Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
Data File : BF114160.D
Acq On : 11 May 2019 16:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 06:52:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	111	0.00
2		1,4-Dioxane	40.000	33.893	15.3	92	0.00
3		Pyridine	40.000	33.560	16.1	95	0.00
4		n-Nitrosodimethylamine	40.000	33.617	16.0	93	0.00
5	S	2-Fluorophenol	80.000	70.487	11.9	96	0.00
6		Aniline	40.000	40.299	-0.7	110	0.00
7	S	Phenol-d6	80.000	80.212	-0.3	112	0.00
8		2-Chlorophenol	40.000	40.515	-1.3	112	0.00
9		Benzaldehyde	40.000	40.588	-1.5	106	0.00
10	C	Phenol	40.000	39.748	0.6	109	0.00
11		bis(2-Chloroethyl)ether	40.000	41.894	-4.7	116	0.00
12		1,3-Dichlorobenzene	40.000	40.605	-1.5	113	0.00
13	C	1,4-Dichlorobenzene	40.000	40.640	-1.6	113	0.00
14		1,2-Dichlorobenzene	40.000	41.708	-4.3	114	0.00
15		Benzyl Alcohol	40.000	41.684	-4.2	114	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	43.542	-8.9	120	0.00
17		2-Methylphenol	40.000	43.393	-8.5	121	0.00
18		Hexachloroethane	40.000	43.075	-7.7	119	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.305	-5.8	120	0.00
20		3+4-Methylphenols	40.000	41.854	-4.6	116	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	119	0.00
22		Acetophenone	40.000	39.678	0.8	117	0.00
23	S	Nitrobenzene-d5	80.000	81.508	-1.9	119	0.00
24		Nitrobenzene	40.000	41.542	-3.9	120	0.00
25		Isophorone	40.000	40.740	-1.9	124	0.00
26	C	2-Nitrophenol	40.000	43.090	-7.7	128	0.00
27		2,4-Dimethylphenol	40.000	40.616	-1.5	120	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.855	-4.6	125	0.00
29	C	2,4-Dichlorophenol	40.000	40.596	-1.5	118	0.00
30		1,2,4-Trichlorobenzene	40.000	40.365	-0.9	118	0.00
31		Naphthalene	40.000	40.287	-0.7	116	0.00
32		Benzoic acid	40.000	41.131	-2.8	130	0.00
33		4-Chloroaniline	40.000	41.074	-2.7	118	0.00
34	C	Hexachlorobutadiene	40.000	40.188	-0.5	116	0.00
35		Caprolactam	40.000	43.050	-7.6	122	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.154	-5.4	120	0.00
37		2-Methylnaphthalene	40.000	42.033	-5.1	120	0.00
38		1-Methylnaphthalene	40.000	41.453	-3.6	117	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	121	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.556	-1.4	118	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.453	3.9	114	0.00
42	S	2,4,6-Tribromophenol	80.000	77.097	3.6	117	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.076	-5.2	122	0.00
44		2,4,5-Trichlorophenol	40.000	40.792	-2.0	118	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
Data File : BF114160.D
Acq On : 11 May 2019 16:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 06:52:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	74.240	7.2	113	0.00
46		1,1'-Biphenyl	40.000	39.551	1.1	117	0.00
47		2-Chloronaphthalene	40.000	39.762	0.6	116	0.00
48		2-Nitroaniline	40.000	40.770	-1.9	120	0.00
49		Acenaphthylene	40.000	39.816	0.5	116	0.00
50		Dimethylphthalate	40.000	40.204	-0.5	120	0.00
51		2,6-Dinitrotoluene	40.000	40.992	-2.5	121	0.00
52	C	Acenaphthene	40.000	39.268	1.8	117	0.00
53		3-Nitroaniline	40.000	40.896	-2.2	120	0.00
54	P	2,4-Dinitrophenol	40.000	42.416	-6.0	139	0.00
55		Dibenzofuran	40.000	38.583	3.5	116	0.00
56	P	4-Nitrophenol	40.000	39.472	1.3	121	0.00
57		2,4-Dinitrotoluene	40.000	39.486	1.3	119	0.00
58		Fluorene	40.000	37.927	5.2	115	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.389	1.5	118	0.00
60		Diethylphthalate	40.000	38.564	3.6	118	0.00
61		4-Chlorophenyl-phenylether	40.000	38.355	4.1	116	0.00
62		4-Nitroaniline	40.000	39.338	1.7	122	0.00
63		Azobenzene	40.000	39.187	2.0	119	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	40.000	45.399	-13.5	130	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.835	-2.1	116	0.00
67		4-Bromophenyl-phenylether	40.000	41.067	-2.7	118	0.00
68		Hexachlorobenzene	40.000	40.119	-0.3	116	0.00
69		Atrazine	40.000	42.119	-5.3	123	0.00
70	C	Pentachlorophenol	40.000	42.725	-6.8	120	0.00
71		Phenanthrene	40.000	39.917	0.2	114	0.00
72		Anthracene	40.000	40.412	-1.0	114	0.00
73		Carbazole	40.000	40.012	-0.0	114	0.00
74		Di-n-butylphthalate	40.000	42.034	-5.1	119	0.00
75	C	Fluoranthene	40.000	39.081	2.3	112	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	113	-0.01
77		Benzidine	40.000	38.725	3.2	112	0.00
78		Pyrene	40.000	38.049	4.9	111	0.00
79	S	Terphenyl-d14	80.000	74.610	6.7	110	0.00
80		Butylbenzylphthalate	40.000	42.578	-6.4	120	0.00
81		Benzo(a)anthracene	40.000	39.982	0.0	110	-0.01
82		3,3'-Dichlorobenzidine	40.000	43.385	-8.5	115	-0.01
83		Chrysene	40.000	40.397	-1.0	112	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.128	-5.3	118	0.00
85	c	Di-n-octyl phthalate	40.000	41.776	-4.4	114	-0.03
86		Indeno(1,2,3-cd)pyrene	40.000	39.503	1.2	115	-0.03
87	I	Perylene-d12	20.000	20.000	0.0	111	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051119\
Data File : BF114160.D
Acq On : 11 May 2019 16:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 13 06:52:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	42.837	-7.1	119	-0.03
89	Benzo(k)fluoranthene	40.000	38.220	4.5	100	-0.03
90 C	Benzo(a)pyrene	40.000	40.834	-2.1	111	-0.03
91	Dibenzo(a,h)anthracene	40.000	40.568	-1.4	115	-0.03
92	Benzo(g,h,i)perylene	40.000	37.940	5.2	110	-0.04

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