Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114122.D
Acq On : 10 May 2019 15:56
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF051019

Quant Time: May 10 18:34:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	98	0.00
2		1,4-Dioxane	40.000	38.347	4.1	92	0.00
3		Pyridine	40.000	38.296	4.3	95	0.00
4		n-Nitrosodimethylamine	40.000	39.158	2.1	95	-0.01
5	S	2-Fluorophenol	80.000	75.078	6.2	90	0.00
6		Aniline	40.000	42.896	-7.2	103	0.00
7	S	Phenol-d6	80.000	81.615	-2.0	100	0.00
8		2-Chlorophenol	40.000	41.087	-2.7	100	0.00
9		Benzaldehyde	40.000	40.406	-1.0	93	0.00
10	C	Phenol	40.000	42.521	-6.3	102	0.00
11		bis(2-Chloroethyl)ether	40.000	41.545	-3.9	101	0.00
12		1,3-Dichlorobenzene	40.000	41.702	-4.3	102	0.00
13	C	1,4-Dichlorobenzene	40.000	40.718	-1.8	100	0.00
14		1,2-Dichlorobenzene	40.000	41.176	-2.9	99	0.00
15		Benzyl Alcohol	40.000	41.187	-3.0	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.734	-1.8	99	0.00
17		2-Methylphenol	40.000	40.712	-1.8	100	0.00
18		Hexachloroethane	40.000	40.647	-1.6	98	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.797	0.5	100	0.00
20		3+4-Methylphenols	40.000	41.155	-2.9	100	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	99	0.00
22		Acetophenone	40.000	40.858	-2.1	100	0.00
23	S	Nitrobenzene-d5	80.000	81.724	-2.2	99	0.00
24		Nitrobenzene	40.000	41.704	-4.3	100	0.00
25		Isophorone	40.000	40.008	-0.0	101	0.00
26	C	2-Nitrophenol	40.000	41.628	-4.1	103	0.00
27		2,4-Dimethylphenol	40.000	40.391	-1.0	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.595	-1.5	100	0.00
29	C	2,4-Dichlorophenol	40.000	41.090	-2.7	99	0.00
30		1,2,4-Trichlorobenzene	40.000	41.459	-3.6	101	0.00
31		Naphthalene	40.000	41.558	-3.9	99	0.00
32		Benzoic acid	40.000	39.906	0.2	104	0.00
33		4-Chloroaniline	40.000	41.728	-4.3	100	0.00
34	C	Hexachlorobutadiene	40.000	41.035	-2.6	99	0.00
35		Caprolactam	40.000	43.943	-9.9	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.692	-6.7	101	0.00
37		2-Methylnaphthalene	40.000	42.598	-6.5	101	0.00
38		1-Methylnaphthalene	40.000	43.021	-7.6	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.571	-1.4	100	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.042	-0.1	101	0.00
42	S	2,4,6-Tribromophenol	80.000	79.326	0.8	102	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.037	-2.6	101	0.00
44		2,4,5-Trichlorophenol	40.000	40.704	-1.8	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114122.D
Acq On : 10 May 2019 15:56
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF051019

Quant Time: May 10 18:34:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	79.093	1.1	102	0.00
46		1,1'-Biphenyl	40.000	40.800	-2.0	102	0.00
47		2-Chloronaphthalene	40.000	40.749	-1.9	101	0.00
48		2-Nitroaniline	40.000	40.746	-1.9	102	0.00
49		Acenaphthylene	40.000	41.312	-3.3	102	0.00
50		Dimethylphthalate	40.000	40.295	-0.7	102	0.00
51		2,6-Dinitrotoluene	40.000	41.636	-4.1	104	0.00
52	C	Acenaphthene	40.000	41.210	-3.0	104	0.00
53		3-Nitroaniline	40.000	41.660	-4.1	104	0.00
54	P	2,4-Dinitrophenol	40.000	42.100	-5.3	117	0.00
55		Dibenzofuran	40.000	40.879	-2.2	104	0.00
56	P	4-Nitrophenol	40.000	41.185	-3.0	107	0.00
57		2,4-Dinitrotoluene	40.000	41.016	-2.5	105	0.00
58		Fluorene	40.000	39.781	0.5	102	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.065	-0.2	102	0.00
60		Diethylphthalate	40.000	40.127	-0.3	104	0.00
61		4-Chlorophenyl-phenylether	40.000	39.983	0.0	102	0.00
62		4-Nitroaniline	40.000	40.550	-1.4	107	0.00
63		Azobenzene	40.000	40.075	-0.2	103	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	44.034	-10.1	109	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.317	-3.3	102	0.00
67		4-Bromophenyl-phenylether	40.000	40.707	-1.8	102	0.00
68		Hexachlorobenzene	40.000	41.251	-3.1	103	0.00
69		Atrazine	40.000	41.034	-2.6	104	0.00
70	C	Pentachlorophenol	40.000	42.752	-6.9	104	0.00
71		Phenanthrene	40.000	41.734	-4.3	103	0.00
72		Anthracene	40.000	42.848	-7.1	105	0.00
73		Carbazole	40.000	41.579	-3.9	103	0.00
74		Di-n-butylphthalate	40.000	42.657	-6.6	105	0.00
75	C	Fluoranthene	40.000	40.954	-2.4	102	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	104	0.00
77		Benzidine	40.000	38.141	4.6	102	0.00
78		Pyrene	40.000	37.731	5.7	102	0.00
79	S	Terphenyl-d14	80.000	75.431	5.7	103	0.00
80		Butylbenzylphthalate	40.000	42.042	-5.1	110	0.00
81		Benzo(a)anthracene	40.000	40.124	-0.3	103	0.00
82		3,3'-Dichlorobenzidine	40.000	42.418	-6.0	104	0.00
83		Chrysene	40.000	40.946	-2.4	105	0.01
84		Bis(2-ethylhexyl)phthalate	40.000	40.246	-0.6	104	0.01
85	c	Di-n-octyl phthalate	40.000	42.900	-7.2	108	0.02
86		Indeno(1,2,3-cd)pyrene	40.000	45.541	-13.9	123	0.03
87	I	Perylene-d12	20.000	20.000	0.0	116	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114122.D
Acq On : 10 May 2019 15:56
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF051019

Quant Time: May 10 18:34:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	40.161	-0.4	116	0.02
89	Benzo(k)fluoranthene	40.000	40.465	-1.2	111	0.02
90 C	Benzo(a)pyrene	40.000	40.638	-1.6	115	0.02
91	Dibenzo(a,h)anthracene	40.000	42.402	-6.0	125	0.03
92	Benzo(g,h,i)perylene	40.000	39.884	0.3	120	0.03

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