Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
Data File : BF112954.D
Acq On : 11 Mar 2019 14:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:47:45 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 01 15:40:43 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	108	0.00
2		1,4-Dioxane	40.000	34.344	14.1	93	0.00
3		Pyridine	40.000	33.263	16.8	91	0.00
4		n-Nitrosodimethylamine	40.000	34.473	13.8	92	0.00
5	S	2-Fluorophenol	80.000	70.782	11.5	97	0.00
6		Aniline	40.000	35.510	11.2	96	0.00
7	S	Phenol-d6	80.000	74.184	7.3	103	0.00
8		2-Chlorophenol	40.000	38.301	4.2	105	0.00
9		Benzaldehyde	40.000	33.173	17.1	91	0.00
10	C	Phenol	40.000	36.222	9.4	97	0.00
11		bis(2-Chloroethyl)ether	40.000	36.927	7.7	101	0.00
12		1,3-Dichlorobenzene	40.000	39.236	1.9	108	0.00
13	C	1,4-Dichlorobenzene	40.000	39.388	1.5	108	0.00
14		1,2-Dichlorobenzene	40.000	39.132	2.2	109	0.00
15		Benzyl Alcohol	40.000	38.085	4.8	103	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.031	9.9	99	0.00
17		2-Methylphenol	40.000	38.433	3.9	106	0.00
18		Hexachloroethane	40.000	38.330	4.2	105	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.686	8.3	102	0.00
20		3+4-Methylphenols	40.000	38.455	3.9	108	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	110	0.00
22		Acetophenone	40.000	39.669	0.8	108	0.00
23	S	Nitrobenzene-d5	80.000	82.274	-2.8	111	0.00
24		Nitrobenzene	40.000	39.757	0.6	108	0.00
25		Isophorone	40.000	37.794	5.5	106	0.00
26	C	2-Nitrophenol	40.000	49.588	-24.0#	128	0.00
27		2,4-Dimethylphenol	40.000	37.817	5.5	106	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.014	5.0	107	0.00
29	C	2,4-Dichlorophenol	40.000	41.467	-3.7	113	0.00
30		1,2,4-Trichlorobenzene	40.000	40.892	-2.2	114	0.00
31		Naphthalene	40.000	38.369	4.1	108	0.00
32		Benzoic acid	40.000	37.171	7.1	101	0.02
33		4-Chloroaniline	40.000	39.466	1.3	109	0.00
34	C	Hexachlorobutadiene	40.000	43.535	-8.8	121	-0.01
35		Caprolactam	40.000	40.931	-2.3	111	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.843	0.4	110	0.00
37		2-Methylnaphthalene	40.000	39.733	0.7	111	0.00
38		1-Methylnaphthalene	40.000	39.049	2.4	110	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	119	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.791	-2.0	122	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.040	-0.1	117	-0.01
42	S	2,4,6-Tribromophenol	80.000	90.779	-13.5	133	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.155	-2.9	120	0.00
44		2,4,5-Trichlorophenol	40.000	41.738	-4.3	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
Data File : BF112954.D
Acq On : 11 Mar 2019 14:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:47:45 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 01 15:40:43 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	76.797	4.0	114	0.00
46		1,1'-Biphenyl	40.000	38.363	4.1	112	0.00
47		2-Chloronaphthalene	40.000	38.388	4.0	116	0.00
48		2-Nitroaniline	40.000	42.950	-7.4	119	0.00
49		Acenaphthylene	40.000	37.081	7.3	113	0.00
50		Dimethylphthalate	40.000	39.745	0.6	120	0.00
51		2,6-Dinitrotoluene	40.000	44.202	-10.5	123	0.00
52	C	Acenaphthene	40.000	35.951	10.1	108	0.00
53		3-Nitroaniline	40.000	41.640	-4.1	117	0.00
54	P	2,4-Dinitrophenol	40.000	54.345	-35.9#	174	0.00
55		Dibenzofuran	40.000	36.829	7.9	113	0.00
56	P	4-Nitrophenol	40.000	41.249	-3.1	113	0.02
57		2,4-Dinitrotoluene	40.000	45.868	-14.7	126	0.00
58		Fluorene	40.000	38.359	4.1	117	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	43.109	-7.8	125	0.00
60		Diethylphthalate	40.000	36.580	8.6	111	0.00
61		4-Chlorophenyl-phenylether	40.000	40.678	-1.7	124	0.00
62		4-Nitroaniline	40.000	44.927	-12.3	130	0.00
63		Azobenzene	40.000	34.694	13.3	105	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	123	0.00
65		4,6-Dinitro-2-methylphenol	40.000	48.703	-21.8	155	0.01
66	c	n-Nitrosodiphenylamine	40.000	36.726	8.2	113	0.00
67		4-Bromophenyl-phenylether	40.000	40.596	-1.5	126	0.00
68		Hexachlorobenzene	40.000	41.569	-3.9	129	0.00
69		Atrazine	40.000	40.012	-0.0	124	0.00
70	C	Pentachlorophenol	40.000	43.996	-10.0	129	0.00
71		Phenanthrene	40.000	38.499	3.8	117	0.00
72		Anthracene	40.000	37.316	6.7	117	0.00
73		Carbazole	40.000	37.136	7.2	117	0.00
74		Di-n-butylphthalate	40.000	36.268	9.3	115	-0.01
75	C	Fluoranthene	40.000	38.721	3.2	117	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	133	0.00
77		Benzidine	40.000	32.773	18.1	106	0.00
78		Pyrene	40.000	35.771	10.6	118	0.00
79	S	Terphenyl-d14	80.000	71.251	10.9	121	0.00
80		Butylbenzylphthalate	40.000	36.288	9.3	117	0.00
81		Benzo(a)anthracene	40.000	33.587	16.0	110	0.00
82		3,3'-Dichlorobenzidine	40.000	37.503	6.2	120	0.00
83		Chrysene	40.000	35.163	12.1	117	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	34.129	14.7	113	-0.01
85	c	Di-n-octyl phthalate	40.000	35.180	12.1	116	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	33.531	16.2	118	0.02
87	I	Perylene-d12	20.000	20.000	0.0	117	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
Data File : BF112954.D
Acq On : 11 Mar 2019 14:57
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 12 04:47:45 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 01 15:40:43 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	41.088	-2.7	112	0.00
89	Benzo(k)fluoranthene	40.000	36.605	8.5	113	0.00
90 C	Benzo(a)pyrene	40.000	38.180	4.6	111	0.00
91	Dibenzo(a,h)anthracene	40.000	39.225	1.9	119	0.00
92	Benzo(g,h,i)perylene	40.000	40.044	-0.1	121	0.01

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