Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061015\
Data File : BF079855.D
Acq On : 10 Jun 2015 10:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 11 14:01:36 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	112	0.00
2		1,4-Dioxane	0.555	0.537	3.2	108	-0.01
3		Pyridine	1.490	1.582	-6.2	111	-0.03
4		n-Nitrosodimethylamine	0.746	0.696	6.7	101	-0.01
5	S	2-Fluorophenol	1.214	1.221	-0.6	109	0.00
6		Aniline	2.262	2.263	-0.0	108	0.00
7	S	Phenol-d6	1.570	1.638	-4.3	111	0.00
8		2-Chlorophenol	1.368	1.364	0.3	114	0.00
9		Benzaldehyde	0.921	0.957	-3.9	111	0.00
10	C	Phenol	1.673	1.727	-3.2	111	0.00
11		bis(2-Chloroethyl)ether	1.306	1.292	1.1	114	0.00
12		1,3-Dichlorobenzene	1.568	1.568	0.0	111	0.00
13	C	1,4-Dichlorobenzene	1.777	1.739	2.1	109	0.00
14		1,2-Dichlorobenzene	1.543	1.495	3.1	105	0.00
15		Benzyl Alcohol	1.266	1.315	-3.9	107	0.00
16		2,2'-oxybis(1-Chloropropane	2.025	2.033	-0.4	113	0.00
17		2-Methylphenol	1.177	1.207	-2.5	108	0.00
18		Hexachloroethane	0.552	0.546	1.1	113	0.00
19	P	n-Nitroso-di-n-propylamine	1.094	1.044	4.6	112	0.00
20		3+4-Methylphenols	1.521	1.524	-0.2	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
22		Acetophenone	0.496	0.523	-5.4	115	0.00
23	S	Nitrobenzene-d5	0.346	0.362	-4.6	106	0.00
24		Nitrobenzene	0.348	0.334	4.0	106	0.00
25		Isophorone	0.722	0.708	1.9	111	0.00
26	C	2-Nitrophenol	0.139	0.130	6.5	99	0.00
27		2,4-Dimethylphenol	0.326	0.322	1.2	112	0.00
28		bis(2-Chloroethoxy)methane	0.437	0.436	0.2	108	0.00
29	C	2,4-Dichlorophenol	0.278	0.284	-2.2	110	0.01
30		1,2,4-Trichlorobenzene	0.357	0.366	-2.5	111	0.00
31		Naphthalene	1.065	1.052	1.2	112	0.00
32		Benzoic acid	0.098	0.104	-6.1	110	0.03
33		4-Chloroaniline	0.430	0.426	0.9	109	0.00
34	C	Hexachlorobutadiene	0.196	0.196	0.0	106	0.00
35		Caprolactam	0.081	0.084	-3.7	111	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.322	-7.0	112	0.00
37		2-Methylnaphthalene	0.757	0.769	-1.6	115	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.01
39		1,2,4,5-Tetrachlorobenzene	0.699	0.697	0.3	108	0.00
40	P	Hexachlorocyclopentadiene	0.334	0.293	12.3	100	0.00
41	S	2,4,6-Tribromophenol	0.173	0.176	-1.7	117	0.00
42	C	2,4,6-Trichlorophenol	0.422	0.436	-3.3	112	0.00
43		2,4,5-Trichlorophenol	0.436	0.442	-1.4	114	0.00
44	S	2-Fluorobiphenyl	1.422	1.354	4.8	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061015\
Data File : BF079855.D
Acq On : 10 Jun 2015 10:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 11 14:01:36 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	1.614	2.7	108	0.01
46		2-Chloronaphthalene	1.354	1.271	6.1	107	0.00
47		2-Nitroaniline	0.297	0.285	4.0	112	0.00
48		Acenaphthylene	2.302	2.182	5.2	110	0.00
49		Dimethylphthalate	1.559	1.476	5.3	112	0.00
50		2,6-Dinitrotoluene	0.270	0.269	0.4	114	0.00
51	C	Acenaphthene	1.347	1.278	5.1	107	0.01
52		3-Nitroaniline	0.333	0.315	5.4	109	0.00
53	P	2,4-Dinitrophenol	0.064	0.048#	25.0	79	0.00
54		Dibenzofuran	1.886	1.838	2.5	109	0.00
55	P	4-Nitrophenol	0.242	0.227	6.2	106	0.00
56		2,4-Dinitrotoluene	0.347	0.345	0.6	106	0.00
57		Fluorene	1.670	1.674	-0.2	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.322	0.332	-3.1	112	0.00
59		Diethylphthalate	1.527	1.502	1.6	112	0.00
60		4-Chlorophenyl-phenylether	0.723	0.712	1.5	114	0.00
61		4-Nitroaniline	0.326	0.327	-0.3	109	0.00
62		Azobenzene	1.503	1.471	2.1	111	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
64		4,6-Dinitro-2-methylphenol	0.057	0.049	14.0	95	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.668	-0.6	119	0.00
66		4-Bromophenyl-phenylether	0.231	0.232	-0.4	121	0.00
67		Hexachlorobenzene	0.243	0.239	1.6	107	0.00
68		Atrazine	0.188	0.227	-20.7	132	0.00
69	C	Pentachlorophenol	0.089	0.093	-4.5	118	0.00
70		Phenanthrene	1.195	1.149	3.8	112	0.00
71		Anthracene	1.170	1.162	0.7	107	0.00
72		Carbazole	1.083	1.072	1.0	112	0.00
73		Di-n-butylphthalate	1.124	1.174	-4.4	112	0.01
74	C	Fluoranthene	1.253	1.282	-2.3	112	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	109	0.01
76		Benzidine	0.596	0.598	-0.3	101	0.00
77		Pyrene	1.246	1.253	-0.6	107	0.00
78	S	Terphenyl-d14	0.887	0.866	2.4	105	0.00
79		Butylbenzylphthalate	0.417	0.460	-10.3	118	0.01
80		Benzo(a)anthracene	1.189	1.264	-6.3	113	0.01
81		3,3'-Dichlorobenzidine	0.371	0.407	-9.7	114	0.01
82		Chrysene	1.116	1.121	-0.4	106	0.01
83		Bis(2-ethylhexyl)phthalate	0.645	0.700	-8.5	118	0.01
84	c	Di-n-octyl phthalate	1.002	1.104	-10.2	119	0.01
85		Indeno(1,2,3-cd)pyrene	1.178	1.318	-11.9	116	0.02
86	I	Perylene-d12	1.000	1.000	0.0	114	0.01
87		Benzo(b)fluoranthene	1.190	1.199	-0.8	114	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061015\
Data File : BF079855.D
Acq On : 10 Jun 2015 10:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 11 14:01:36 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	1.320	-6.0	111	0.01
89 C	Benzo(a)pyrene	1.125	1.196	-6.3	116	0.02
90	Dibenzo(a,h)anthracene	1.042	1.136	-9.0	115	0.02
91	Benzo(g,h,i)perylene	1.100	1.172	-6.5	116	0.02

(#) = Out of Range SPCC's out = 1 CCC's out = 0