Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091516\
Data File : BF090082.D
Acq On : 15 Sep 2016 9:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 11:31:34 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 15 11:18:43 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	102	0.00
2		1,4-Dioxane	40.000	39.823	0.4	107	0.01
3		Pyridine	40.000	42.980	-7.4	124	0.00
4		n-Nitrosodimethylamine	40.000	41.086	-2.7	119	0.00
5	S	2-Fluorophenol	80.000	89.480	-11.9	106	0.00
6		Aniline	40.000	42.327	-5.8	107	0.00
7	S	Phenol-d6	80.000	78.506	1.9	105	0.00
8		2-Chlorophenol	40.000	41.540	-3.8	107	0.01
9		Benzaldehyde	40.000	42.315	-5.8	105	0.00
10	C	Phenol	40.000	37.092	7.3	103	0.00
11		bis(2-Chloroethyl)ether	40.000	39.941	0.1	105	0.00
12		1,3-Dichlorobenzene	40.000	40.179	-0.4	102	0.00
13	C	1,4-Dichlorobenzene	40.000	40.731	-1.8	103	0.00
14		1,2-Dichlorobenzene	40.000	37.500	6.3	101	0.00
15		Benzyl Alcohol	40.000	40.338	-0.8	103	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.711	-4.3	102	0.00
17		2-Methylphenol	40.000	41.063	-2.7	104	0.00
18		Hexachloroethane	40.000	38.693	3.3	102	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.614	3.5	100	0.00
20		3+4-Methylphenols	40.000	41.741	-4.4	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	42.382	-6.0	103	0.00
23	S	Nitrobenzene-d5	80.000	81.827	-2.3	100	0.00
24		Nitrobenzene	40.000	39.161	2.1	102	0.00
25		Isophorone	40.000	39.903	0.2	98	0.00
26	C	2-Nitrophenol	40.000	42.721	-6.8	98	0.00
27		2,4-Dimethylphenol	40.000	42.550	-6.4	97	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.180	2.1	95	0.00
29	C	2,4-Dichlorophenol	40.000	41.427	-3.6	97	0.00
30		1,2,4-Trichlorobenzene	40.000	39.572	1.1	96	0.00
31		Naphthalene	40.000	38.807	3.0	102	0.00
32		Benzoic acid	40.000	39.720	0.7	94	0.00
33		4-Chloroaniline	40.000	39.323	1.7	102	0.00
34	C	Hexachlorobutadiene	40.000	39.234	1.9	103	0.00
35		Caprolactam	40.000	40.326	-0.8	105	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.749	0.6	104	0.00
37		2-Methylnaphthalene	40.000	40.172	-0.4	103	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	36.864	7.8	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	38.534	3.7	100	0.00
41	S	2,4,6-Tribromophenol	80.000	78.655	1.7	103	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.230	4.4	102	0.00
43		2,4,5-Trichlorophenol	40.000	39.117	2.2	103	0.00
44	S	2-Fluorobiphenyl	80.000	72.116	9.9	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091516\
Data File : BF090082.D
Acq On : 15 Sep 2016 9:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 11:31:34 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 15 11:18:43 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	35.824	10.4	101	0.00
46		2-Chloronaphthalene	40.000	38.884	2.8	104	0.00
47		2-Nitroaniline	40.000	42.410	-6.0	105	0.00
48		Acenaphthylene	40.000	39.950	0.1	104	0.00
49		Dimethylphthalate	40.000	40.692	-1.7	103	0.00
50		2,6-Dinitrotoluene	40.000	40.981	-2.5	109	0.01
51	C	Acenaphthene	40.000	40.068	-0.2	101	0.00
52		3-Nitroaniline	40.000	39.989	0.0	102	0.00
53	P	2,4-Dinitrophenol	40.000	40.800	-2.0	99	0.00
54		Dibenzofuran	40.000	40.326	-0.8	102	0.00
55	P	4-Nitrophenol	40.000	41.180	-2.9	103	0.00
56		2,4-Dinitrotoluene	40.000	35.862	10.3	103	0.00
57		Fluorene	40.000	37.037	7.4	102	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	36.524	8.7	103	0.00
59		Diethylphthalate	40.000	36.386	9.0	103	0.00
60		4-Chlorophenyl-phenylether	40.000	38.848	2.9	101	0.00
61		4-Nitroaniline	40.000	40.813	-2.0	103	0.00
62		Azobenzene	40.000	36.379	9.1	98	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	40.000	41.977	-4.9	100	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.409	6.5	99	0.00
66		4-Bromophenyl-phenylether	40.000	42.339	-5.8	102	0.00
67		Hexachlorobenzene	40.000	40.054	-0.1	100	0.00
68		Atrazine	40.000	44.451	-11.1	102	0.00
69	C	Pentachlorophenol	40.000	44.463	-11.2	103	0.00
70		Phenanthrene	40.000	39.311	1.7	102	0.00
71		Anthracene	40.000	38.706	3.2	102	0.00
72		Carbazole	40.000	40.536	-1.3	100	0.00
73		Di-n-butylphthalate	40.000	36.909	7.7	104	0.00
74	C	Fluoranthene	40.000	37.191	7.0	102	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
76		Benzidine	40.000	40.683	-1.7	102	0.00
77		Pyrene	40.000	38.049	4.9	105	0.00
78	S	Terphenyl-d14	80.000	73.244	8.4	103	0.00
79		Butylbenzylphthalate	40.000	38.768	3.1	103	0.00
80		Benzo(a)anthracene	40.000	38.136	4.7	102	0.00
81		3,3'-Dichlorobenzidine	40.000	39.447	1.4	102	0.00
82		Chrysene	40.000	32.933	17.7	104	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.016	7.5	104	0.00
84	c	Di-n-octyl phthalate	40.000	36.572	8.6	103	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	35.167	12.1	107	0.01
86	I	Perylene-d12	20.000	20.000	0.0	108	0.00
87		Benzo(b)fluoranthene	40.000	47.043	-17.6	148	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091516\
Data File : BF090082.D
Acq On : 15 Sep 2016 9:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 15 11:31:34 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 15 11:18:43 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	32.563	18.6	75	0.00
89 C	Benzo(a)pyrene	40.000	39.890	0.3	108	0.00
90	Dibenzo(a,h)anthracene	40.000	38.007	5.0	106	0.00
91	Benzo(g,h,i)perylene	40.000	37.601	6.0	107	0.00

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