Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092716\
Data File : BF090249.D
Acq On : 27 Sep 2016 13:39
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 27 16:52:54 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 26 15:58:38 2016
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	96	0.00
2		1,4-Dioxane	0.666	0.566	15.0	89	-0.01
3		Pyridine	1.444	1.380	4.4	96	-0.01
4		n-Nitrosodimethylamine	0.431	0.445	-3.2	105	0.00
5	S	2-Fluorophenol	1.227	1.275	-3.9	103	0.00
6		Aniline	1.889	1.883	0.3	100	0.00
7	S	Phenol-d6	1.515	1.444	4.7	93	0.00
8		2-Chlorophenol	1.412	1.390	1.6	99	0.00
9		Benzaldehyde	0.660	0.727	-10.2	111	0.00
10	C	Phenol	1.791	1.561	12.8	83	0.00
11		bis(2-Chloroethyl)ether	1.397	1.381	1.1	97	0.00
12		1,3-Dichlorobenzene	1.475	1.391	5.7	98	-0.01
13	C	1,4-Dichlorobenzene	1.506	1.402	6.9	98	-0.01
14		1,2-Dichlorobenzene	1.342	1.304	2.8	97	0.00
15		Benzyl Alcohol	0.904	0.984	-8.8	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.669	1.487	10.9	90	0.00
17		2-Methylphenol	1.112	1.135	-2.1	101	0.00
18		Hexachloroethane	0.537	0.502	6.5	93	-0.01
19	P	n-Nitroso-di-n-propylamine	0.890	0.934	-4.9	104	0.00
20		3+4-Methylphenols	1.335	1.408	-5.5	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.482	0.448	7.1	92	0.00
23	S	Nitrobenzene-d5	0.345	0.337	2.3	100	-0.01
24		Nitrobenzene	0.359	0.373	-3.9	100	0.00
25		Isophorone	0.721	0.693	3.9	99	0.00
26	C	2-Nitrophenol	0.177	0.195	-10.2	110	0.00
27		2,4-Dimethylphenol	0.315	0.335	-6.3	109	0.00
28		bis(2-Chloroethoxy)methane	0.436	0.442	-1.4	102	0.00
29	C	2,4-Dichlorophenol	0.276	0.268	2.9	95	-0.01
30		1,2,4-Trichlorobenzene	0.276	0.278	-0.7	100	0.00
31		Naphthalene	0.952	0.958	-0.6	100	-0.01
32		Benzoic acid	0.216	0.257	-19.0	109	0.02
33		4-Chloroaniline	0.395	0.428	-8.4	105	0.00
34	C	Hexachlorobutadiene	0.145	0.145	0.0	102	0.00
35		Caprolactam	0.091	0.096	-5.5	110	0.00
36	C	4-Chloro-3-methylphenol	0.312	0.325	-4.2	108	0.00
37		2-Methylnaphthalene	0.592	0.614	-3.7	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	0.459	0.482	-5.0	109	0.00
40	P	Hexachlorocyclopentadiene	0.227	0.243	-7.0	101	0.00
41	S	2,4,6-Tribromophenol	0.172	0.185	-7.6	107	0.00
42	C	2,4,6-Trichlorophenol	0.327	0.350	-7.0	102	0.00
43		2,4,5-Trichlorophenol	0.339	0.337	0.6	102	-0.01
44	S	2-Fluorobiphenyl	1.019	0.986	3.2	102	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092716\
Data File : BF090249.D
Acq On : 27 Sep 2016 13:39
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 27 16:52:54 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 26 15:58:38 2016
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	1.431	1.506	-5.2	103	0.00
46		2-Chloronaphthalene	1.056	0.965	8.6	100	-0.01
47		2-Nitroaniline	0.318	0.309	2.8	103	-0.01
48		Acenaphthylene	1.714	1.617	5.7	101	0.00
49		Dimethylphthalate	1.274	1.281	-0.5	109	-0.01
50		2,6-Dinitrotoluene	0.284	0.258	9.2	88	0.00
51	C	Acenaphthene	1.018	0.911	10.5	91	0.00
52		3-Nitroaniline	0.333	0.357	-7.2	111	0.00
53	P	2,4-Dinitrophenol	0.124	0.176	-41.9#	122	0.00
54		Dibenzofuran	1.339	1.296	3.2	105	0.00
55	P	4-Nitrophenol	0.228	0.233	-2.2	98	0.00
56		2,4-Dinitrotoluene	0.336	0.319	5.1	90	0.00
57		Fluorene	1.009	1.096	-8.6	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.254	0.271	-6.7	103	0.00
59		Diethylphthalate	1.230	1.247	-1.4	102	0.00
60		4-Chlorophenyl-phenylether	0.460	0.454	1.3	106	0.00
61		4-Nitroaniline	0.339	0.348	-2.7	98	0.00
62		Azobenzene	1.281	1.299	-1.4	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.01
64		4,6-Dinitro-2-methylphenol	0.125	0.147	-17.6	124	0.00
65	c	n-Nitrosodiphenylamine	0.658	0.662	-0.6	109	0.00
66		4-Bromophenyl-phenylether	0.197	0.179	9.1	99	-0.01
67		Hexachlorobenzene	0.229	0.214	6.6	108	-0.01
68		Atrazine	0.180	0.184	-2.2	118	-0.01
69	C	Pentachlorophenol	0.130	0.136	-4.6	105	-0.01
70		Phenanthrene	0.935	0.990	-5.9	106	0.00
71		Anthracene	0.981	0.901	8.2	100	-0.01
72		Carbazole	1.037	0.911	12.2	98	-0.01
73		Di-n-butylphthalate	1.206	1.234	-2.3	109	0.00
74	C	Fluoranthene	1.004	0.941	6.3	102	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	103	-0.01
76		Benzidine	0.669	0.600	10.3	92	-0.01
77		Pyrene	1.369	1.351	1.3	109	0.00
78	S	Terphenyl-d14	0.788	0.711	9.8	106	0.00
79		Butylbenzylphthalate	0.669	0.672	-0.4	107	-0.01
80		Benzo(a)anthracene	1.076	1.154	-7.2	110	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.456	-2.7	109	-0.01
82		Chrysene	1.048	1.084	-3.4	108	0.00
83		Bis(2-ethylhexyl)phthalate	0.857	0.929	-8.4	110	0.00
84	c	Di-n-octyl phthalate	1.476	1.668	-13.0	110	0.00
85		Indeno(1,2,3-cd)pyrene	0.847	0.955	-12.8	112	0.00
86	I	Perylene-d12	1.000	1.000	0.0	108	0.00
87		Benzo(b)fluoranthene	1.107	1.123	-1.4	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092716\
Data File : BF090249.D
Acq On : 27 Sep 2016 13:39
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 27 16:52:54 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 26 15:58:38 2016
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	1.039	0.933	10.2	107	0.00
89 C	Benzo(a)pyrene	1.042	0.970	6.9	105	-0.01
90	Dibenzo(a,h)anthracene	0.813	0.795	2.2	110	-0.01
91	Benzo(g,h,i)perylene	0.796	0.767	3.6	110	0.00

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