Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
Data File : BF084466.D
Acq On : 14 Jan 2016 2:42
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 14 15:47:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 14 14:15:52 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	78	0.00
2		1,4-Dioxane	40.000	41.156	-2.9	77	0.01
3		Pyridine	40.000	39.881	0.3	71	0.00
4		n-Nitrosodimethylamine	40.000	35.930	10.2	66	0.01
5	S	2-Fluorophenol	80.000	81.423	-1.8	84	0.00
6		Aniline	40.000	38.199	4.5	72	0.00
7	S	Phenol-d6	80.000	75.961	5.0	73	0.00
8		2-Chlorophenol	40.000	40.228	-0.6	79	0.00
9		Benzaldehyde	40.000	34.687	13.3	68	-0.01
10	C	Phenol	40.000	35.371	11.6	64	-0.01
11		bis(2-Chloroethyl)ether	40.000	40.371	-0.9	82	-0.01
12		1,3-Dichlorobenzene	40.000	40.718	-1.8	81	0.00
13	C	1,4-Dichlorobenzene	40.000	41.799	-4.5	78	0.00
14		1,2-Dichlorobenzene	40.000	36.373	9.1	76	0.00
15		Benzyl Alcohol	40.000	33.296	16.8	62	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.174	12.1	70	0.00
17		2-Methylphenol	40.000	40.916	-2.3	74	0.00
18		Hexachloroethane	40.000	38.219	4.5	72	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	34.302	14.2	69	0.00
20		3+4-Methylphenols	40.000	35.975	10.1	75	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	79	0.00
22		Acetophenone	40.000	38.954	2.6	72	0.00
23	S	Nitrobenzene-d5	80.000	70.630	11.7	72	0.00
24		Nitrobenzene	40.000	42.950	-7.4	77	0.00
25		Isophorone	40.000	37.699	5.8	74	0.00
26	C	2-Nitrophenol	40.000	41.032	-2.6	83	-0.01
27		2,4-Dimethylphenol	40.000	35.587	11.0	66	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.454	3.9	82	0.00
29	C	2,4-Dichlorophenol	40.000	40.597	-1.5	82	0.00
30		1,2,4-Trichlorobenzene	40.000	40.757	-1.9	78	0.00
31		Naphthalene	40.000	40.757	-1.9	81	0.00
32		Benzoic acid	40.000	31.295	21.8	59	0.01
33		4-Chloroaniline	40.000	39.306	1.7	80	-0.01
34	C	Hexachlorobutadiene	40.000	36.796	8.0	73	0.00
35		Caprolactam	40.000	40.839	-2.1	72	0.00
36	C	4-Chloro-3-methylphenol	40.000	36.713	8.2	76	0.00
37		2-Methylnaphthalene	40.000	36.390	9.0	75	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	76	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	42.098	-5.2	79	0.00
40	P	Hexachlorocyclopentadiene	40.000	35.462	11.3	65	0.00
41	S	2,4,6-Tribromophenol	80.000	77.341	3.3	69	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.314	4.2	74	0.00
43		2,4,5-Trichlorophenol	40.000	39.143	2.1	71	0.00
44	S	2-Fluorobiphenyl	80.000	76.107	4.9	71	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
Data File : BF084466.D
Acq On : 14 Jan 2016 2:42
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 14 15:47:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 14 14:15:52 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	39.544	1.1	80	0.00
46		2-Chloronaphthalene	40.000	37.108	7.2	78	0.00
47		2-Nitroaniline	40.000	45.760	-14.4	90	0.00
48		Acenaphthylene	40.000	39.446	1.4	71	0.00
49		Dimethylphthalate	40.000	42.145	-5.4	76	0.00
50		2,6-Dinitrotoluene	40.000	42.200	-5.5	72	0.00
51	C	Acenaphthene	40.000	38.823	2.9	68	0.00
52		3-Nitroaniline	40.000	44.060	-10.2	77	0.00
53	P	2,4-Dinitrophenol	40.000	45.546	-13.9	83	0.00
54		Dibenzofuran	40.000	39.490	1.3	69	0.00
55	P	4-Nitrophenol	40.000	41.010	-2.5	64	0.00
56		2,4-Dinitrotoluene	40.000	40.454	-1.1	65	0.00
57		Fluorene	40.000	37.426	6.4	67	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.948	-2.4	73	0.00
59		Diethylphthalate	40.000	41.525	-3.8	76	0.00
60		4-Chlorophenyl-phenylether	40.000	46.722	-16.8	82	0.00
61		4-Nitroaniline	40.000	45.376	-13.4	88	0.00
62		Azobenzene	40.000	41.929	-4.8	78	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	80	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.030	-0.1	86	0.00
65	c	n-Nitrosodiphenylamine	40.000	35.782	10.5	71	0.00
66		4-Bromophenyl-phenylether	40.000	39.053	2.4	74	0.00
67		Hexachlorobenzene	40.000	35.413	11.5	75	0.00
68		Atrazine	40.000	40.515	-1.3	72	0.00
69	C	Pentachlorophenol	40.000	33.695	15.8	59	0.00
70		Phenanthrene	40.000	36.266	9.3	67	0.00
71		Anthracene	40.000	42.362	-5.9	88	0.00
72		Carbazole	40.000	39.518	1.2	76	0.00
73		Di-n-butylphthalate	40.000	43.171	-7.9	82	0.00
74	C	Fluoranthene	40.000	37.057	7.4	77	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	82	0.00
76		Benzidine	40.000	33.144	17.1	67	0.00
77		Pyrene	40.000	33.245	16.9	77	0.00
78	S	Terphenyl-d14	80.000	63.889	20.1	76	0.00
79		Butylbenzylphthalate	40.000	35.983	10.0	72	0.00
80		Benzo(a)anthracene	40.000	34.303	14.2	74	0.00
81		3,3'-Dichlorobenzidine	40.000	38.710	3.2	78	0.00
82		Chrysene	40.000	33.607	16.0	70	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	34.324	14.2	73	0.00
84	c	Di-n-octyl phthalate	40.000	33.269	16.8	65	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	36.419	9.0	76	0.00
86	I	Perylene-d12	20.000	20.000	0.0	77	0.00
87		Benzo(b)fluoranthene	40.000	43.725	-9.3	81	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
Data File : BF084466.D
Acq On : 14 Jan 2016 2:42
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 14 15:47:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 14 14:15:52 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	33.753	15.6	66	0.00
89 C	Benzo(a)pyrene	40.000	39.674	0.8	74	0.00
90	Dibenzo(a,h)anthracene	40.000	37.451	6.4	76	0.00
91	Benzo(g,h,i)perylene	40.000	37.926	5.2	79	-0.01

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