Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052019\
Data File : BF114428.D
Acq On : 20 May 2019 10:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 10:49:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 20 04:47:40 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	69	-0.01
2		1,4-Dioxane	40.000	34.505	13.7	58	0.01
3		Pyridine	40.000	35.859	10.4	63	0.00
4		n-Nitrosodimethylamine	40.000	35.501	11.2	61	0.00
5	S	2-Fluorophenol	80.000	77.046	3.7	65	0.00
6		Aniline	40.000	39.429	1.4	67	0.00
7	S	Phenol-d6	80.000	81.798	-2.2	70	0.00
8		2-Chlorophenol	40.000	41.453	-3.6	71	-0.01
9		Benzaldehyde	40.000	33.487	16.3	54	-0.01
10	C	Phenol	40.000	39.814	0.5	67	-0.01
11		bis(2-Chloroethyl)ether	40.000	41.535	-3.8	71	0.00
12		1,3-Dichlorobenzene	40.000	40.554	-1.4	70	0.00
13	C	1,4-Dichlorobenzene	40.000	40.608	-1.5	70	-0.01
14		1,2-Dichlorobenzene	40.000	40.882	-2.2	69	0.00
15		Benzyl Alcohol	40.000	38.707	3.2	65	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.955	2.6	66	-0.01
17		2-Methylphenol	40.000	40.744	-1.9	70	-0.01
18		Hexachloroethane	40.000	40.283	-0.7	69	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.208	2.0	69	-0.01
20		3+4-Methylphenols	40.000	42.626	-6.6	73	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	72	0.00
22		Acetophenone	40.000	40.805	-2.0	72	-0.01
23	S	Nitrobenzene-d5	80.000	80.477	-0.6	71	0.00
24		Nitrobenzene	40.000	40.280	-0.7	70	-0.01
25		Isophorone	40.000	39.360	1.6	72	0.00
26	C	2-Nitrophenol	40.000	41.360	-3.4	74	0.00
27		2,4-Dimethylphenol	40.000	39.316	1.7	70	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.974	0.1	71	0.00
29	C	2,4-Dichlorophenol	40.000	40.480	-1.2	71	0.00
30		1,2,4-Trichlorobenzene	40.000	39.592	1.0	70	0.00
31		Naphthalene	40.000	41.216	-3.0	71	-0.01
32		Benzoic acid	40.000	40.459	-1.1	77	-0.01
33		4-Chloroaniline	40.000	41.483	-3.7	72	0.00
34	C	Hexachlorobutadiene	40.000	40.196	-0.5	70	0.00
35		Caprolactam	40.000	43.920	-9.8	74	-0.02
36	C	4-Chloro-3-methylphenol	40.000	42.581	-6.5	73	0.00
37		2-Methylnaphthalene	40.000	41.370	-3.4	71	0.00
38		1-Methylnaphthalene	40.000	41.534	-3.8	71	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	73	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.663	-1.7	71	0.00
41	P	Hexachlorocyclopentadiene	40.000	35.819	10.5	63	0.00
42	S	2,4,6-Tribromophenol	80.000	74.045	7.4	68	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.570	1.1	69	0.00
44		2,4,5-Trichlorophenol	40.000	39.628	0.9	69	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052019\
Data File : BF114428.D
Acq On : 20 May 2019 10:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 10:49:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 20 04:47:40 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	77.011	3.7	71	0.00
46		1,1'-Biphenyl	40.000	40.048	-0.1	71	0.00
47		2-Chloronaphthalene	40.000	39.894	0.3	70	0.00
48		2-Nitroaniline	40.000	41.085	-2.7	73	0.00
49		Acenaphthylene	40.000	39.954	0.1	70	0.00
50		Dimethylphthalate	40.000	40.154	-0.4	72	0.00
51		2,6-Dinitrotoluene	40.000	40.377	-0.9	72	0.00
52	C	Acenaphthene	40.000	38.741	3.1	69	0.00
53		3-Nitroaniline	40.000	41.696	-4.2	74	0.00
54	P	2,4-Dinitrophenol	40.000	35.638	10.9	67	0.00
55		Dibenzofuran	40.000	37.913	5.2	68	0.00
56	P	4-Nitrophenol	40.000	37.690	5.8	69	0.00
57		2,4-Dinitrotoluene	40.000	38.137	4.7	69	0.00
58		Fluorene	40.000	39.984	0.0	72	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.157	4.6	69	0.00
60		Diethylphthalate	40.000	38.893	2.8	72	0.00
61		4-Chlorophenyl-phenylether	40.000	39.974	0.1	72	0.00
62		4-Nitroaniline	40.000	41.340	-3.4	77	0.00
63		Azobenzene	40.000	39.176	2.1	72	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	73	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.618	-4.0	74	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.324	-0.8	71	0.00
67		4-Bromophenyl-phenylether	40.000	38.902	2.7	70	0.00
68		Hexachlorobenzene	40.000	38.840	2.9	69	0.00
69		Atrazine	40.000	39.172	2.1	71	0.00
70	C	Pentachlorophenol	40.000	34.378	14.1	60	0.00
71		Phenanthrene	40.000	40.556	-1.4	72	0.00
72		Anthracene	40.000	40.881	-2.2	72	0.00
73		Carbazole	40.000	41.607	-4.0	73	0.00
74		Di-n-butylphthalate	40.000	41.244	-3.1	72	0.00
75	C	Fluoranthene	40.000	41.103	-2.8	73	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	73	0.00
77		Benzidine	40.000	33.410	16.5	63	0.00
78		Pyrene	40.000	37.993	5.0	73	0.00
79	S	Terphenyl-d14	80.000	74.653	6.7	72	0.00
80		Butylbenzylphthalate	40.000	40.669	-1.7	75	0.00
81		Benzo(a)anthracene	40.000	42.008	-5.0	76	0.00
82		3,3'-Dichlorobenzidine	40.000	42.072	-5.2	73	0.00
83		Chrysene	40.000	40.920	-2.3	74	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.852	-4.6	76	0.00
85	c	Di-n-octyl phthalate	40.000	42.680	-6.7	76	0.02
86		Indeno(1,2,3-cd)pyrene	40.000	40.683	-1.7	77	0.02
87	I	Perylene-d12	20.000	20.000	0.0	76	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052019\
Data File : BF114428.D
Acq On : 20 May 2019 10:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 10:49:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 20 04:47:40 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	41.521	-3.8	79	0.02
89	Benzo(k)fluoranthene	40.000	39.702	0.7	72	0.02
90 C	Benzo(a)pyrene	40.000	40.947	-2.4	77	0.02
91	Dibenzo(a,h)anthracene	40.000	39.952	0.1	78	0.02
92	Benzo(g,h,i)perylene	40.000	40.012	-0.0	79	0.02

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