Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM113018\
Data File : BM017816.D
Acq On : 30 Nov 2018 10:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 30 14:09:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 30 13:21:52 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	96	-0.03
2		1,4-Dioxane	40.000	34.418	14.0	85	-0.02
3		Pyridine	40.000	34.943	12.6	83	-0.02
4		n-Nitrosodimethylamine	40.000	31.005	22.5	73	-0.02
5	S	2-Fluorophenol	80.000	73.647	7.9	88	-0.03
6		Aniline	40.000	36.948	7.6	86	-0.03
7	S	Phenol-d6	80.000	73.532	8.1	86	-0.03
8		2-Chlorophenol	40.000	38.112	4.7	91	-0.03
9		Benzaldehyde	40.000	32.219	19.5	76	-0.03
10	C	Phenol	40.000	36.563	8.6	86	-0.03
11		bis(2-Chloroethyl)ether	40.000	37.918	5.2	89	-0.03
12		1,3-Dichlorobenzene	40.000	38.633	3.4	94	-0.03
13	C	1,4-Dichlorobenzene	40.000	38.778	3.1	93	-0.03
14		1,2-Dichlorobenzene	40.000	38.302	4.2	92	-0.03
15		Benzyl Alcohol	40.000	37.427	6.4	86	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	30.290	24.3	72	-0.02
17		2-Methylphenol	40.000	37.740	5.6	88	-0.03
18		Hexachloroethane	40.000	40.606	-1.5	95	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	38.131	4.7	88	-0.03
20		3+4-Methylphenols	40.000	37.546	6.1	87	-0.03
21	I	Naphthalene-d8	20.000	20.000	0.0	90	-0.03
22		Acetophenone	40.000	39.564	1.1	87	-0.03
23	S	Nitrobenzene-d5	80.000	78.645	1.7	86	-0.03
24		Nitrobenzene	40.000	38.462	3.8	86	-0.03
25		Isophorone	40.000	40.505	-1.3	88	-0.04
26	C	2-Nitrophenol	40.000	40.524	-1.3	87	-0.03
27		2,4-Dimethylphenol	40.000	39.536	1.2	86	-0.03
28		bis(2-Chloroethoxy)methane	40.000	39.615	1.0	86	-0.03
29	C	2,4-Dichlorophenol	40.000	40.318	-0.8	86	-0.04
30		1,2,4-Trichlorobenzene	40.000	40.044	-0.1	89	-0.03
31		Naphthalene	40.000	39.446	1.4	87	-0.04
32		Benzoic acid	40.000	39.291	1.8	86	-0.04
33		4-Chloroaniline	40.000	38.380	4.0	82	-0.04
34	C	Hexachlorobutadiene	40.000	41.067	-2.7	92	-0.04
35		Caprolactam	40.000	36.826	7.9	82	-0.04
36	C	4-Chloro-3-methylphenol	40.000	38.031	4.9	82	-0.03
37		2-Methylnaphthalene	40.000	38.672	3.3	85	-0.03
38		1-Methylnaphthalene	40.000	38.297	4.3	84	-0.03
39	I	Acenaphthene-d10	20.000	20.000	0.0	82	-0.03
40		1,2,4,5-Tetrachlorobenzene	40.000	41.523	-3.8	85	-0.03
41	P	Hexachlorocyclopentadiene	40.000	42.687	-6.7	86	-0.03
42	S	2,4,6-Tribromophenol	80.000	75.710	5.4	75	-0.02
43	C	2,4,6-Trichlorophenol	40.000	42.248	-5.6	84	-0.03
44		2,4,5-Trichlorophenol	40.000	40.013	-0.0	80	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM113018\
Data File : BM017816.D
Acq On : 30 Nov 2018 10:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 30 14:09:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 30 13:21:52 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	81.461	-1.8	83	-0.03
46		1,1'-Biphenyl	40.000	40.476	-1.2	83	-0.04
47		2-Chloronaphthalene	40.000	40.084	-0.2	82	-0.03
48		2-Nitroaniline	40.000	38.141	4.6	74	-0.03
49		Acenaphthylene	40.000	39.709	0.7	81	-0.03
50		Dimethylphthalate	40.000	38.825	2.9	80	-0.02
51		2,6-Dinitrotoluene	40.000	39.003	2.5	78	-0.03
52	C	Acenaphthene	40.000	39.073	2.3	80	-0.03
53		3-Nitroaniline	40.000	37.596	6.0	77	-0.02
54	P	2,4-Dinitrophenol	40.000	37.520	6.2	78	-0.02
55		Dibenzofuran	40.000	38.405	4.0	79	-0.02
56	P	4-Nitrophenol	40.000	36.427	8.9	76	-0.03
57		2,4-Dinitrotoluene	40.000	39.074	2.3	76	-0.02
58		Fluorene	40.000	38.026	4.9	78	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	38.573	3.6	77	-0.03
60		Diethylphthalate	40.000	38.283	4.3	81	-0.02
61		4-Chlorophenyl-phenylether	40.000	38.079	4.8	78	-0.02
62		4-Nitroaniline	40.000	36.571	8.6	74	-0.02
63		Azobenzene	40.000	39.227	1.9	79	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	73	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	42.638	-6.6	75	-0.02
66	c	n-Nitrosodiphenylamine	40.000	42.975	-7.4	77	-0.03
67		4-Bromophenyl-phenylether	40.000	42.584	-6.5	76	-0.03
68		Hexachlorobenzene	40.000	40.692	-1.7	74	-0.02
69		Atrazine	40.000	41.083	-2.7	72	-0.02
70	C	Pentachlorophenol	40.000	40.716	-1.8	73	-0.02
71		Phenanthrene	40.000	39.478	1.3	72	-0.02
72		Anthracene	40.000	40.248	-0.6	72	-0.03
73		Carbazole	40.000	39.450	1.4	70	-0.02
74		Di-n-butylphthalate	40.000	43.675	-9.2	76	-0.02
75	C	Fluoranthene	40.000	35.000	12.5	62	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	55	-0.02
77		Benzidine	40.000	43.499	-8.7	57	-0.02
78		Pyrene	40.000	43.863	-9.7	60	-0.02
79	S	Terphenyl-d14	80.000	88.481	-10.6	61	-0.02
80		Butylbenzylphthalate	40.000	47.556	-18.9	61	-0.02
81		Benzo(a)anthracene	40.000	39.138	2.2	53	-0.02
82		3,3'-Dichlorobenzidine	40.000	42.131	-5.3	54	-0.02
83		Chrysene	40.000	39.176	2.1	53	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	44.952	-12.4	60	-0.02
85	c	Di-n-octyl phthalate	40.000	43.007	-7.5	57	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	55.290	-38.2#	73	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	61	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM113018\
Data File : BM017816.D
Acq On : 30 Nov 2018 10:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Nov 30 14:09:54 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 30 13:21:52 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	37.862	5.3	58	-0.03
89	Benzo(k)fluoranthene	40.000	35.913	10.2	53	-0.03
90 C	Benzo(a)pyrene	40.000	39.194	2.0	59	-0.03
91	Dibenzo(a,h)anthracene	40.000	48.443	-21.1	74	-0.05
92	Benzo(g,h,i)perylene	40.000	50.178	-25.4#	76	-0.05

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