Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080317\
Data File : BF097322.D
Acq On : 3 Aug 2017 17:47
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 04 01:15:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
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	77	-0.05
2		1,4-Dioxane	40.000	47.417	-18.5	99	-0.11
3		Pyridine	40.000	35.308	11.7	62	-0.12
4		n-Nitrosodimethylamine	40.000	38.656	3.4	72	-0.11
5	S	2-Fluorophenol	80.000	84.660	-5.8	84	-0.06
6		Aniline	40.000	42.012	-5.0	80	-0.05
7	S	Phenol-d6	80.000	83.009	-3.8	79	-0.05
8		2-Chlorophenol	40.000	42.854	-7.1	83	-0.05
9		Benzaldehyde	40.000	42.734	-6.8	85	-0.05
10	C	Phenol	40.000	44.124	-10.3	83	-0.05
11		bis(2-Chloroethyl)ether	40.000	42.587	-6.5	81	-0.05
12		1,3-Dichlorobenzene	40.000	41.304	-3.3	80	-0.05
13	C	1,4-Dichlorobenzene	40.000	42.124	-5.3	86	-0.05
14		1,2-Dichlorobenzene	40.000	40.521	-1.3	82	-0.06
15		Benzyl Alcohol	40.000	43.990	-10.0	92	-0.05
16		2,2'-oxybis(1-Chloropropane	40.000	40.959	-2.4	84	-0.05
17		2-Methylphenol	40.000	41.201	-3.0	84	-0.05
18		Hexachloroethane	40.000	39.748	0.6	79	-0.06
19	P	n-Nitroso-di-n-propylamine	40.000	39.107	2.2	81	-0.05
20		3+4-Methylphenols	40.000	41.890	-4.7	86	-0.05
21	I	Naphthalene-d8	20.000	20.000	0.0	78	-0.06
22		Acetophenone	40.000	39.423	1.4	79	-0.05
23	S	Nitrobenzene-d5	80.000	78.720	1.6	80	-0.05
24		Nitrobenzene	40.000	40.544	-1.4	79	-0.05
25		Isophorone	40.000	36.825	7.9	75	-0.05
26	C	2-Nitrophenol	40.000	41.035	-2.6	80	-0.05
27		2,4-Dimethylphenol	40.000	38.556	3.6	72	-0.05
28		bis(2-Chloroethoxy)methane	40.000	37.813	5.5	73	-0.05
29	C	2,4-Dichlorophenol	40.000	41.174	-2.9	78	-0.05
30		1,2,4-Trichlorobenzene	40.000	39.903	0.2	80	-0.06
31		Naphthalene	40.000	41.647	-4.1	85	-0.06
32		Benzoic acid	40.000	40.600	-1.5	75	-0.05
33		4-Chloroaniline	40.000	40.873	-2.2	78	-0.05
34	C	Hexachlorobutadiene	40.000	39.981	0.0	79	-0.06
35		Caprolactam	40.000	44.963	-12.4	84	-0.06
36	C	4-Chloro-3-methylphenol	40.000	41.953	-4.9	80	-0.05
37		2-Methylnaphthalene	40.000	40.718	-1.8	80	-0.06
38	I	Acenaphthene-d10	20.000	20.000	0.0	79	-0.06
39		1,2,4,5-Tetrachlorobenzene	40.000	40.556	-1.4	78	-0.06
40	P	Hexachlorocyclopentadiene	40.000	47.113	-17.8	93	-0.06
41	S	2,4,6-Tribromophenol	80.000	83.011	-3.8	85	-0.06
42	C	2,4,6-Trichlorophenol	40.000	39.495	1.3	75	-0.06
43		2,4,5-Trichlorophenol	40.000	38.513	3.7	73	-0.05
44	S	2-Fluorobiphenyl	80.000	78.446	1.9	77	-0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080317\
Data File : BF097322.D
Acq On : 3 Aug 2017 17:47
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 04 01:15:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
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	39.531	1.2	77	-0.06
46		2-Chloronaphthalene	40.000	40.393	-1.0	79	-0.06
47		2-Nitroaniline	40.000	42.632	-6.6	77	-0.06
48		Acenaphthylene	40.000	38.436	3.9	80	-0.06
49		Dimethylphthalate	40.000	39.106	2.2	82	-0.06
50		2,6-Dinitrotoluene	40.000	39.685	0.8	80	-0.06
51	C	Acenaphthene	40.000	40.020	-0.1	83	-0.06
52		3-Nitroaniline	40.000	40.763	-1.9	85	-0.06
53	P	2,4-Dinitrophenol	40.000	38.060	4.8	72	-0.05
54		Dibenzofuran	40.000	41.112	-2.8	84	-0.06
55	P	4-Nitrophenol	40.000	34.491	13.8	65	-0.05
56		2,4-Dinitrotoluene	40.000	44.397	-11.0	91	-0.05
57		Fluorene	40.000	41.252	-3.1	83	-0.06
58		2,3,4,6-Tetrachlorophenol	40.000	41.332	-3.3	83	-0.06
59		Diethylphthalate	40.000	42.695	-6.7	89	-0.06
60		4-Chlorophenyl-phenylether	40.000	41.337	-3.3	83	-0.06
61		4-Nitroaniline	40.000	40.645	-1.6	80	-0.05
62		Azobenzene	40.000	43.812	-9.5	82	-0.06
63	I	Phenanthrene-d10	20.000	20.000	0.0	79	-0.06
64		4,6-Dinitro-2-methylphenol	40.000	43.177	-7.9	84	-0.05
65	c	n-Nitrosodiphenylamine	40.000	38.424	3.9	85	-0.06
66		4-Bromophenyl-phenylether	40.000	39.998	0.0	86	-0.06
67		Hexachlorobenzene	40.000	38.985	2.5	85	-0.06
68		Atrazine	40.000	37.868	5.3	84	-0.06
69	C	Pentachlorophenol	40.000	39.615	1.0	78	-0.06
70		Phenanthrene	40.000	38.093	4.8	81	-0.06
71		Anthracene	40.000	37.572	6.1	81	-0.06
72		Carbazole	40.000	40.282	-0.7	89	-0.06
73		Di-n-butylphthalate	40.000	39.998	0.0	86	-0.06
74	C	Fluoranthene	40.000	38.210	4.5	86	-0.06
75	I	Chrysene-d12	20.000	20.000	0.0	83	-0.06
76		Benzidine	40.000	25.139	37.2#	48	-0.06
77		Pyrene	40.000	39.966	0.1	88	-0.06
78	S	Terphenyl-d14	80.000	79.801	0.2	89	-0.06
79		Butylbenzylphthalate	40.000	41.254	-3.1	87	-0.07
80		Benzo(a)anthracene	40.000	41.384	-3.5	90	-0.06
81		3,3'-Dichlorobenzidine	40.000	42.527	-6.3	92	-0.06
82		Chrysene	40.000	42.682	-6.7	93	-0.06
83		Bis(2-ethylhexyl)phthalate	40.000	39.699	0.8	86	-0.07
84	c	Di-n-octyl phthalate	40.000	41.357	-3.4	88	-0.07
85		Indeno(1,2,3-cd)pyrene	40.000	35.310	11.7	81	-0.11
86	I	Perylene-d12	20.000	20.000	0.0	85	-0.08
87		Benzo(b)fluoranthene	40.000	41.929	-4.8	94	-0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080317\
Data File : BF097322.D
Acq On : 3 Aug 2017 17:47
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 04 01:15:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
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	36.848	7.9	83	-0.07
89 C	Benzo(a)pyrene	40.000	40.711	-1.8	92	-0.07
90	Dibenzo(a,h)anthracene	40.000	35.483	11.3	83	-0.11
91	Benzo(g,h,i)perylene	40.000	35.390	11.5	82	-0.12

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