Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF040320\
Data File : BF119708.D
Acq On : 3 Apr 2020 10:12
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 03 11:15:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 01 11:59:45 2020
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	81	-0.01
2		1,4-Dioxane	40.000	36.858	7.9	72	0.00
3		Pyridine	40.000	36.988	7.5	72	0.00
4		n-Nitrosodimethylamine	40.000	33.647	15.9	66	0.00
5	S	2-Fluorophenol	80.000	78.830	1.5	76	0.00
6		Aniline	40.000	38.161	4.6	73	-0.01
7	S	Phenol-d6	80.000	79.058	1.2	76	0.00
8		2-Chlorophenol	40.000	40.118	-0.3	77	0.00
9		Benzaldehyde	40.000	33.761	15.6	66	-0.01
10	C	Phenol	40.000	41.968	-4.9	81	0.00
11		bis(2-Chloroethyl)ether	40.000	38.799	3.0	75	0.00
12		1,3-Dichlorobenzene	40.000	39.742	0.6	77	-0.01
13	C	1,4-Dichlorobenzene	40.000	39.952	0.1	78	-0.01
14		1,2-Dichlorobenzene	40.000	40.344	-0.9	77	0.00
15		Benzyl Alcohol	40.000	38.378	4.1	73	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	34.274	14.3	67	0.00
17		2-Methylphenol	40.000	39.367	1.6	76	0.00
18		Hexachloroethane	40.000	40.197	-0.5	79	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.663	5.8	73	-0.01
20		3+4-Methylphenols	40.000	38.966	2.6	75	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	80	-0.01
22		Acetophenone	40.000	38.166	4.6	74	0.00
23	S	Nitrobenzene-d5	80.000	81.636	-2.0	78	0.00
24		Nitrobenzene	40.000	39.596	1.0	77	-0.01
25		Isophorone	40.000	37.853	5.4	74	0.00
26	C	2-Nitrophenol	40.000	49.629	-24.1#	95	0.00
27		2,4-Dimethylphenol	40.000	36.981	7.5	72	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.004	2.5	76	0.00
29	C	2,4-Dichlorophenol	40.000	40.085	-0.2	78	0.00
30		1,2,4-Trichlorobenzene	40.000	40.575	-1.4	79	-0.01
31		Naphthalene	40.000	40.758	-1.9	78	-0.01
32		Benzoic acid	40.000	38.309	4.2	74	-0.02
33		4-Chloroaniline	40.000	38.948	2.6	75	-0.01
34	C	Hexachlorobutadiene	40.000	41.258	-3.1	81	-0.01
35		Caprolactam	40.000	39.209	2.0	74	-0.01
36	C	4-Chloro-3-methylphenol	40.000	39.739	0.7	77	0.00
37		2-Methylnaphthalene	40.000	40.023	-0.1	78	0.00
38		1-Methylnaphthalene	40.000	40.362	-0.9	77	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	81	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	40.944	-2.4	80	-0.01
41	P	Hexachlorocyclopentadiene	40.000	40.264	-0.7	79	-0.01
42	S	2,4,6-Tribromophenol	80.000	89.575	-12.0	88	-0.01
43	C	2,4,6-Trichlorophenol	40.000	40.896	-2.2	81	0.00
44		2,4,5-Trichlorophenol	40.000	40.298	-0.7	79	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF040320\
Data File : BF119708.D
Acq On : 3 Apr 2020 10:12
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 03 11:15:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 01 11:59:45 2020
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	77.424	3.2	79	0.00
46		1,1'-Biphenyl	40.000	39.965	0.1	78	-0.01
47		2-Chloronaphthalene	40.000	39.810	0.5	78	-0.01
48		2-Nitroaniline	40.000	41.323	-3.3	79	0.00
49		Acenaphthylene	40.000	39.532	1.2	77	0.00
50		Dimethylphthalate	40.000	40.280	-0.7	79	-0.01
51		2,6-Dinitrotoluene	40.000	43.269	-8.2	83	-0.01
52	C	Acenaphthene	40.000	40.654	-1.6	80	-0.01
53		3-Nitroaniline	40.000	41.308	-3.3	80	-0.01
54	P	2,4-Dinitrophenol	40.000	56.555	-41.4#	131	0.00
55		Dibenzofuran	40.000	40.052	-0.1	78	-0.01
56	P	4-Nitrophenol	40.000	42.293	-5.7	80	0.00
57		2,4-Dinitrotoluene	40.000	45.571	-13.9	88	-0.01
58		Fluorene	40.000	41.032	-2.6	80	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.234	-3.1	79	-0.01
60		Diethylphthalate	40.000	40.473	-1.2	79	-0.01
61		4-Chlorophenyl-phenylether	40.000	41.464	-3.7	81	-0.01
62		4-Nitroaniline	40.000	44.060	-10.2	84	-0.02
63		Azobenzene	40.000	38.219	4.5	75	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	82	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	59.009	-47.5#	117	-0.01
66	c	n-Nitrosodiphenylamine	40.000	39.344	1.6	79	-0.01
67		4-Bromophenyl-phenylether	40.000	40.021	-0.1	79	0.00
68		Hexachlorobenzene	40.000	40.692	-1.7	81	-0.01
69		Atrazine	40.000	39.821	0.4	80	-0.01
70	C	Pentachlorophenol	40.000	42.212	-5.5	84	-0.01
71		Phenanthrene	40.000	39.643	0.9	79	-0.01
72		Anthracene	40.000	39.848	0.4	79	-0.01
73		Carbazole	40.000	39.566	1.1	78	-0.01
74		Di-n-butylphthalate	40.000	41.593	-4.0	83	-0.01
75	C	Fluoranthene	40.000	40.286	-0.7	79	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	86	-0.02
77		Benzidine	40.000	35.151	12.1	75	-0.01
78		Pyrene	40.000	36.885	7.8	80	-0.01
79	S	Terphenyl-d14	80.000	75.867	5.2	83	-0.01
80		Butylbenzylphthalate	40.000	40.055	-0.1	87	-0.01
81		Benzo(a)anthracene	40.000	39.013	2.5	82	-0.01
82		3,3'-Dichlorobenzidine	40.000	38.434	3.9	79	-0.02
83		Chrysene	40.000	37.403	6.5	78	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	41.575	-3.9	90	-0.01
85	c	Di-n-octyl phthalate	40.000	42.773	-6.9	89	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	37.337	6.7	84	-0.03
87	I	Perylene-d12	20.000	20.000	0.0	86	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF040320\
Data File : BF119708.D
Acq On : 3 Apr 2020 10:12
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 03 11:15:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 01 11:59:45 2020
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	38.994	2.5	83	-0.01
89	Benzo(k)fluoranthene	40.000	38.959	2.6	80	-0.02
90 C	Benzo(a)pyrene	40.000	38.798	3.0	81	-0.02
91	Dibenzo(a,h)anthracene	40.000	37.132	7.2	83	-0.03
92	Benzo(g,h,i)perylene	40.000	36.087	9.8	82	-0.04

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