Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061918\
Data File : BF106581.D
Acq On : 19 Jun 2018 14:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 16:17:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 19 14:32:09 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	105	0.00
2		1,4-Dioxane	40.000	38.283	4.3	102	0.00
3		Pyridine	40.000	38.024	4.9	101	0.00
4		n-Nitrosodimethylamine	40.000	38.709	3.2	101	0.00
5	S	2-Fluorophenol	80.000	76.363	4.5	103	0.00
6		Aniline	40.000	38.367	4.1	103	0.00
7	S	Phenol-d6	80.000	75.254	5.9	102	0.00
8		2-Chlorophenol	40.000	38.063	4.8	102	0.00
9		Benzaldehyde	40.000	37.000	7.5	102	0.00
10	C	Phenol	40.000	37.953	5.1	103	0.00
11		bis(2-Chloroethyl)ether	40.000	37.551	6.1	101	0.00
12		1,3-Dichlorobenzene	40.000	37.960	5.1	102	0.00
13	C	1,4-Dichlorobenzene	40.000	38.281	4.3	104	0.00
14		1,2-Dichlorobenzene	40.000	38.107	4.7	102	0.00
15		Benzyl Alcohol	40.000	38.350	4.1	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.817	5.5	102	0.00
17		2-Methylphenol	40.000	38.148	4.6	103	0.00
18		Hexachloroethane	40.000	37.968	5.1	103	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.080	9.8	102	0.00
20		3+4-Methylphenols	40.000	37.930	5.2	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	38.221	4.4	101	0.00
23	S	Nitrobenzene-d5	80.000	75.946	5.1	101	0.00
24		Nitrobenzene	40.000	38.699	3.3	102	0.00
25		Isophorone	40.000	38.328	4.2	103	0.00
26	C	2-Nitrophenol	40.000	39.706	0.7	101	0.00
27		2,4-Dimethylphenol	40.000	39.269	1.8	102	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.277	4.3	103	0.00
29	C	2,4-Dichlorophenol	40.000	39.024	2.4	102	0.00
30		1,2,4-Trichlorobenzene	40.000	38.488	3.8	102	0.00
31		Naphthalene	40.000	38.074	4.8	103	0.00
32		Benzoic acid	40.000	37.252	6.9	100	0.00
33		4-Chloroaniline	40.000	38.162	4.6	102	0.00
34	C	Hexachlorobutadiene	40.000	39.037	2.4	103	0.00
35		Caprolactam	40.000	39.186	2.0	101	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.398	4.0	102	0.00
37		2-Methylnaphthalene	40.000	37.802	5.5	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	38.295	4.3	102	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.886	0.3	102	0.00
41	S	2,4,6-Tribromophenol	80.000	78.548	1.8	102	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.671	3.3	103	0.00
43		2,4,5-Trichlorophenol	40.000	39.245	1.9	103	0.00
44	S	2-Fluorobiphenyl	80.000	78.701	1.6	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061918\
Data File : BF106581.D
Acq On : 19 Jun 2018 14:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 16:17:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 19 14:32:09 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		1,1'-Biphenyl	40.000	37.455	6.4	101	0.00
46		2-Chloronaphthalene	40.000	37.668	5.8	101	0.00
47		2-Nitroaniline	40.000	39.242	1.9	102	0.00
48		Acenaphthylene	40.000	37.868	5.3	102	0.00
49		Dimethylphthalate	40.000	37.547	6.1	102	0.00
50		2,6-Dinitrotoluene	40.000	38.407	4.0	102	0.00
51	C	Acenaphthene	40.000	36.213	9.5	96	0.00
52		3-Nitroaniline	40.000	39.681	0.8	104	0.00
53	P	2,4-Dinitrophenol	40.000	38.990	2.5	105	0.00
54		Dibenzofuran	40.000	38.319	4.2	103	0.00
55	P	4-Nitrophenol	40.000	40.919	-2.3	103	0.00
56		2,4-Dinitrotoluene	40.000	39.981	0.0	103	0.00
57		Fluorene	40.000	37.062	7.3	102	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.779	0.6	104	0.00
59		Diethylphthalate	40.000	37.970	5.1	103	0.00
60		4-Chlorophenyl-phenylether	40.000	37.645	5.9	101	0.00
61		4-Nitroaniline	40.000	39.143	2.1	103	0.00
62		Azobenzene	40.000	37.760	5.6	101	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.405	-1.0	101	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.570	3.6	102	0.00
66		4-Bromophenyl-phenylether	40.000	38.825	2.9	102	0.00
67		Hexachlorobenzene	40.000	39.209	2.0	102	0.00
68		Atrazine	40.000	39.432	1.4	103	0.00
69	C	Pentachlorophenol	40.000	41.129	-2.8	103	0.00
70		Phenanthrene	40.000	39.347	1.6	102	0.00
71		Anthracene	40.000	39.111	2.2	102	0.00
72		Carbazole	40.000	38.019	5.0	101	0.00
73		Di-n-butylphthalate	40.000	38.137	4.7	100	0.00
74	C	Fluoranthene	40.000	39.081	2.3	100	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	100	0.00
76		Benzidine	40.000	41.741	-4.4	103	0.00
77		Pyrene	40.000	38.194	4.5	102	0.00
78	S	Terphenyl-d14	80.000	76.743	4.1	98	0.00
79		Butylbenzylphthalate	40.000	38.960	2.6	101	0.00
80		Benzo(a)anthracene	40.000	38.148	4.6	100	0.00
81		3,3'-Dichlorobenzidine	40.000	38.072	4.8	99	0.00
82		Chrysene	40.000	37.945	5.1	101	0.01
83		Bis(2-ethylhexyl)phthalate	40.000	38.900	2.8	102	0.00
84	c	Di-n-octyl phthalate	40.000	38.233	4.4	100	0.02
85		Indeno(1,2,3-cd)pyrene	40.000	36.544	8.6	102	0.03
86	I	Perylene-d12	20.000	20.000	0.0	99	0.02
87		Benzo(b)fluoranthene	40.000	37.917	5.2	97	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061918\
Data File : BF106581.D
Acq On : 19 Jun 2018 14:48
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 16:17:19 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 19 14:32:09 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(k)fluoranthene	40.000	40.661	-1.7	102	0.02
89 C	Benzo(a)pyrene	40.000	38.426	3.9	96	0.02
90	Dibenzo(a,h)anthracene	40.000	38.370	4.1	100	0.03
91	Benzo(g,h,i)perylene	40.000	38.634	3.4	101	0.02

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