Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060915\
Data File : BF079834.D
Acq On : 9 Jun 2015 16:17
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 10 01:06:14 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 15:11:40 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	101	0.00
2		1,4-Dioxane	0.555	0.535	3.6	97	-0.01
3		Pyridine	1.490	1.579	-6.0	100	-0.03
4		n-Nitrosodimethylamine	0.746	0.698	6.4	91	-0.01
5	S	2-Fluorophenol	1.214	1.219	-0.4	99	0.00
6		Aniline	2.262	2.283	-0.9	98	0.00
7	S	Phenol-d6	1.570	1.614	-2.8	99	0.00
8		2-Chlorophenol	1.368	1.370	-0.1	103	0.00
9		Benzaldehyde	0.921	0.942	-2.3	98	0.00
10	C	Phenol	1.673	1.770	-5.8	103	0.00
11		bis(2-Chloroethyl)ether	1.306	1.328	-1.7	106	0.00
12		1,3-Dichlorobenzene	1.568	1.567	0.1	100	0.00
13	C	1,4-Dichlorobenzene	1.777	1.722	3.1	97	0.00
14		1,2-Dichlorobenzene	1.543	1.494	3.2	95	0.00
15		Benzyl Alcohol	1.266	1.269	-0.2	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.025	1.987	1.9	99	0.00
17		2-Methylphenol	1.177	1.206	-2.5	98	0.00
18		Hexachloroethane	0.552	0.561	-1.6	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.094	1.051	3.9	102	0.00
20		3+4-Methylphenols	1.521	1.546	-1.6	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.496	0.512	-3.2	101	0.00
23	S	Nitrobenzene-d5	0.346	0.374	-8.1	99	0.00
24		Nitrobenzene	0.348	0.349	-0.3	100	0.00
25		Isophorone	0.722	0.721	0.1	102	0.00
26	C	2-Nitrophenol	0.139	0.134	3.6	92	0.00
27		2,4-Dimethylphenol	0.326	0.330	-1.2	103	0.00
28		bis(2-Chloroethoxy)methane	0.437	0.421	3.7	94	0.00
29	C	2,4-Dichlorophenol	0.278	0.274	1.4	96	0.00
30		1,2,4-Trichlorobenzene	0.357	0.352	1.4	96	0.00
31		Naphthalene	1.065	1.117	-4.9	108	0.00
32		Benzoic acid	0.098	0.104	-6.1	99	0.02
33		4-Chloroaniline	0.430	0.426	0.9	98	0.00
34	C	Hexachlorobutadiene	0.196	0.198	-1.0	96	0.00
35		Caprolactam	0.081	0.084	-3.7	100	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.315	-4.7	99	0.00
37		2-Methylnaphthalene	0.757	0.769	-1.6	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	0.699	0.693	0.9	93	0.00
40	P	Hexachlorocyclopentadiene	0.334	0.336	-0.6	99	0.00
41	S	2,4,6-Tribromophenol	0.173	0.189	-9.2	108	0.00
42	C	2,4,6-Trichlorophenol	0.422	0.422	0.0	94	0.00
43		2,4,5-Trichlorophenol	0.436	0.478	-9.6	107	0.00
44	S	2-Fluorobiphenyl	1.422	1.414	0.6	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060915\
Data File : BF079834.D
Acq On : 9 Jun 2015 16:17
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 10 01:06:14 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 15:11:40 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	1.658	1.636	1.3	95	0.00
46		2-Chloronaphthalene	1.354	1.370	-1.2	100	0.00
47		2-Nitroaniline	0.297	0.313	-5.4	106	0.00
48		Acenaphthylene	2.302	2.354	-2.3	102	0.00
49		Dimethylphthalate	1.559	1.523	2.3	100	0.00
50		2,6-Dinitrotoluene	0.270	0.288	-6.7	106	0.00
51	C	Acenaphthene	1.347	1.326	1.6	96	0.00
52		3-Nitroaniline	0.333	0.362	-8.7	108	0.00
53	P	2,4-Dinitrophenol	0.064	0.065	-1.6	94	0.00
54		Dibenzofuran	1.886	1.904	-1.0	98	0.00
55	P	4-Nitrophenol	0.242	0.263	-8.7	107	0.00
56		2,4-Dinitrotoluene	0.347	0.348	-0.3	92	0.00
57		Fluorene	1.670	1.671	-0.1	97	0.00
58		2,3,4,6-Tetrachlorophenol	0.322	0.358	-11.2	105	0.00
59		Diethylphthalate	1.527	1.619	-6.0	104	0.00
60		4-Chlorophenyl-phenylether	0.723	0.748	-3.5	104	0.00
61		4-Nitroaniline	0.326	0.364	-11.7	105	0.00
62		Azobenzene	1.503	1.572	-4.6	102	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	-0.01
64		4,6-Dinitro-2-methylphenol	0.057	0.060	-5.3	104	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.679	-2.3	108	0.00
66		4-Bromophenyl-phenylether	0.231	0.221	4.3	102	0.00
67		Hexachlorobenzene	0.243	0.256	-5.3	102	0.00
68		Atrazine	0.188	0.182	3.2	95	-0.01
69	C	Pentachlorophenol	0.089	0.095	-6.7	107	-0.01
70		Phenanthrene	1.195	1.138	4.8	98	-0.01
71		Anthracene	1.170	1.204	-2.9	98	-0.01
72		Carbazole	1.083	1.134	-4.7	106	-0.01
73		Di-n-butylphthalate	1.124	1.187	-5.6	101	-0.02
74	C	Fluoranthene	1.253	1.234	1.5	96	-0.05
75	I	Chrysene-d12	1.000	1.000	0.0	99	-0.13
76		Benzidine	0.596	0.614	-3.0	95	-0.06
77		Pyrene	1.246	1.253	-0.6	98	-0.06
78	S	Terphenyl-d14	0.887	0.863	2.7	96	-0.07
79		Butylbenzylphthalate	0.417	0.422	-1.2	99	-0.09
80		Benzo(a)anthracene	1.189	1.148	3.4	94	-0.11
81		3,3'-Dichlorobenzidine	0.371	0.382	-3.0	98	-0.11
82		Chrysene	1.116	1.133	-1.5	98	-0.11
83		Bis(2-ethylhexyl)phthalate	0.645	0.659	-2.2	101	-0.13
84	c	Di-n-octyl phthalate	1.002	1.016	-1.4	100	-0.17
85		Indeno(1,2,3-cd)pyrene	1.178	1.206	-2.4	97	-0.18
86	I	Perylene-d12	1.000	1.000	0.0	96	-0.17
87		Benzo(b)fluoranthene	1.190	1.160	2.5	94	-0.16

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060915\
Data File : BF079834.D
Acq On : 9 Jun 2015 16:17
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 10 01:06:14 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 15:11:40 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	1.245	1.398	-12.3	100	-0.17
89 C	Benzo(a)pyrene	1.125	1.186	-5.4	97	-0.16
90	Dibenzo(a,h)anthracene	1.042	1.094	-5.0	94	-0.18
91	Benzo(g,h,i)perylene	1.100	1.145	-4.1	96	-0.18

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