Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082516\
Data File : BF089911.D
Acq On : 25 Aug 2016 15:06
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 25 19:30:01 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 19 14:02:57 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	124	-0.02
2		1,4-Dioxane	40.000	40.709	-1.8	127	-0.06
3		Pyridine	40.000	40.425	-1.1	118	-0.07
4		n-Nitrosodimethylamine	40.000	39.739	0.7	118	-0.06
5	S	2-Fluorophenol	80.000	76.590	4.3	125	-0.02
6		Aniline	40.000	38.656	3.4	119	-0.01
7	S	Phenol-d6	80.000	73.877	7.7	109	-0.01
8		2-Chlorophenol	40.000	38.285	4.3	115	-0.02
9		Benzaldehyde	40.000	37.186	7.0	107	-0.02
10	C	Phenol	40.000	32.878	17.8	96	-0.02
11		bis(2-Chloroethyl)ether	40.000	42.367	-5.9	122	-0.02
12		1,3-Dichlorobenzene	40.000	41.983	-5.0	120	-0.02
13	C	1,4-Dichlorobenzene	40.000	39.189	2.0	113	-0.02
14		1,2-Dichlorobenzene	40.000	37.641	5.9	118	-0.02
15		Benzyl Alcohol	40.000	33.460	16.3	101	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	38.976	2.6	114	-0.03
17		2-Methylphenol	40.000	42.159	-5.4	132	-0.01
18		Hexachloroethane	40.000	40.057	-0.1	126	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	33.185	17.0	96	-0.02
20		3+4-Methylphenols	40.000	39.840	0.4	134	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	123	-0.02
22		Acetophenone	40.000	37.231	6.9	111	-0.02
23	S	Nitrobenzene-d5	80.000	69.871	12.7	113	-0.02
24		Nitrobenzene	40.000	40.388	-1.0	120	-0.02
25		Isophorone	40.000	37.741	5.6	116	-0.02
26	C	2-Nitrophenol	40.000	38.685	3.3	127	-0.02
27		2,4-Dimethylphenol	40.000	40.512	-1.3	136	-0.01
28		bis(2-Chloroethoxy)methane	40.000	34.351	14.1	99	-0.02
29	C	2,4-Dichlorophenol	40.000	43.675	-9.2	142	-0.01
30		1,2,4-Trichlorobenzene	40.000	38.880	2.8	117	-0.02
31		Naphthalene	40.000	38.413	4.0	119	-0.02
32		Benzoic acid	40.000	40.260	-0.6	125	0.01
33		4-Chloroaniline	40.000	34.559	13.6	104	-0.02
34	C	Hexachlorobutadiene	40.000	38.915	2.7	120	-0.02
35		Caprolactam	40.000	40.809	-2.0	128	0.00
36	C	4-Chloro-3-methylphenol	40.000	34.727	13.2	118	-0.01
37		2-Methylnaphthalene	40.000	34.755	13.1	103	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	130	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	36.873	7.8	116	-0.02
40	P	Hexachlorocyclopentadiene	40.000	47.002	-17.5	136	-0.02
41	S	2,4,6-Tribromophenol	80.000	78.835	1.5	129	-0.01
42	C	2,4,6-Trichlorophenol	40.000	41.536	-3.8	127	-0.01
43		2,4,5-Trichlorophenol	40.000	40.180	-0.4	129	-0.01
44	S	2-Fluorobiphenyl	80.000	73.180	8.5	127	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082516\
Data File : BF089911.D
Acq On : 25 Aug 2016 15:06
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 25 19:30:01 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 19 14:02:57 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	33.642	15.9	102	-0.02
46		2-Chloronaphthalene	40.000	40.248	-0.6	130	-0.01
47		2-Nitroaniline	40.000	36.356	9.1	104	-0.02
48		Acenaphthylene	40.000	39.791	0.5	129	-0.01
49		Dimethylphthalate	40.000	37.224	6.9	114	-0.02
50		2,6-Dinitrotoluene	40.000	37.499	6.3	129	-0.01
51	C	Acenaphthene	40.000	37.127	7.2	114	-0.01
52		3-Nitroaniline	40.000	45.046	-12.6	130	-0.02
53	P	2,4-Dinitrophenol	40.000	41.001	-2.5	146	-0.01
54		Dibenzofuran	40.000	38.334	4.2	122	-0.01
55	P	4-Nitrophenol	40.000	45.131	-12.8	163	0.00
56		2,4-Dinitrotoluene	40.000	39.581	1.0	114	-0.01
57		Fluorene	40.000	38.625	3.4	132	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	37.944	5.1	118	-0.02
59		Diethylphthalate	40.000	37.112	7.2	113	-0.02
60		4-Chlorophenyl-phenylether	40.000	35.396	11.5	119	-0.01
61		4-Nitroaniline	40.000	47.091	-17.7	132	-0.01
62		Azobenzene	40.000	35.596	11.0	108	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	128	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	51.654	-29.1#	179	0.00
65	c	n-Nitrosodiphenylamine	40.000	35.111	12.2	109	-0.01
66		4-Bromophenyl-phenylether	40.000	35.094	12.3	104	-0.02
67		Hexachlorobenzene	40.000	38.142	4.6	116	-0.02
68		Atrazine	40.000	42.788	-7.0	134	-0.01
69	C	Pentachlorophenol	40.000	40.239	-0.6	120	-0.02
70		Phenanthrene	40.000	34.083	14.8	103	-0.02
71		Anthracene	40.000	39.832	0.4	136	-0.01
72		Carbazole	40.000	39.449	1.4	115	-0.02
73		Di-n-butylphthalate	40.000	40.183	-0.5	130	-0.01
74	C	Fluoranthene	40.000	43.021	-7.6	132	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	146	-0.03
76		Benzidine	40.000	49.022	-22.6	163	-0.01
77		Pyrene	40.000	32.040	19.9	113	-0.02
78	S	Terphenyl-d14	80.000	55.616	30.5#	101	-0.02
79		Butylbenzylphthalate	40.000	44.443	-11.1	162	-0.02
80		Benzo(a)anthracene	40.000	41.653	-4.1	153	-0.03
81		3,3'-Dichlorobenzidine	40.000	49.014	-22.5	176	-0.03
82		Chrysene	40.000	41.949	-4.9	156	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	41.356	-3.4	148	-0.03
84	c	Di-n-octyl phthalate	40.000	49.272	-23.2#	169	-0.06
85		Indeno(1,2,3-cd)pyrene	40.000	32.992	17.5	125	-0.08
86	I	Perylene-d12	20.000	20.000	0.0	159	-0.08
87		Benzo(b)fluoranthene	40.000	41.300	-3.2	155	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082516\
Data File : BF089911.D
Acq On : 25 Aug 2016 15:06
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 25 19:30:01 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 19 14:02:57 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	44.378	-10.9	180	-0.06
89 C	Benzo(a)pyrene	40.000	40.557	-1.4	154	-0.07
90	Dibenzo(a,h)anthracene	40.000	31.953	20.1	123	-0.09
91	Benzo(g,h,i)perylene	40.000	30.676	23.3	119	-0.08

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