Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022720\
Data File : BF119101.D
Acq On : 27 Feb 2020 9:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 28 02:51:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 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	127	0.00
2		1,4-Dioxane	40.000	37.119	7.2	125	0.02
3		Pyridine	40.000	38.359	4.1	129	0.02
4		n-Nitrosodimethylamine	40.000	43.641	-9.1	147	0.02
5	S	2-Fluorophenol	80.000	78.835	1.5	130	0.00
6		Aniline	40.000	39.328	1.7	128	0.00
7	S	Phenol-d6	80.000	80.851	-1.1	133	0.00
8		2-Chlorophenol	40.000	40.763	-1.9	134	0.00
9		Benzaldehyde	40.000	34.359	14.1	114	0.00
10	C	Phenol	40.000	39.733	0.7	131	0.00
11		bis(2-Chloroethyl)ether	40.000	40.242	-0.6	134	0.00
12		1,3-Dichlorobenzene	40.000	39.866	0.3	133	0.00
13	C	1,4-Dichlorobenzene	40.000	39.963	0.1	132	0.00
14		1,2-Dichlorobenzene	40.000	39.261	1.8	129	0.00
15		Benzyl Alcohol	40.000	41.360	-3.4	135	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.927	2.7	131	0.00
17		2-Methylphenol	40.000	40.452	-1.1	135	0.00
18		Hexachloroethane	40.000	40.300	-0.7	135	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.435	-1.1	136	0.00
20		3+4-Methylphenols	40.000	39.496	1.3	134	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	128	0.00
22		Acetophenone	40.000	40.470	-1.2	137	0.00
23	S	Nitrobenzene-d5	80.000	80.571	-0.7	136	0.00
24		Nitrobenzene	40.000	40.264	-0.7	137	0.00
25		Isophorone	40.000	39.641	0.9	136	0.00
26	C	2-Nitrophenol	40.000	40.653	-1.6	138	0.00
27		2,4-Dimethylphenol	40.000	39.501	1.2	134	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.489	3.8	131	0.00
29	C	2,4-Dichlorophenol	40.000	39.534	1.2	132	0.00
30		1,2,4-Trichlorobenzene	40.000	38.809	3.0	130	0.00
31		Naphthalene	40.000	38.655	3.4	129	0.00
32		Benzoic acid	40.000	46.290	-15.7	169	0.00
33		4-Chloroaniline	40.000	39.312	1.7	132	0.00
34	C	Hexachlorobutadiene	40.000	38.510	3.7	131	0.00
35		Caprolactam	40.000	40.953	-2.4	136	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.843	-2.1	138	0.00
37		2-Methylnaphthalene	40.000	38.420	3.9	130	0.00
38		1-Methylnaphthalene	40.000	38.559	3.6	129	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	128	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.219	4.5	130	0.00
41	P	Hexachlorocyclopentadiene	40.000	35.399	11.5	125	0.00
42	S	2,4,6-Tribromophenol	80.000	78.931	1.3	134	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.727	0.7	135	0.00
44		2,4,5-Trichlorophenol	40.000	38.480	3.8	130	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022720\
Data File : BF119101.D
Acq On : 27 Feb 2020 9:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 28 02:51:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 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	72.230	9.7	126	0.00
46		1,1'-Biphenyl	40.000	37.794	5.5	127	0.00
47		2-Chloronaphthalene	40.000	39.013	2.5	133	0.00
48		2-Nitroaniline	40.000	42.240	-5.6	143	0.00
49		Acenaphthylene	40.000	38.076	4.8	128	0.00
50		Dimethylphthalate	40.000	39.442	1.4	136	0.00
51		2,6-Dinitrotoluene	40.000	39.801	0.5	135	0.00
52	C	Acenaphthene	40.000	38.547	3.6	130	0.00
53		3-Nitroaniline	40.000	40.565	-1.4	136	0.00
54	P	2,4-Dinitrophenol	40.000	43.457	-8.6	156	0.00
55		Dibenzofuran	40.000	37.193	7.0	123	0.00
56	P	4-Nitrophenol	40.000	37.491	6.3	124	0.00
57		2,4-Dinitrotoluene	40.000	38.354	4.1	127	0.00
58		Fluorene	40.000	38.736	3.2	130	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.942	2.6	134	0.00
60		Diethylphthalate	40.000	39.189	2.0	133	0.00
61		4-Chlorophenyl-phenylether	40.000	38.216	4.5	128	0.00
62		4-Nitroaniline	40.000	40.929	-2.3	138	0.00
63		Azobenzene	40.000	39.435	1.4	134	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	128	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.878	-7.2	146	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.865	2.8	132	0.00
67		4-Bromophenyl-phenylether	40.000	38.337	4.2	132	0.00
68		Hexachlorobenzene	40.000	38.769	3.1	133	0.00
69		Atrazine	40.000	36.179	9.6	124	0.00
70	C	Pentachlorophenol	40.000	39.671	0.8	137	0.00
71		Phenanthrene	40.000	38.567	3.6	131	0.00
72		Anthracene	40.000	38.626	3.4	129	0.00
73		Carbazole	40.000	39.448	1.4	133	0.00
74		Di-n-butylphthalate	40.000	38.856	2.9	132	0.00
75	C	Fluoranthene	40.000	38.310	4.2	128	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	116	0.00
77		Benzidine	40.000	30.883	22.8	93	0.00
78		Pyrene	40.000	40.396	-1.0	128	0.00
79	S	Terphenyl-d14	80.000	80.058	-0.1	128	0.00
80		Butylbenzylphthalate	40.000	42.070	-5.2	131	0.00
81		Benzo(a)anthracene	40.000	39.969	0.1	123	0.00
82		3,3'-Dichlorobenzidine	40.000	38.871	2.8	114	0.00
83		Chrysene	40.000	38.952	2.6	121	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.733	-4.3	128	0.00
85	c	Di-n-octyl phthalate	40.000	38.674	3.3	113	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	35.785	10.5	85	0.00
87	I	Perylene-d12	20.000	20.000	0.0	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022720\
Data File : BF119101.D
Acq On : 27 Feb 2020 9:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 28 02:51:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 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	41.350	-3.4	113	0.00
89	Benzo(k)fluoranthene	40.000	39.222	1.9	98	0.00
90 C	Benzo(a)pyrene	40.000	38.863	2.8	76	0.00
91	Dibenzo(a,h)anthracene	40.000	40.035	-0.1	83	0.00
92	Benzo(g,h,i)perylene	40.000	40.977	-2.4	86	0.01

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