Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020317\
Data File : BF092780.D
Acq On : 3 Feb 2017 17:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 23:20:07 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 2017
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	111	0.05
2		1,4-Dioxane	0.630	0.506	19.7	91	0.08
3		Pyridine	1.602	1.376	14.1	94	0.10
4		n-Nitrosodimethylamine	0.752	0.604	19.7	89	0.09
5	S	2-Fluorophenol	1.166	0.948	18.7	86	0.05
6		Aniline	2.046	2.103	-2.8	117	0.03
7	S	Phenol-d6	1.500	1.537	-2.5	114	0.03
8		2-Chlorophenol	1.286	1.263	1.8	108	0.05
9		Benzaldehyde	1.075	1.069	0.6	107	0.05
10	C	Phenol	1.755	1.667	5.0	111	0.05
11		bis(2-Chloroethyl)ether	1.401	1.508	-7.6	125	0.05
12		1,3-Dichlorobenzene	1.506	1.454	3.5	109	0.03
13	C	1,4-Dichlorobenzene	1.525	1.443	5.4	108	0.05
14		1,2-Dichlorobenzene	1.458	1.459	-0.1	110	0.05
15		Benzyl Alcohol	1.110	1.092	1.6	112	0.05
16		2,2'-oxybis(1-Chloropropane	2.434	2.500	-2.7	115	0.05
17		2-Methylphenol	1.103	1.175	-6.5	112	0.05
18		Hexachloroethane	0.520	0.445	14.4	94	0.05
19	P	n-Nitroso-di-n-propylamine	0.955	0.841	11.9	107	0.03
20		3+4-Methylphenols	1.319	1.196	9.3	98	0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.05
22		Acetophenone	0.457	0.477	-4.4	100	0.05
23	S	Nitrobenzene-d5	0.359	0.373	-3.9	96	0.05
24		Nitrobenzene	0.369	0.344	6.8	93	0.03
25		Isophorone	0.669	0.660	1.3	94	0.05
26	C	2-Nitrophenol	0.175	0.206	-17.7	101	0.05
27		2,4-Dimethylphenol	0.308	0.310	-0.6	90	0.05
28		bis(2-Chloroethoxy)methane	0.443	0.452	-2.0	96	0.05
29	C	2,4-Dichlorophenol	0.287	0.280	2.4	95	0.05
30		1,2,4-Trichlorobenzene	0.309	0.304	1.6	94	0.05
31		Naphthalene	1.014	0.965	4.8	92	0.05
32		Benzoic acid	0.156	0.167	-7.1	90	0.03
33		4-Chloroaniline	0.398	0.440	-10.6	101	0.05
34	C	Hexachlorobutadiene	0.170	0.163	4.1	90	0.05
35		Caprolactam	0.090	0.100	-11.1	96	0.05
36	C	4-Chloro-3-methylphenol	0.299	0.332	-11.0	102	0.06
37		2-Methylnaphthalene	0.651	0.673	-3.4	97	0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.06
39		1,2,4,5-Tetrachlorobenzene	0.561	0.564	-0.5	102	0.06
40	P	Hexachlorocyclopentadiene	0.253	0.221	12.6	86	0.05
41	S	2,4,6-Tribromophenol	0.145	0.152	-4.8	97	0.06
42	C	2,4,6-Trichlorophenol	0.369	0.388	-5.1	99	0.06
43		2,4,5-Trichlorophenol	0.404	0.425	-5.2	109	0.06
44	S	2-Fluorobiphenyl	1.104	1.050	4.9	90	0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020317\
Data File : BF092780.D
Acq On : 3 Feb 2017 17:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 23:20:07 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 2017
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	1.448	1.521	-5.0	101	0.06
46		2-Chloronaphthalene	1.133	1.129	0.4	94	0.06
47		2-Nitroaniline	0.381	0.398	-4.5	112	0.06
48		Acenaphthylene	1.957	1.904	2.7	105	0.05
49		Dimethylphthalate	1.347	1.367	-1.5	101	0.06
50		2,6-Dinitrotoluene	0.291	0.306	-5.2	103	0.06
51	C	Acenaphthene	1.170	1.194	-2.1	108	0.06
52		3-Nitroaniline	0.333	0.339	-1.8	96	0.06
53	P	2,4-Dinitrophenol	0.130	0.145	-11.5	106	0.05
54		Dibenzofuran	1.565	1.491	4.7	102	0.06
55	P	4-Nitrophenol	0.240	0.263	-9.6	111	0.06
56		2,4-Dinitrotoluene	0.387	0.365	5.7	84	0.05
57		Fluorene	1.204	1.289	-7.1	110	0.06
58		2,3,4,6-Tetrachlorophenol	0.270	0.286	-5.9	95	0.06
59		Diethylphthalate	1.270	1.312	-3.3	101	0.06
60		4-Chlorophenyl-phenylether	0.578	0.602	-4.2	107	0.06
61		4-Nitroaniline	0.341	0.340	0.3	101	0.06
62		Azobenzene	1.312	1.349	-2.8	103	0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.07
64		4,6-Dinitro-2-methylphenol	0.103	0.111	-7.8	96	0.05
65	c	n-Nitrosodiphenylamine	0.639	0.608	4.9	98	0.05
66		4-Bromophenyl-phenylether	0.209	0.212	-1.4	109	0.06
67		Hexachlorobenzene	0.206	0.210	-1.9	109	0.06
68		Atrazine	0.205	0.217	-5.9	118	0.06
69	C	Pentachlorophenol	0.089	0.102	-14.6	113	0.06
70		Phenanthrene	1.146	1.065	7.1	96	0.06
71		Anthracene	1.151	1.130	1.8	106	0.07
72		Carbazole	1.100	0.982	10.7	88	0.06
73		Di-n-butylphthalate	1.190	1.236	-3.9	109	0.06
74	C	Fluoranthene	1.182	1.072	9.3	92	0.06
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.06
76		Benzidine	0.852	0.839	1.5	100	0.07
77		Pyrene	1.497	1.355	9.5	87	0.06
78	S	Terphenyl-d14	0.851	0.911	-7.1	113	0.07
79		Butylbenzylphthalate	0.608	0.659	-8.4	110	0.07
80		Benzo(a)anthracene	1.223	1.199	2.0	98	0.06
81		3,3'-Dichlorobenzidine	0.436	0.430	1.4	97	0.07
82		Chrysene	1.145	1.208	-5.5	99	0.07
83		Bis(2-ethylhexyl)phthalate	0.831	0.868	-4.5	102	0.07
84	c	Di-n-octyl phthalate	1.354	1.464	-8.1	104	0.07
85		Indeno(1,2,3-cd)pyrene	0.887	0.918	-3.5	110	0.14
86	I	Perylene-d12	1.000	1.000	0.0	105	0.09
87		Benzo(b)fluoranthene	1.316	1.247	5.2	94	0.08

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020317\
Data File : BF092780.D
Acq On : 3 Feb 2017 17:09
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 23:20:07 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 30 14:50:05 2017
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	1.137	1.237	-8.8	124	0.08
89 C	Benzo(a)pyrene	1.139	1.136	0.3	105	0.09
90	Dibenzo(a,h)anthracene	0.920	0.905	1.6	109	0.13
91	Benzo(g,h,i)perylene	0.930	0.900	3.2	108	0.15

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