Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120815\
Data File : BF083593.D
Acq On : 8 Dec 2015 13:49
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Dec 09 01:10:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 00:23:35 2015
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	97	0.00
2		1,4-Dioxane	0.651	0.683	-4.9	106	0.00
3		Pyridine	1.583	1.715	-8.3	106	0.00
4		n-Nitrosodimethylamine	0.833	0.910	-9.2	98	0.01
5	S	2-Fluorophenol	1.319	1.450	-9.9	101	0.00
6		Aniline	2.171	2.277	-4.9	94	0.01
7	S	Phenol-d6	1.762	1.889	-7.2	100	0.00
8		2-Chlorophenol	1.442	1.535	-6.4	100	0.00
9		Benzaldehyde	0.936	0.974	-4.1	97	0.00
10	C	Phenol	2.134	2.287	-7.2	98	0.00
11		bis(2-Chloroethyl)ether	1.570	1.591	-1.3	95	0.01
12		1,3-Dichlorobenzene	1.607	1.749	-8.8	102	0.11
13	C	1,4-Dichlorobenzene	1.649	1.753	-6.3	102	0.00
14		1,2-Dichlorobenzene	1.530	1.622	-6.0	100	0.00
15		Benzyl Alcohol	1.254	1.412	-12.6	101	0.00
16		2,2'-oxybis(1-Chloropropane	2.903	3.083	-6.2	100	0.00
17		2-Methylphenol	1.173	1.245	-6.1	98	0.00
18		Hexachloroethane	0.582	0.624	-7.2	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.230	1.282	-4.2	100	0.00
20		3+4-Methylphenols	1.545	1.605	-3.9	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	-0.01
22		Acetophenone	0.561	0.603	-7.5	100	0.00
23	S	Nitrobenzene-d5	0.429	0.455	-6.1	101	0.00
24		Nitrobenzene	0.474	0.496	-4.6	102	0.00
25		Isophorone	0.832	0.873	-4.9	100	0.01
26	C	2-Nitrophenol	0.188	0.207	-10.1	100	0.00
27		2,4-Dimethylphenol	0.329	0.349	-6.1	100	0.00
28		bis(2-Chloroethoxy)methane	0.462	0.484	-4.8	99	0.00
29	C	2,4-Dichlorophenol	0.299	0.315	-5.4	98	0.00
30		1,2,4-Trichlorobenzene	0.331	0.344	-3.9	99	0.00
31		Naphthalene	1.059	1.096	-3.5	100	0.00
32		Benzoic acid	0.185	0.216	-16.8	104	0.01
33		4-Chloroaniline	0.386	0.406	-5.2	97	0.00
34	C	Hexachlorobutadiene	0.193	0.203	-5.2	100	0.00
35		Caprolactam	0.090	0.098	-8.9	99	0.00
36	C	4-Chloro-3-methylphenol	0.336	0.359	-6.8	100	0.01
37		2-Methylnaphthalene	0.665	0.691	-3.9	100	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
39		1,2,4,5-Tetrachlorobenzene	0.657	0.704	-7.2	100	0.00
40	P	Hexachlorocyclopentadiene	0.089	0.087	2.2	94	0.00
41	S	2,4,6-Tribromophenol	0.178	0.194	-9.0	100	0.00
42	C	2,4,6-Trichlorophenol	0.391	0.426	-9.0	99	0.00
43		2,4,5-Trichlorophenol	0.387	0.427	-10.3	99	0.01
44	S	2-Fluorobiphenyl	1.421	1.464	-3.0	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120815\
Data File : BF083593.D
Acq On : 8 Dec 2015 13:49
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Dec 09 01:10:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 00:23:35 2015
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	1.738	1.833	-5.5	99	0.00
46		2-Chloronaphthalene	1.312	1.407	-7.2	100	0.00
47		2-Nitroaniline	0.473	0.509	-7.6	99	0.00
48		Acenaphthylene	2.164	2.310	-6.7	101	0.00
49		Dimethylphthalate	1.592	1.685	-5.8	100	0.00
50		2,6-Dinitrotoluene	0.329	0.360	-9.4	101	0.00
51	C	Acenaphthene	1.239	1.307	-5.5	100	0.00
52		3-Nitroaniline	0.349	0.393	-12.6	100	0.00
53	P	2,4-Dinitrophenol	0.140	0.145	-3.6	92	-0.01
54		Dibenzofuran	1.718	1.819	-5.9	99	0.00
55	P	4-Nitrophenol	0.154	0.164	-6.5	99	0.01
56		2,4-Dinitrotoluene	0.436	0.494	-13.3	100	0.00
57		Fluorene	1.500	1.602	-6.8	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.286	0.320	-11.9	96	0.00
59		Diethylphthalate	1.548	1.652	-6.7	102	0.00
60		4-Chlorophenyl-phenylether	0.714	0.783	-9.7	102	0.00
61		4-Nitroaniline	0.317	0.339	-6.9	94	0.00
62		Azobenzene	1.704	1.857	-9.0	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
64		4,6-Dinitro-2-methylphenol	0.120	0.139	-15.8	97	0.00
65	c	n-Nitrosodiphenylamine	0.660	0.700	-6.1	100	0.00
66		4-Bromophenyl-phenylether	0.223	0.237	-6.3	100	0.00
67		Hexachlorobenzene	0.237	0.251	-5.9	101	0.00
68		Atrazine	0.196	0.210	-7.1	97	0.00
69	C	Pentachlorophenol	0.052	0.051	1.9	91	0.00
70		Phenanthrene	1.149	1.212	-5.5	101	0.00
71		Anthracene	1.190	1.261	-6.0	101	0.00
72		Carbazole	1.021	1.099	-7.6	103	0.00
73		Di-n-butylphthalate	1.284	1.382	-7.6	100	0.00
74	C	Fluoranthene	1.176	1.248	-6.1	101	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.622	0.641	-3.1	94	0.01
77		Pyrene	1.440	1.497	-4.0	101	0.00
78	S	Terphenyl-d14	0.850	0.857	-0.8	101	0.00
79		Butylbenzylphthalate	0.639	0.695	-8.8	103	0.01
80		Benzo(a)anthracene	1.171	1.198	-2.3	101	0.00
81		3,3'-Dichlorobenzidine	0.395	0.425	-7.6	101	0.00
82		Chrysene	1.174	1.226	-4.4	102	0.00
83		Bis(2-ethylhexyl)phthalate	0.887	0.939	-5.9	99	0.00
84	c	Di-n-octyl phthalate	1.355	1.461	-7.8	103	0.00
85		Indeno(1,2,3-cd)pyrene	0.792	0.795	-0.4	104	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Benzo(b)fluoranthene	1.139	2.550	-123.9#	225#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120815\
Data File : BF083593.D
Acq On : 8 Dec 2015 13:49
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Dec 09 01:10:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 00:23:35 2015
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	1.255	2.554	-103.5#	195#	-0.03
89 C	Benzo(a)pyrene	1.111	1.151	-3.6	104	0.00
90	Dibenzo(a,h)anthracene	0.850	0.862	-1.4	105	0.00
91	Benzo(g,h,i)perylene	0.828	0.818	1.2	103	0.00

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