Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF020819\
Data File : BF112456.D
Acq On : 8 Feb 2019 13:37
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 09 04:55:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 10:38:18 2019
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	63	-0.01
2		1,4-Dioxane	40.000	51.832	-29.6#	85	-0.03
3		Pyridine	40.000	48.159	-20.4	79	-0.02
4		n-Nitrosodimethylamine	40.000	48.923	-22.3	77	-0.02
5	S	2-Fluorophenol	80.000	97.048	-21.3	70	-0.01
6		Aniline	40.000	43.123	-7.8	66	-0.01
7	S	Phenol-d6	80.000	86.166	-7.7	66	0.00
8		2-Chlorophenol	40.000	41.499	-3.7	65	-0.01
9		Benzaldehyde	40.000	40.335	-0.8	63	0.00
10	C	Phenol	40.000	42.615	-6.5	65	-0.01
11		bis(2-Chloroethyl)ether	40.000	40.564	-1.4	63	-0.01
12		1,3-Dichlorobenzene	40.000	40.682	-1.7	66	-0.01
13	C	1,4-Dichlorobenzene	40.000	40.049	-0.1	66	-0.01
14		1,2-Dichlorobenzene	40.000	40.056	-0.1	66	-0.02
15		Benzyl Alcohol	40.000	39.800	0.5	65	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	38.660	3.4	62	-0.01
17		2-Methylphenol	40.000	39.273	1.8	64	0.00
18		Hexachloroethane	40.000	39.835	0.4	64	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	39.288	1.8	64	-0.01
20		3+4-Methylphenols	40.000	41.431	-3.6	68	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	71	-0.01
22		Acetophenone	40.000	39.924	0.2	67	-0.01
23	S	Nitrobenzene-d5	80.000	79.662	0.4	66	0.00
24		Nitrobenzene	40.000	40.328	-0.8	66	0.00
25		Isophorone	40.000	38.866	2.8	66	-0.01
26	C	2-Nitrophenol	40.000	41.117	-2.8	70	0.00
27		2,4-Dimethylphenol	40.000	40.125	-0.3	71	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.915	2.7	69	-0.01
29	C	2,4-Dichlorophenol	40.000	40.928	-2.3	73	0.00
30		1,2,4-Trichlorobenzene	40.000	39.797	0.5	72	-0.01
31		Naphthalene	40.000	39.792	0.5	73	-0.01
32		Benzoic acid	40.000	35.746	10.6	62	0.00
33		4-Chloroaniline	40.000	40.975	-2.4	75	0.00
34	C	Hexachlorobutadiene	40.000	38.240	4.4	70	-0.01
35		Caprolactam	40.000	41.821	-4.6	77	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.002	-0.0	73	0.00
37		2-Methylnaphthalene	40.000	39.396	1.5	73	-0.01
38		1-Methylnaphthalene	40.000	39.035	2.4	72	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	70	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	38.434	3.9	72	-0.01
41	P	Hexachlorocyclopentadiene	40.000	32.974	17.6	61	-0.02
42	S	2,4,6-Tribromophenol	80.000	82.491	-3.1	82	-0.01
43	C	2,4,6-Trichlorophenol	40.000	39.036	2.4	71	-0.01
44		2,4,5-Trichlorophenol	40.000	38.894	2.8	70	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF020819\
Data File : BF112456.D
Acq On : 8 Feb 2019 13:37
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 09 04:55:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 10:38:18 2019
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	S	2-Fluorobiphenyl	80.000	76.762	4.0	74	-0.01
46		1,1'-Biphenyl	40.000	38.975	2.6	73	-0.01
47		2-Chloronaphthalene	40.000	38.990	2.5	73	-0.01
48		2-Nitroaniline	40.000	40.640	-1.6	74	0.00
49		Acenaphthylene	40.000	39.710	0.7	74	-0.01
50		Dimethylphthalate	40.000	38.617	3.5	73	-0.01
51		2,6-Dinitrotoluene	40.000	40.164	-0.4	76	0.00
52	C	Acenaphthene	40.000	39.525	1.2	73	-0.01
53		3-Nitroaniline	40.000	41.486	-3.7	77	0.00
54	P	2,4-Dinitrophenol	40.000	36.349	9.1	66	0.00
55		Dibenzofuran	40.000	39.554	1.1	73	-0.01
56	P	4-Nitrophenol	40.000	33.799	15.5	60	0.00
57		2,4-Dinitrotoluene	40.000	40.976	-2.4	75	0.00
58		Fluorene	40.000	40.499	-1.2	81	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	39.297	1.8	68	0.00
60		Diethylphthalate	40.000	38.306	4.2	71	-0.02
61		4-Chlorophenyl-phenylether	40.000	38.476	3.8	73	-0.01
62		4-Nitroaniline	40.000	44.235	-10.6	83	0.00
63		Azobenzene	40.000	39.598	1.0	72	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	73	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	36.350	9.1	68	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.651	3.4	75	-0.01
67		4-Bromophenyl-phenylether	40.000	37.090	7.3	71	-0.01
68		Hexachlorobenzene	40.000	38.095	4.8	74	0.00
69		Atrazine	40.000	34.912	12.7	67	-0.01
70	C	Pentachlorophenol	40.000	36.608	8.5	72	0.00
71		Phenanthrene	40.000	38.663	3.3	74	-0.01
72		Anthracene	40.000	39.990	0.0	85	-0.01
73		Carbazole	40.000	40.777	-1.9	78	-0.01
74		Di-n-butylphthalate	40.000	36.721	8.2	71	-0.01
75	C	Fluoranthene	40.000	38.532	3.7	69	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	65	0.00
77		Benzidine	40.000	39.765	0.6	63	0.00
78		Pyrene	40.000	39.015	2.5	69	-0.01
79	S	Terphenyl-d14	80.000	74.132	7.3	68	0.00
80		Butylbenzylphthalate	40.000	35.851	10.4	63	-0.01
81		Benzo(a)anthracene	40.000	39.184	2.0	67	0.00
82		3,3'-Dichlorobenzidine	40.000	39.049	2.4	67	0.00
83		Chrysene	40.000	38.858	2.9	69	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	33.586	16.0	59	0.00
85	c	Di-n-octyl phthalate	40.000	33.409	16.5	58	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	35.139	12.2	71	0.00
87	I	Perylene-d12	20.000	20.000	0.0	64	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF020819\
Data File : BF112456.D
Acq On : 8 Feb 2019 13:37
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 09 04:55:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 10:38:18 2019
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(b)fluoranthene	40.000	39.840	0.4	65	0.00
89	Benzo(k)fluoranthene	40.000	39.939	0.2	68	0.00
90 C	Benzo(a)pyrene	40.000	39.923	0.2	67	0.00
91	Dibenzo(a,h)anthracene	40.000	38.028	4.9	71	0.00
92	Benzo(g,h,i)perylene	40.000	38.965	2.6	79	0.01

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