Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112618\
Data File : BF110959.D
Acq On : 26 Nov 2018 13:05
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF112618

Quant Time: Nov 26 15:28:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 26 15:26:09 2018
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	41.885	-4.7	102	0.00
3		Pyridine	40.000	39.639	0.9	95	0.00
4		n-Nitrosodimethylamine	40.000	41.866	-4.7	104	0.00
5	S	2-Fluorophenol	80.000	84.750	-5.9	99	0.00
6		Aniline	40.000	41.833	-4.6	99	0.00
7	S	Phenol-d6	80.000	81.427	-1.8	100	0.00
8		2-Chlorophenol	40.000	41.305	-3.3	101	0.00
9		Benzaldehyde	40.000	41.845	-4.6	101	0.00
10	C	Phenol	40.000	41.890	-4.7	99	0.00
11		bis(2-Chloroethyl)ether	40.000	40.692	-1.7	101	0.00
12		1,3-Dichlorobenzene	40.000	41.310	-3.3	101	0.00
13	C	1,4-Dichlorobenzene	40.000	40.841	-2.1	100	0.00
14		1,2-Dichlorobenzene	40.000	40.670	-1.7	100	0.00
15		Benzyl Alcohol	40.000	42.332	-5.8	101	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.701	-1.8	100	0.00
17		2-Methylphenol	40.000	40.836	-2.1	100	0.00
18		Hexachloroethane	40.000	41.283	-3.2	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.647	-4.1	102	0.00
20		3+4-Methylphenols	40.000	40.406	-1.0	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	101	0.00
22		Acetophenone	40.000	40.399	-1.0	100	0.00
23	S	Nitrobenzene-d5	80.000	80.845	-1.1	101	0.00
24		Nitrobenzene	40.000	40.266	-0.7	102	0.00
25		Isophorone	40.000	39.690	0.8	99	0.00
26	C	2-Nitrophenol	40.000	41.827	-4.6	101	0.00
27		2,4-Dimethylphenol	40.000	41.056	-2.6	100	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.589	-1.5	101	0.00
29	C	2,4-Dichlorophenol	40.000	41.065	-2.7	99	0.00
30		1,2,4-Trichlorobenzene	40.000	40.637	-1.6	100	0.00
31		Naphthalene	40.000	40.269	-0.7	101	0.00
32		Benzoic acid	40.000	35.942	10.1	87	0.00
33		4-Chloroaniline	40.000	40.285	-0.7	99	0.00
34	C	Hexachlorobutadiene	40.000	40.523	-1.3	101	0.00
35		Caprolactam	40.000	39.753	0.6	98	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.333	-0.8	98	0.00
37		2-Methylnaphthalene	40.000	39.803	0.5	99	0.00
38		1-Methylnaphthalene	40.000	39.273	1.8	99	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.706	-1.8	100	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.613	-4.0	104	0.00
42	S	2,4,6-Tribromophenol	80.000	83.090	-3.9	101	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.003	-5.0	100	0.00
44		2,4,5-Trichlorophenol	40.000	41.769	-4.4	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112618\
Data File : BF110959.D
Acq On : 26 Nov 2018 13:05
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF112618

Quant Time: Nov 26 15:28:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 26 15:26:09 2018
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	80.888	-1.1	102	0.00
46		1,1'-Biphenyl	40.000	40.630	-1.6	100	0.00
47		2-Chloronaphthalene	40.000	40.382	-1.0	99	0.00
48		2-Nitroaniline	40.000	40.794	-2.0	100	0.00
49		Acenaphthylene	40.000	40.544	-1.4	100	0.00
50		Dimethylphthalate	40.000	40.151	-0.4	100	0.00
51		2,6-Dinitrotoluene	40.000	41.322	-3.3	100	0.00
52	C	Acenaphthene	40.000	40.345	-0.9	100	0.00
53		3-Nitroaniline	40.000	40.869	-2.2	100	0.00
54	P	2,4-Dinitrophenol	40.000	42.842	-7.1	100	0.00
55		Dibenzofuran	40.000	40.142	-0.4	101	0.00
56	P	4-Nitrophenol	40.000	39.684	0.8	96	0.00
57		2,4-Dinitrotoluene	40.000	40.876	-2.2	99	0.00
58		Fluorene	40.000	40.379	-0.9	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.899	0.3	98	0.00
60		Diethylphthalate	40.000	39.865	0.3	100	0.00
61		4-Chlorophenyl-phenylether	40.000	40.278	-0.7	100	0.00
62		4-Nitroaniline	40.000	41.155	-2.9	100	0.00
63		Azobenzene	40.000	40.311	-0.8	100	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.553	-6.4	102	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.015	-2.5	101	0.00
67		4-Bromophenyl-phenylether	40.000	40.792	-2.0	99	0.00
68		Hexachlorobenzene	40.000	40.706	-1.8	100	0.00
69		Atrazine	40.000	43.159	-7.9	99	0.00
70	C	Pentachlorophenol	40.000	41.691	-4.2	96	0.00
71		Phenanthrene	40.000	40.140	-0.4	100	0.00
72		Anthracene	40.000	40.140	-0.4	99	0.00
73		Carbazole	40.000	40.192	-0.5	99	0.00
74		Di-n-butylphthalate	40.000	40.511	-1.3	100	0.00
75	C	Fluoranthene	40.000	39.774	0.6	99	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	101	0.00
77		Benzidine	40.000	32.626	18.4	91	0.00
78		Pyrene	40.000	39.479	1.3	100	0.00
79	S	Terphenyl-d14	80.000	77.431	3.2	99	0.00
80		Butylbenzylphthalate	40.000	39.882	0.3	99	0.00
81		Benzo(a)anthracene	40.000	39.594	1.0	99	0.00
82		3,3'-Dichlorobenzidine	40.000	39.504	1.2	97	0.00
83		Chrysene	40.000	39.744	0.6	100	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.509	-1.3	101	0.00
85	c	Di-n-octyl phthalate	40.000	41.561	-3.9	103	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	38.113	4.7	101	0.01
87	I	Perylene-d12	20.000	20.000	0.0	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF112618\
Data File : BF110959.D
Acq On : 26 Nov 2018 13:05
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF112618

Quant Time: Nov 26 15:28:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 26 15:26:09 2018
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	40.326	-0.8	100	0.00
89	Benzo(k)fluoranthene	40.000	41.926	-4.8	104	0.00
90 C	Benzo(a)pyrene	40.000	39.820	0.4	101	0.00
91	Dibenzo(a,h)anthracene	40.000	39.980	0.1	101	0.00
92	Benzo(g,h,i)perylene	40.000	39.887	0.3	100	0.00

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