Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF101118\
Data File : BF109811.D
Acq On : 12 Oct 2018 3:10
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 12 08:58:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 08 16:02:11 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	67	0.00
2		1,4-Dioxane	40.000	45.192	-13.0	75	-0.06
3		Pyridine	40.000	36.368	9.1	58	-0.05
4		n-Nitrosodimethylamine	40.000	37.122	7.2	60	-0.05
5	S	2-Fluorophenol	80.000	87.911	-9.9	67	-0.01
6		Aniline	40.000	39.565	1.1	63	0.00
7	S	Phenol-d6	80.000	79.429	0.7	64	0.00
8		2-Chlorophenol	40.000	41.282	-3.2	66	0.00
9		Benzaldehyde	40.000	40.795	-2.0	62	0.00
10	C	Phenol	40.000	36.917	7.7	63	0.00
11		bis(2-Chloroethyl)ether	40.000	33.982	15.0	62	0.00
12		1,3-Dichlorobenzene	40.000	40.659	-1.6	66	0.00
13	C	1,4-Dichlorobenzene	40.000	41.107	-2.8	69	0.00
14		1,2-Dichlorobenzene	40.000	41.712	-4.3	69	0.00
15		Benzyl Alcohol	40.000	37.130	7.2	58	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.740	5.6	63	0.00
17		2-Methylphenol	40.000	36.705	8.2	61	0.00
18		Hexachloroethane	40.000	35.167	12.1	58	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.910	7.7	62	0.00
20		3+4-Methylphenols	40.000	35.349	11.6	56	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	59	0.00
22		Acetophenone	40.000	41.660	-4.1	61	0.00
23	S	Nitrobenzene-d5	80.000	85.917	-7.4	63	0.00
24		Nitrobenzene	40.000	42.022	-5.1	62	0.00
25		Isophorone	40.000	38.791	3.0	58	0.00
26	C	2-Nitrophenol	40.000	21.519	46.2#	30	0.00
27		2,4-Dimethylphenol	40.000	37.903	5.2	55	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.426	1.4	58	0.00
29	C	2,4-Dichlorophenol	40.000	37.289	6.8	54	0.00
30		1,2,4-Trichlorobenzene	40.000	40.173	-0.4	59	0.00
31		Naphthalene	40.000	37.805	5.5	56	0.00
32		Benzoic acid	40.000	20.041	49.9#	32	-0.04
33		4-Chloroaniline	40.000	38.240	4.4	57	0.00
34	C	Hexachlorobutadiene	40.000	41.906	-4.8	62	0.00
35		Caprolactam	40.000	35.190	12.0	51	-0.02
36	C	4-Chloro-3-methylphenol	40.000	35.140	12.1	51	-0.01
37		2-Methylnaphthalene	40.000	35.489	11.3	53	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	51	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	42.656	-6.6	54	0.00
40	P	Hexachlorocyclopentadiene	40.000	8.140	79.7#	1	0.00
41	S	2,4,6-Tribromophenol	80.000	88.788	-11.0	57	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.815	3.0	48	0.00
43		2,4,5-Trichlorophenol	40.000	39.954	0.1	50	0.00
44	S	2-Fluorobiphenyl	80.000	85.701	-7.1	56	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF101118\
Data File : BF109811.D
Acq On : 12 Oct 2018 3:10
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 12 08:58:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 08 16:02:11 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		1,1'-Biphenyl	40.000	41.848	-4.6	54	0.00
46		2-Chloronaphthalene	40.000	40.763	-1.9	52	0.00
47		2-Nitroaniline	40.000	42.898	-7.2	54	0.00
48		Acenaphthylene	40.000	40.434	-1.1	52	0.00
49		Dimethylphthalate	40.000	41.240	-3.1	54	-0.01
50		2,6-Dinitrotoluene	40.000	36.905	7.7	46	0.00
51	C	Acenaphthene	40.000	40.220	-0.5	53	0.00
52		3-Nitroaniline	40.000	46.613	-16.5	59	0.00
53	P	2,4-Dinitrophenol	40.000	0.000	100.0#	0	-9.80#
54		Dibenzofuran	40.000	40.303	-0.8	53	0.00
55	P	4-Nitrophenol	40.000	31.156	22.1	38	0.00
56		2,4-Dinitrotoluene	40.000	35.091	12.3	44	0.00
57		Fluorene	40.000	40.579	-1.4	53	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	43.368	-8.4	54	0.00
59		Diethylphthalate	40.000	41.587	-4.0	55	-0.01
60		4-Chlorophenyl-phenylether	40.000	40.767	-1.9	54	0.00
61		4-Nitroaniline	40.000	49.766	-24.4	61	-0.01
62		Azobenzene	40.000	40.602	-1.5	52	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	59	0.00
64		4,6-Dinitro-2-methylphenol	40.000	0.466	98.8#	1	0.00
65	c	n-Nitrosodiphenylamine	40.000	36.090	9.8	54	-0.01
66		4-Bromophenyl-phenylether	40.000	34.877	12.8	52	0.00
67		Hexachlorobenzene	40.000	35.957	10.1	53	0.00
68		Atrazine	40.000	34.324	14.2	51	0.00
69	C	Pentachlorophenol	40.000	41.111	-2.8	58	0.00
70		Phenanthrene	40.000	38.369	4.1	58	0.00
71		Anthracene	40.000	41.610	-4.0	62	0.00
72		Carbazole	40.000	42.841	-7.1	63	0.00
73		Di-n-butylphthalate	40.000	39.800	0.5	59	0.00
74	C	Fluoranthene	40.000	47.142	-17.9	69	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	86	0.00
76		Benzidine	40.000	37.139	7.2	81	0.00
77		Pyrene	40.000	33.143	17.1	71	0.00
78	S	Terphenyl-d14	80.000	67.198	16.0	74	0.00
79		Butylbenzylphthalate	40.000	36.530	8.7	77	0.00
80		Benzo(a)anthracene	40.000	37.781	5.5	82	0.00
81		3,3'-Dichlorobenzidine	40.000	41.914	-4.8	88	0.00
82		Chrysene	40.000	39.565	1.1	84	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.280	-0.7	86	0.00
84	c	Di-n-octyl phthalate	40.000	46.691	-16.7	97	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	31.858	20.4	71	0.00
86	I	Perylene-d12	20.000	20.000	0.0	75	0.00
87		Benzo(b)fluoranthene	40.000	41.935	-4.8	81	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF101118\
Data File : BF109811.D
Acq On : 12 Oct 2018 3:10
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 12 08:58:44 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 08 16:02:11 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(k)fluoranthene	40.000	40.522	-1.3	72	0.00
89 C	Benzo(a)pyrene	40.000	38.658	3.4	72	0.00
90	Dibenzo(a,h)anthracene	40.000	37.691	5.8	72	0.00
91	Benzo(g,h,i)perylene	40.000	37.239	6.9	69	0.00

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