Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011215\
Data File : BF076808.D
Acq On : 12 Jan 2015 10:41
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 09:47:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jan 10 06:22:22 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	106	0.00
2		1,4-Dioxane	0.511	0.491	3.9	95	0.00
3		Pyridine	1.360	1.460	-7.4	127	-0.01
4		n-Nitrosodimethylamine	0.684	0.770	-12.6	109	0.00
5	S	2-Fluorophenol	1.199	1.277	-6.5	105	-0.01
6		Aniline	2.083	2.215	-6.3	104	0.00
7	S	Phenol-d6	1.509	1.576	-4.4	104	-0.01
8		2-Chlorophenol	1.417	1.420	-0.2	102	0.00
9		Benzaldehyde	0.896	0.755	15.7	102	0.00
10	C	Phenol	1.644	1.690	-2.8	102	-0.01
11		bis(2-Chloroethyl)ether	1.317	1.307	0.8	101	0.00
12		1,3-Dichlorobenzene	1.574	1.676	-6.5	109	0.00
13	C	1,4-Dichlorobenzene	1.650	1.734	-5.1	106	0.00
14		1,2-Dichlorobenzene	1.538	1.622	-5.5	106	0.00
15		Benzyl Alcohol	1.255	1.288	-2.6	100	0.00
16		2,2'-oxybis(1-Chloropropane	1.770	1.803	-1.9	105	0.00
17		2-Methylphenol	1.135	1.161	-2.3	103	-0.01
18		Hexachloroethane	0.597	0.623	-4.4	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.072	1.110	-3.5	104	0.00
20		3+4-Methylphenols	1.499	1.553	-3.6	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.491	0.530	-7.9	102	0.00
23	S	Nitrobenzene-d5	0.360	0.394	-9.4	103	0.00
24		Nitrobenzene	0.364	0.410	-12.6	104	0.00
25		Isophorone	0.655	0.702	-7.2	101	0.00
26	C	2-Nitrophenol	0.174	0.191	-9.8	103	0.00
27		2,4-Dimethylphenol	0.282	0.309	-9.6	103	0.00
28		bis(2-Chloroethoxy)methane	0.384	0.414	-7.8	104	0.00
29	C	2,4-Dichlorophenol	0.269	0.300	-11.5	100	0.00
30		1,2,4-Trichlorobenzene	0.303	0.330	-8.9	103	0.00
31		Naphthalene	1.026	1.117	-8.9	104	0.00
32		Benzoic acid	0.138	0.144	-4.3	97	0.00
33		4-Chloroaniline	0.402	0.443	-10.2	104	0.00
34	C	Hexachlorobutadiene	0.173	0.187	-8.1	102	0.00
35		Caprolactam	0.080	0.084	-5.0	100	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.328	-10.1	101	-0.01
37		2-Methylnaphthalene	0.720	0.764	-6.1	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	0.497	0.514	-3.4	97	0.00
40	P	Hexachlorocyclopentadiene	0.247	0.261	-5.7	100	0.00
41	S	2,4,6-Tribromophenol	0.167	0.177	-6.0	98	-0.01
42	C	2,4,6-Trichlorophenol	0.361	0.401	-11.1	101	0.00
43		2,4,5-Trichlorophenol	0.379	0.422	-11.3	101	-0.01
44	S	2-Fluorobiphenyl	1.359	1.429	-5.2	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011215\
Data File : BF076808.D
Acq On : 12 Jan 2015 10:41
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 09:47:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jan 10 06:22:22 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	1.538	1.632	-6.1	100	0.00
46		2-Chloronaphthalene	1.236	1.331	-7.7	99	0.00
47		2-Nitroaniline	0.368	0.425	-15.5	106	-0.01
48		Acenaphthylene	2.094	2.217	-5.9	100	0.00
49		Dimethylphthalate	1.454	1.530	-5.2	97	0.00
50		2,6-Dinitrotoluene	0.308	0.310	-0.6	96	0.00
51	C	Acenaphthene	1.183	1.234	-4.3	97	0.00
52		3-Nitroaniline	0.352	0.380	-8.0	104	0.00
53	P	2,4-Dinitrophenol	0.075	0.076	-1.3	102	0.00
54		Dibenzofuran	1.731	1.856	-7.2	102	0.00
55	P	4-Nitrophenol	0.227	0.246	-8.4	105	-0.01
56		2,4-Dinitrotoluene	0.394	0.433	-9.9	98	0.00
57		Fluorene	1.445	1.574	-8.9	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.274	0.298	-8.8	104	0.00
59		Diethylphthalate	1.497	1.564	-4.5	95	0.00
60		4-Chlorophenyl-phenylether	0.620	0.634	-2.3	97	-0.01
61		4-Nitroaniline	0.326	0.377	-15.6	102	-0.01
62		Azobenzene	1.535	1.609	-4.8	97	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.113	0.120	-6.2	99	0.00
65	c	n-Nitrosodiphenylamine	0.726	0.748	-3.0	97	0.00
66		4-Bromophenyl-phenylether	0.208	0.218	-4.8	96	0.00
67		Hexachlorobenzene	0.217	0.226	-4.1	99	-0.01
68		Atrazine	0.220	0.228	-3.6	96	0.00
69	C	Pentachlorophenol	0.073	0.078	-6.8	99	0.00
70		Phenanthrene	1.246	1.316	-5.6	101	0.00
71		Anthracene	1.211	1.296	-7.0	100	0.00
72		Carbazole	1.187	1.240	-4.5	98	-0.01
73		Di-n-butylphthalate	1.410	1.433	-1.6	95	0.00
74	C	Fluoranthene	1.277	1.290	-1.0	98	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.626	0.646	-3.2	116	0.00
77		Pyrene	1.439	1.450	-0.8	100	0.00
78	S	Terphenyl-d14	0.911	0.891	2.2	98	0.00
79		Butylbenzylphthalate	0.668	0.641	4.0	92	0.00
80		Benzo(a)anthracene	1.274	1.375	-7.9	107	0.00
81		3,3'-Dichlorobenzidine	0.407	0.433	-6.4	107	0.00
82		Chrysene	1.153	1.155	-0.2	98	0.01
83		Bis(2-ethylhexyl)phthalate	0.926	0.909	1.8	99	0.01
84	c	Di-n-octyl phthalate	1.562	1.587	-1.6	100	0.02
85		Indeno(1,2,3-cd)pyrene	1.141	1.275	-11.7	109	0.10
86	I	Perylene-d12	1.000	1.000	0.0	110	0.06
87		Benzo(b)fluoranthene	1.385	1.390	-0.4	95	0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011215\
Data File : BF076808.D
Acq On : 12 Jan 2015 10:41
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 13 09:47:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jan 10 06:22:22 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	1.211	1.278	-5.5	117	0.03
89 C	Benzo(a)pyrene	1.190	1.269	-6.6	107	0.06
90	Dibenzo(a,h)anthracene	1.085	1.128	-4.0	106	0.09
91	Benzo(g,h,i)perylene	1.107	1.180	-6.6	110	0.10

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