Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070215\
Data File : BF080340.D
Acq On : 3 Jul 2015 12:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 03 23:40:44 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 03 22:57:58 2015
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	90	-0.03
2		1,4-Dioxane	40.000	141.829	-254.6#	394	-0.06
3		Pyridine	40.000	35.938	10.2	77	-0.04
4		n-Nitrosodimethylamine	40.000	35.613	11.0	78	-0.04
5	S	2-Fluorophenol	80.000	71.543	10.6	85	-0.03
6		Aniline	40.000	36.100	9.7	85	-0.02
7	S	Phenol-d6	80.000	69.804	12.7	85	-0.03
8		2-Chlorophenol	40.000	37.221	6.9	91	-0.03
9		Benzaldehyde	40.000	32.626	18.4	86	-0.03
10	C	Phenol	40.000	32.201	19.5	77	-0.03
11		bis(2-Chloroethyl)ether	40.000	34.054	14.9	77	-0.03
12		1,3-Dichlorobenzene	40.000	37.455	6.4	92	-0.03
13	C	1,4-Dichlorobenzene	40.000	35.766	10.6	84	-0.03
14		1,2-Dichlorobenzene	40.000	36.431	8.9	86	-0.03
15		Benzyl Alcohol	40.000	36.403	9.0	83	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	33.289	16.8	79	-0.03
17		2-Methylphenol	40.000	33.540	16.2	78	-0.03
18		Hexachloroethane	40.000	34.270	14.3	80	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	30.935	22.7	75	-0.03
20		3+4-Methylphenols	40.000	34.689	13.3	81	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	91	-0.03
22		Acetophenone	40.000	38.208	4.5	87	-0.03
23	S	Nitrobenzene-d5	80.000	69.892	12.6	85	-0.02
24		Nitrobenzene	40.000	35.647	10.9	79	-0.03
25		Isophorone	40.000	35.434	11.4	81	-0.03
26	C	2-Nitrophenol	40.000	33.548	16.1	75	-0.02
27		2,4-Dimethylphenol	40.000	34.298	14.3	80	-0.02
28		bis(2-Chloroethoxy)methane	40.000	35.500	11.3	85	-0.03
29	C	2,4-Dichlorophenol	40.000	34.572	13.6	78	-0.03
30		1,2,4-Trichlorobenzene	40.000	34.354	14.1	81	-0.02
31		Naphthalene	40.000	37.258	6.9	90	-0.03
32		Benzoic acid	40.000	54.655	-36.6#	162	-0.02
33		4-Chloroaniline	40.000	35.531	11.2	84	-0.03
34	C	Hexachlorobutadiene	40.000	34.287	14.3	82	-0.03
35		Caprolactam	40.000	34.975	12.6	84	-0.02
36	C	4-Chloro-3-methylphenol	40.000	35.482	11.3	81	-0.03
37		2-Methylnaphthalene	40.000	33.765	15.6	80	-0.03
38	I	Acenaphthene-d10	20.000	20.000	0.0	85	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	36.758	8.1	81	-0.03
40	P	Hexachlorocyclopentadiene	40.000	30.457	23.9	65	-0.03
41	S	2,4,6-Tribromophenol	80.000	77.818	2.7	84	-0.03
42	C	2,4,6-Trichlorophenol	40.000	34.457	13.9	76	-0.03
43		2,4,5-Trichlorophenol	40.000	38.967	2.6	89	-0.03
44	S	2-Fluorobiphenyl	80.000	79.030	1.2	87	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070215\
Data File : BF080340.D
Acq On : 3 Jul 2015 12:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 03 23:40:44 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 03 22:57:58 2015
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	39.422	1.4	89	-0.03
46		2-Chloronaphthalene	40.000	35.908	10.2	79	-0.02
47		2-Nitroaniline	40.000	39.542	1.1	88	-0.03
48		Acenaphthylene	40.000	36.700	8.2	80	-0.03
49		Dimethylphthalate	40.000	38.067	4.8	81	-0.03
50		2,6-Dinitrotoluene	40.000	38.667	3.3	87	-0.03
51	C	Acenaphthene	40.000	37.126	7.2	83	-0.03
52		3-Nitroaniline	40.000	39.403	1.5	87	-0.03
53	P	2,4-Dinitrophenol	40.000	42.935	-7.3	111	-0.03
54		Dibenzofuran	40.000	37.007	7.5	85	-0.03
55	P	4-Nitrophenol	40.000	36.319	9.2	84	-0.02
56		2,4-Dinitrotoluene	40.000	36.265	9.3	76	-0.03
57		Fluorene	40.000	39.496	1.3	87	-0.03
58		2,3,4,6-Tetrachlorophenol	40.000	38.660	3.4	85	-0.02
59		Diethylphthalate	40.000	37.924	5.2	89	-0.02
60		4-Chlorophenyl-phenylether	40.000	35.376	11.6	79	-0.03
61		4-Nitroaniline	40.000	35.437	11.4	82	-0.03
62		Azobenzene	40.000	36.537	8.7	80	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	87	-0.03
64		4,6-Dinitro-2-methylphenol	40.000	39.984	0.0	94	-0.02
65	c	n-Nitrosodiphenylamine	40.000	34.436	13.9	79	-0.03
66		4-Bromophenyl-phenylether	40.000	35.590	11.0	83	-0.03
67		Hexachlorobenzene	40.000	36.802	8.0	88	-0.03
68		Atrazine	40.000	38.872	2.8	88	-0.03
69	C	Pentachlorophenol	40.000	35.001	12.5	83	-0.03
70		Phenanthrene	40.000	38.281	4.3	88	-0.03
71		Anthracene	40.000	37.486	6.3	88	-0.02
72		Carbazole	40.000	39.516	1.2	94	-0.03
73		Di-n-butylphthalate	40.000	36.072	9.8	81	-0.02
74	C	Fluoranthene	40.000	36.922	7.7	83	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	94	0.03
76		Benzidine	40.000	37.176	7.1	95	0.00
77		Pyrene	40.000	36.026	9.9	90	0.00
78	S	Terphenyl-d14	80.000	72.917	8.9	88	0.00
79		Butylbenzylphthalate	40.000	31.485	21.3	78	0.00
80		Benzo(a)anthracene	40.000	37.653	5.9	93	0.03
81		3,3'-Dichlorobenzidine	40.000	36.585	8.5	90	0.03
82		Chrysene	40.000	37.798	5.5	92	0.03
83		Bis(2-ethylhexyl)phthalate	40.000	32.412	19.0	79	0.04
84	c	Di-n-octyl phthalate	40.000	31.563	21.1#	77	0.06
85		Indeno(1,2,3-cd)pyrene	40.000	44.143	-10.4	115	0.05
86	I	Perylene-d12	20.000	20.000	0.0	101	0.06
87		Benzo(b)fluoranthene	40.000	34.935	12.7	90	0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070215\
Data File : BF080340.D
Acq On : 3 Jul 2015 12:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 03 23:40:44 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 03 22:57:58 2015
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	37.557	6.1	115	0.07
89 C	Benzo(a)pyrene	40.000	37.206	7.0	102	0.06
90	Dibenzo(a,h)anthracene	40.000	40.872	-2.2	118	0.06
91	Benzo(g,h,i)perylene	40.000	39.959	0.1	112	0.05

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