Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088365.D
Acq On : 25 Jun 2016 9:17
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 99 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jun 27 08:43:48 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 27 04:27:12 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	429	-0.13
2		1,4-Dioxane	40.000	0.000	100.0#	0	-2.41#
3		Pyridine	40.000	0.018	100.0#	0	0.16
4		n-Nitrosodimethylamine	40.000	0.028	99.9#	0	-0.07
5	S	2-Fluorophenol	80.000	0.000	100.0#	0	-5.43#
6		Aniline	40.000	6.806	83.0#	73	-0.27
7	S	Phenol-d6	80.000	0.325	99.6#	2	-0.13
8		2-Chlorophenol	40.000	7.461	81.3#	82	-0.27
9		Benzaldehyde	40.000	5.497	86.3#	77	-0.29
10	C	Phenol	40.000	6.909	82.7#	101	-0.25
11		bis(2-Chloroethyl)ether	40.000	8.390	79.0#	97	-0.27
12		1,3-Dichlorobenzene	40.000	5.727	85.7#	62	-0.06
13	C	1,4-Dichlorobenzene	40.000	7.571	81.1#	85	-0.27
14		1,2-Dichlorobenzene	40.000	6.507	83.7#	69	-0.29
15		Benzyl Alcohol	40.000	6.579	83.6#	68	-0.26
16		2,2'-oxybis(1-Chloropropane	40.000	5.603	86.0#	60	-0.27
17		2-Methylphenol	40.000	8.827	77.9#	92	-0.09
18		Hexachloroethane	40.000	7.448	81.4#	79	-0.29
19	P	n-Nitroso-di-n-propylamine	40.000	6.989	82.5#	82	-0.26
20		3+4-Methylphenols	40.000	7.457	81.4#	86	-0.24
21	I	Naphthalene-d8	20.000	20.000	0.0	0	-0.39
22		Acetophenone	40.000	13982.358	-34855.9#	84	-0.26
23	S	Nitrobenzene-d5	80.000	47082.423	-58753.0#	135	-0.26
24		Nitrobenzene	40.000	12493.501	-31133.8#	79	-0.26
25		Isophorone	40.000	13777.128	-34342.8#	79	-0.27
26	C	2-Nitrophenol	40.000	14781.944	-36854.9#	75	-0.27
27		2,4-Dimethylphenol	40.000	12191.370	-30378.4#	69	-0.25
28		bis(2-Chloroethoxy)methane	40.000	649.830	-1524.6#	4	-0.41
29	C	2,4-Dichlorophenol	40.000	13912.649	-34681.6#	80	-0.25
30		1,2,4-Trichlorobenzene	40.000	13126.025	-32715.1#	79	-0.27
31		Naphthalene	40.000	13420.941	-33452.4#	83	-0.27
32		Benzoic acid	40.000	6639.844	-16499.6#	33	-0.22
33		4-Chloroaniline	40.000	3963.621	-9809.1#	22	-0.33
34	C	Hexachlorobutadiene	40.000	12911.285	-32178.2#	73	-0.29
35		Caprolactam	40.000	12007.234	-29918.1#	64	-0.24
36	C	4-Chloro-3-methylphenol	40.000	14053.913	-35034.8#	84	-0.24
37		2-Methylnaphthalene	40.000	5174.191	-12835.5#	29	-0.40
38	I	Acenaphthene-d10	20.000	20.000	0.0	1	-0.48
39		1,2,4,5-Tetrachlorobenzene	40.000	-20.000	150.0#	82	-0.27
40	P	Hexachlorocyclopentadiene	40.000	1319.173	-3197.9#	26	-0.27
41	S	2,4,6-Tribromophenol	80.000	0.000	100.0#	0	-10.67#
42	C	2,4,6-Trichlorophenol	40.000	3617.034	-8942.6#	70	-0.26
43		2,4,5-Trichlorophenol	40.000	3476.641	-8591.6#	71	-0.31
44	S	2-Fluorobiphenyl	80.000	12384.599	-15380.7#	117	-0.27

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088365.D
Acq On : 25 Jun 2016 9:17
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 99 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jun 27 08:43:48 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 27 04:27:12 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	4388.479	-10871.2#	75	-0.27
46		2-Chloronaphthalene	40.000	3625.908	-8964.8#	77	-0.27
47		2-Nitroaniline	40.000	142.911	-257.3#	3	-0.40
48		Acenaphthylene	40.000	3673.188	-9083.0#	74	-0.29
49		Dimethylphthalate	40.000	4098.405	-10146.0#	89	-0.26
50		2,6-Dinitrotoluene	40.000	493.251	-1133.1#	11	-0.62#
51	C	Acenaphthene	40.000	3534.888	-8737.2#	70	-0.29
52		3-Nitroaniline	40.000	3380.836	-8352.1#	64	-0.26
53	P	2,4-Dinitrophenol	40.000	0.000	100.0#	0	-9.94#
54		Dibenzofuran	40.000	332.946	-732.4#	6	-0.62#
55	P	4-Nitrophenol	40.000	348.467	-771.2#	6	-0.42
56		2,4-Dinitrotoluene	40.000	2861.244	-7053.1#	61	-0.66#
57		Fluorene	40.000	4674.722	-11586.8#	80	-0.29
58		2,3,4,6-Tetrachlorophenol	40.000	3156.776	-7791.9#	59	-0.31
59		Diethylphthalate	40.000	3631.569	-8978.9#	75	-0.27
60		4-Chlorophenyl-phenylether	40.000	3349.021	-8272.6#	72	-0.29
61		4-Nitroaniline	40.000	337.900	-744.8#	6	-0.65#
62		Azobenzene	40.000	3454.020	-8535.0#	68	-0.29
63	I	Phenanthrene-d10	20.000	20.000	0.0	0	-0.56#
64		4,6-Dinitro-2-methylphenol	40.000	1018.922	-2447.3#	6	-0.55#
65	c	n-Nitrosodiphenylamine	40.000	292.261	-630.7#	2	-0.62#
66		4-Bromophenyl-phenylether	40.000	0.000	100.0#	0	-10.91#
67		Hexachlorobenzene	40.000	0.000	100.0#	0	-10.97#
68		Atrazine	40.000	0.000	100.0#	0	-11.07#
69	C	Pentachlorophenol	40.000	0.000	100.0#	0	-11.17#
70		Phenanthrene	40.000	0.000	100.0#	0	-11.38#
71		Anthracene	40.000	11482.502	-28606.3#	72	-0.34
72		Carbazole	40.000	419.879	-949.7#	3	-0.70#
73		Di-n-butylphthalate	40.000	53.153	-32.9#	0	-0.49
74	C	Fluoranthene	40.000	0.000	100.0#	0	-12.57#
75	I	Chrysene-d12	20.000	0.000	100.0#	0	-13.99#
76		Benzidine	40.000	0.000	100.0#	6	-0.02
77		Pyrene	40.000	0.000	100.0#	0	-12.80#
78	S	Terphenyl-d14	80.000	0.000	100.0#	0	-12.95#
79		Butylbenzylphthalate	40.000	0.000	100.0#	0	-13.42#
80		Benzo(a)anthracene	40.000	0.000	100.0#	81	-0.25
81		3,3'-Dichlorobenzidine	40.000	0.000	100.0#	0	-13.94#
82		Chrysene	40.000	0.000	100.0#	0	-14.01#
83		Bis(2-ethylhexyl)phthalate	40.000	0.000	100.0#	0	-13.98#
84	c	Di-n-octyl phthalate	40.000	0.000	100.0#	0	-14.58#
85		Indeno(1,2,3-cd)pyrene	40.000	0.000	100.0#	62	-0.15
86	I	Perylene-d12	20.000	0.000	100.0#	0	-15.39#
87		Benzo(b)fluoranthene	40.000	0.000	100.0#	52	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088365.D
Acq On : 25 Jun 2016 9:17
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 99 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jun 27 08:43:48 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 27 04:27:12 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	0.000	100.0#	90	-0.03
89 C	Benzo(a)pyrene	40.000	0.000	100.0#	0	-15.35#
90	Dibenzo(a,h)anthracene	40.000	0.000	100.0#	2	-0.16
91	Benzo(g,h,i)perylene	40.000	0.000	100.0#	0	-0.21

(#) = Out of Range SPCC's out = 0 CCC's out = 13