Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071219\
Data File : BF115525.D
Acq On : 12 Jul 2019 13:53
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF071219

Quant Time: Jul 12 14:18:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 14:03:52 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	104	0.00
2		1,4-Dioxane	40.000	37.507	6.2	101	0.00
3		Pyridine	40.000	38.630	3.4	105	0.00
4		n-Nitrosodimethylamine	40.000	38.314	4.2	102	0.00
5	S	2-Fluorophenol	80.000	77.578	3.0	103	0.00
6		Aniline	40.000	39.229	1.9	104	0.00
7	S	Phenol-d6	80.000	78.326	2.1	105	0.00
8		2-Chlorophenol	40.000	39.415	1.5	104	0.00
9		Benzaldehyde	40.000	36.322	9.2	104	0.00
10	C	Phenol	40.000	39.737	0.7	105	0.00
11		bis(2-Chloroethyl)ether	40.000	38.767	3.1	104	0.00
12		1,3-Dichlorobenzene	40.000	38.669	3.3	103	0.00
13	C	1,4-Dichlorobenzene	40.000	39.387	1.5	104	0.00
14		1,2-Dichlorobenzene	40.000	39.355	1.6	103	0.00
15		Benzyl Alcohol	40.000	40.165	-0.4	106	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.381	1.5	104	0.00
17		2-Methylphenol	40.000	39.385	1.5	106	0.00
18		Hexachloroethane	40.000	39.495	1.3	104	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.760	3.1	105	0.00
20		3+4-Methylphenols	40.000	38.278	4.3	104	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	38.970	2.6	105	0.00
23	S	Nitrobenzene-d5	80.000	78.641	1.7	106	0.00
24		Nitrobenzene	40.000	39.757	0.6	108	0.00
25		Isophorone	40.000	38.543	3.6	107	0.00
26	C	2-Nitrophenol	40.000	40.185	-0.5	108	0.00
27		2,4-Dimethylphenol	40.000	39.431	1.4	106	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.120	2.2	105	0.00
29	C	2,4-Dichlorophenol	40.000	39.333	1.7	106	0.00
30		1,2,4-Trichlorobenzene	40.000	38.903	2.7	104	0.00
31		Naphthalene	40.000	39.016	2.5	104	0.00
32		Benzoic acid	40.000	41.893	-4.7	134	0.00
33		4-Chloroaniline	40.000	38.551	3.6	106	0.00
34	C	Hexachlorobutadiene	40.000	39.041	2.4	105	0.00
35		Caprolactam	40.000	38.828	2.9	108	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.646	3.4	106	0.00
37		2-Methylnaphthalene	40.000	38.653	3.4	105	0.00
38		1-Methylnaphthalene	40.000	38.975	2.6	106	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.350	-0.9	107	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.195	2.0	101	0.00
42	S	2,4,6-Tribromophenol	80.000	79.212	1.0	108	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.804	0.5	106	0.00
44		2,4,5-Trichlorophenol	40.000	40.336	-0.8	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071219\
Data File : BF115525.D
Acq On : 12 Jul 2019 13:53
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF071219

Quant Time: Jul 12 14:18:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 14:03:52 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	82.458	-3.1	104	0.00
46		1,1'-Biphenyl	40.000	39.864	0.3	106	0.00
47		2-Chloronaphthalene	40.000	39.745	0.6	105	0.00
48		2-Nitroaniline	40.000	40.087	-0.2	109	0.00
49		Acenaphthylene	40.000	39.439	1.4	105	0.00
50		Dimethylphthalate	40.000	38.890	2.8	106	0.00
51		2,6-Dinitrotoluene	40.000	40.402	-1.0	108	0.00
52	C	Acenaphthene	40.000	39.923	0.2	107	0.00
53		3-Nitroaniline	40.000	39.764	0.6	106	0.00
54	P	2,4-Dinitrophenol	40.000	41.841	-4.6	110	0.00
55		Dibenzofuran	40.000	38.061	4.8	105	0.00
56	P	4-Nitrophenol	40.000	40.596	-1.5	108	0.00
57		2,4-Dinitrotoluene	40.000	40.439	-1.1	108	0.00
58		Fluorene	40.000	39.185	2.0	105	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.925	0.2	107	0.00
60		Diethylphthalate	40.000	39.581	1.0	107	0.00
61		4-Chlorophenyl-phenylether	40.000	36.862	7.8	105	0.00
62		4-Nitroaniline	40.000	39.160	2.1	105	0.00
63		Azobenzene	40.000	39.535	1.2	106	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.751	-4.4	110	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.466	1.3	105	0.00
67		4-Bromophenyl-phenylether	40.000	39.628	0.9	105	0.00
68		Hexachlorobenzene	40.000	39.400	1.5	106	0.00
69		Atrazine	40.000	37.302	6.7	105	0.00
70	C	Pentachlorophenol	40.000	40.530	-1.3	106	0.00
71		Phenanthrene	40.000	39.016	2.5	104	0.00
72		Anthracene	40.000	38.198	4.5	105	0.00
73		Carbazole	40.000	39.000	2.5	105	0.00
74		Di-n-butylphthalate	40.000	37.813	5.5	104	0.00
75	C	Fluoranthene	40.000	37.759	5.6	104	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	111	0.00
77		Benzidine	40.000	33.863	15.3	96	0.00
78		Pyrene	40.000	37.297	6.8	106	0.00
79	S	Terphenyl-d14	80.000	73.097	8.6	106	0.00
80		Butylbenzylphthalate	40.000	38.167	4.6	108	0.00
81		Benzo(a)anthracene	40.000	38.838	2.9	111	0.00
82		3,3'-Dichlorobenzidine	40.000	39.770	0.6	109	0.00
83		Chrysene	40.000	38.243	4.4	109	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	39.184	2.0	110	0.00
85	c	Di-n-octyl phthalate	40.000	41.295	-3.2	117	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	37.611	6.0	112	0.00
87	I	Perylene-d12	20.000	20.000	0.0	114	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071219\
Data File : BF115525.D
Acq On : 12 Jul 2019 13:53
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF071219

Quant Time: Jul 12 14:18:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 14:03:52 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	41.854	-4.6	128	0.00
89	Benzo(k)fluoranthene	40.000	37.437	6.4	105	0.00
90 C	Benzo(a)pyrene	40.000	39.701	0.7	115	0.00
91	Dibenzo(a,h)anthracene	40.000	39.034	2.4	113	0.00
92	Benzo(g,h,i)perylene	40.000	38.033	4.9	111	0.00

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