Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF122118\
Data File : BF111681.D
Acq On : 21 Dec 2018 12:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 22:46:41 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 19 17:39:39 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	80	0.00
2		1,4-Dioxane	40.000	39.746	0.6	82	-0.03
3		Pyridine	40.000	40.140	-0.4	83	-0.02
4		n-Nitrosodimethylamine	40.000	40.962	-2.4	83	-0.02
5	S	2-Fluorophenol	80.000	83.494	-4.4	85	0.00
6		Aniline	40.000	42.055	-5.1	85	0.00
7	S	Phenol-d6	80.000	84.479	-5.6	86	0.00
8		2-Chlorophenol	40.000	42.392	-6.0	86	0.00
9		Benzaldehyde	40.000	40.873	-2.2	84	0.00
10	C	Phenol	40.000	42.240	-5.6	86	0.00
11		bis(2-Chloroethyl)ether	40.000	40.552	-1.4	82	0.00
12		1,3-Dichlorobenzene	40.000	42.258	-5.6	85	0.00
13	C	1,4-Dichlorobenzene	40.000	42.002	-5.0	85	0.00
14		1,2-Dichlorobenzene	40.000	42.450	-6.1	86	0.00
15		Benzyl Alcohol	40.000	43.103	-7.8	86	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.460	3.8	78	0.00
17		2-Methylphenol	40.000	42.026	-5.1	85	0.00
18		Hexachloroethane	40.000	42.570	-6.4	85	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.102	-2.8	85	0.00
20		3+4-Methylphenols	40.000	42.377	-5.9	85	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	83	0.00
22		Acetophenone	40.000	41.360	-3.4	86	0.00
23	S	Nitrobenzene-d5	80.000	92.469	-15.6	91	0.00
24		Nitrobenzene	40.000	44.842	-12.1	88	0.00
25		Isophorone	40.000	40.936	-2.3	84	0.00
26	C	2-Nitrophenol	40.000	53.828	-34.6#	102	0.00
27		2,4-Dimethylphenol	40.000	41.472	-3.7	84	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.799	-2.0	84	0.00
29	C	2,4-Dichlorophenol	40.000	42.332	-5.8	86	0.00
30		1,2,4-Trichlorobenzene	40.000	41.671	-4.2	86	0.00
31		Naphthalene	40.000	41.908	-4.8	86	0.00
32		Benzoic acid	40.000	48.972	-22.4	98	0.00
33		4-Chloroaniline	40.000	43.138	-7.8	88	0.00
34	C	Hexachlorobutadiene	40.000	41.829	-4.6	86	0.00
35		Caprolactam	40.000	44.643	-11.6	92	0.00
36	C	4-Chloro-3-methylphenol	40.000	43.697	-9.2	89	0.00
37		2-Methylnaphthalene	40.000	42.453	-6.1	87	0.00
38		1-Methylnaphthalene	40.000	42.525	-6.3	87	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	84	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.834	-4.6	88	0.00
41	P	Hexachlorocyclopentadiene	40.000	46.933	-17.3	97	0.00
42	S	2,4,6-Tribromophenol	80.000	92.175	-15.2	95	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.771	-6.9	91	0.00
44		2,4,5-Trichlorophenol	40.000	43.594	-9.0	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF122118\
Data File : BF111681.D
Acq On : 21 Dec 2018 12:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 22:46:41 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 19 17:39:39 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	81.724	-2.2	88	0.00
46		1,1'-Biphenyl	40.000	42.072	-5.2	89	0.00
47		2-Chloronaphthalene	40.000	41.958	-4.9	89	0.00
48		2-Nitroaniline	40.000	50.416	-26.0#	103	0.00
49		Acenaphthylene	40.000	42.707	-6.8	91	0.00
50		Dimethylphthalate	40.000	42.207	-5.5	89	0.00
51		2,6-Dinitrotoluene	40.000	49.095	-22.7	100	0.00
52	C	Acenaphthene	40.000	43.581	-9.0	93	0.00
53		3-Nitroaniline	40.000	51.861	-29.7#	99	0.00
54	P	2,4-Dinitrophenol	40.000	61.795	-54.5#	147	0.00
55		Dibenzofuran	40.000	42.549	-6.4	89	0.00
56	P	4-Nitrophenol	40.000	52.815	-32.0#	101	0.01
57		2,4-Dinitrotoluene	40.000	54.004	-35.0#	108	0.00
58		Fluorene	40.000	42.071	-5.2	90	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	44.495	-11.2	92	0.00
60		Diethylphthalate	40.000	42.002	-5.0	88	0.00
61		4-Chlorophenyl-phenylether	40.000	41.763	-4.4	90	0.00
62		4-Nitroaniline	40.000	52.232	-30.6#	107	0.00
63		Azobenzene	40.000	41.532	-3.8	87	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	40.000	55.210	-38.0#	130	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.565	-3.9	89	0.00
67		4-Bromophenyl-phenylether	40.000	42.115	-5.3	92	0.00
68		Hexachlorobenzene	40.000	42.225	-5.6	91	0.00
69		Atrazine	40.000	41.329	-3.3	89	0.00
70	C	Pentachlorophenol	40.000	44.861	-12.2	91	0.00
71		Phenanthrene	40.000	42.313	-5.8	92	0.00
72		Anthracene	40.000	42.252	-5.6	92	0.00
73		Carbazole	40.000	43.191	-8.0	93	0.00
74		Di-n-butylphthalate	40.000	40.870	-2.2	88	0.00
75	C	Fluoranthene	40.000	43.597	-9.0	94	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	93	0.00
77		Benzidine	40.000	37.488	6.3	86	0.00
78		Pyrene	40.000	41.312	-3.3	95	0.00
79	S	Terphenyl-d14	80.000	79.572	0.5	93	0.00
80		Butylbenzylphthalate	40.000	42.703	-6.8	95	0.00
81		Benzo(a)anthracene	40.000	42.445	-6.1	100	0.00
82		3,3'-Dichlorobenzidine	40.000	44.015	-10.0	99	0.00
83		Chrysene	40.000	42.013	-5.0	97	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.770	-4.4	95	0.00
85	c	Di-n-octyl phthalate	40.000	40.124	-0.3	91	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	35.355	11.6	82	0.01
87	I	Perylene-d12	20.000	20.000	0.0	78	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF122118\
Data File : BF111681.D
Acq On : 21 Dec 2018 12:15
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 22:46:41 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 19 17:39:39 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	45.657	-14.1	86	0.00
89	Benzo(k)fluoranthene	40.000	44.564	-11.4	89	0.00
90 C	Benzo(a)pyrene	40.000	43.017	-7.5	84	0.00
91	Dibenzo(a,h)anthracene	40.000	42.882	-7.2	81	0.01
92	Benzo(g,h,i)perylene	40.000	44.900	-12.2	84	0.02

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