alcohol	1.003	0.994	0.9	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.605	1.542	3.9	107	0.00
17		2-Methylphenol	1.120	1.098	2.0	107	0.00
18		Hexachloroethane	0.555	0.537	3.2	108	0.00
19	P	n-Nitroso-di-n-propylamine	0.916	0.890	2.8	107	0.00
20		3+4-Methylphenols	1.429	1.441	-0.8	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.484	0.458	5.4	105	0.00
23	S	Nitrobenzene-d5	0.327	0.321	1.8	108	0.00
24		Nitrobenzene	0.344	0.338	1.7	106	0.00
25		Isophorone	0.603	0.568	5.8	107	0.00
26	C	2-Nitrophenol	0.180	0.174	3.3	103	0.00
27		2,4-Dimethylphenol	0.259	0.245	5.4	106	0.00
28		bis(2-Chloroethoxy)methane	0.378	0.363	4.0	107	0.00
29	C	2,4-Dichlorophenol	0.271	0.265	2.2	107	0.00
30		1,2,4-Trichlorobenzene	0.307	0.299	2.6	108	0.00
31		Naphthalene	0.977	0.933	4.5	106	0.00
32		Benzoic acid	0.190	0.194	-2.1	115	0.00
33		4-Chloroaniline	0.412	0.408	1.0	107	0.00
34	C	Hexachlorobutadiene	0.167	0.161	3.6	107	0.00
35		Caprolactam	0.086	0.082	4.7	104	0.00
36	C	4-Chloro-3-methylphenol	0.286	0.281	1.7	106	0.00
37		2-Methylnaphthalene	0.644	0.616	4.3	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.613	0.590	3.8	106	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.265	-4.7	112	0.00
41	S	2,4,6-Tribromophenol	0.223	0.216	3.1	106	0.00
42	C	2,4,6-Trichlorophenol	0.388	0.382	1.5	107	0.00
43		2,4,5-Trichlorophenol	0.469	0.453	3.4	106	0.00
44	S	2-Fluorobiphenyl	1.268	1.233	2.8	107	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032818\
Data File : BF104008.D
Acq On : 28 Mar 2018 14:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 16:55:01 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 28 13:41:47 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	1.717	1.645	4.2	107	0.00
46		2-Chloronaphthalene	1.354	1.301	3.9	107	0.00
47		2-Nitroaniline	0.386	0.370	4.1	104	0.00
48		Acenaphthylene	2.051	1.934	5.7	105	0.00
49		Dimethylphthalate	1.579	1.494	5.4	105	0.00
50		2,6-Dinitrotoluene	0.340	0.331	2.6	107	0.00
51	C	Acenaphthene	1.187	1.111	6.4	106	0.00
52		3-Nitroaniline	0.391	0.372	4.9	103	0.00
53	P	2,4-Dinitrophenol	0.128	0.135	-5.5	117	0.00
54		Dibenzofuran	1.733	1.626	6.2	104	0.00
55	P	4-Nitrophenol	0.234	0.226	3.4	102	0.00
56		2,4-Dinitrotoluene	0.433	0.424	2.1	104	0.00
57		Fluorene	1.429	1.332	6.8	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.351	0.338	3.7	107	0.00
59		Diethylphthalate	1.513	1.436	5.1	105	0.00
60		4-Chlorophenyl-phenylether	0.657	0.620	5.6	105	0.00
61		4-Nitroaniline	0.365	0.343	6.0	102	0.00
62		Azobenzene	1.368	1.311	4.2	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.124	-2.5	107	0.00
65	c	n-Nitrosodiphenylamine	0.677	0.639	5.6	104	0.00
66		4-Bromophenyl-phenylether	0.214	0.206	3.7	106	0.00
67		Hexachlorobenzene	0.246	0.233	5.3	104	0.00
68		Atrazine	0.208	0.198	4.8	104	0.00
69	C	Pentachlorophenol	0.135	0.135	0.0	105	0.00
70		Phenanthrene	1.083	1.007	7.0	103	0.00
71		Anthracene	1.131	1.061	6.2	104	0.00
72		Carbazole	1.080	1.015	6.0	104	0.00
73		Di-n-butylphthalate	1.286	1.203	6.5	103	0.00
74	C	Fluoranthene	1.151	1.066	7.4	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
76		Benzidine	0.570	0.601	-5.4	105	0.00
77		Pyrene	1.438	1.386	3.6	103	0.00
78	S	Terphenyl-d14	0.866	0.806	6.9	102	0.00
79		Butylbenzylphthalate	0.677	0.650	4.0	101	0.00
80		Benzo(a)anthracene	1.251	1.194	4.6	102	0.00
81		3,3'-Dichlorobenzidine	0.414	0.387	6.5	100	0.00
82		Chrysene	1.226	1.156	5.7	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.875	0.824	5.8	102	0.00
84	c	Di-n-octyl phthalate	1.505	1.433	4.8	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.270	1.128	11.2	93	0.00
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Benzo(b)fluoranthene	1.217	1.102	9.4	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF032818\
Data File : BF104008.D
Acq On : 28 Mar 2018 14:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 28 16:55:01 2018
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 28 13:41:47 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	1.147	1.136	1.0	101	0.00
89 C	Benzo(a)pyrene	1.128	1.064	5.7	99	0.00
90	Dibenzo(a,h)anthracene	1.043	0.928	11.0	93	0.00
91	Benzo(g,h,i)perylene	1.045	0.912	12.7	91	0.00

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