Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051617\
Data File : BM010055.D
Acq On : 16 May 2017 16:08
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 17 05:39:55 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 11 19:50:16 2017
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	72	-0.02
2		1,4-Dioxane	40.000	46.709	-16.8	81	0.00
3		Pyridine	40.000	45.443	-13.6	78	-0.01
4		n-Nitrosodimethylamine	40.000	65.646	-64.1#	111	-0.01
5	S	2-Fluorophenol	80.000	88.218	-10.3	76	-0.02
6		Aniline	40.000	43.435	-8.6	75	-0.02
7	S	Phenol-d6	80.000	86.755	-8.4	75	-0.04
8		2-Chlorophenol	40.000	42.428	-6.1	72	-0.03
9		Benzaldehyde	40.000	45.216	-13.0	77	-0.02
10	C	Phenol	40.000	42.609	-6.5	74	-0.04
11		bis(2-Chloroethyl)ether	40.000	41.317	-3.3	71	-0.03
12		1,3-Dichlorobenzene	40.000	42.242	-5.6	74	-0.03
13	C	1,4-Dichlorobenzene	40.000	41.224	-3.1	72	-0.04
14		1,2-Dichlorobenzene	40.000	43.215	-8.0	75	-0.03
15		Benzyl Alcohol	40.000	45.325	-13.3	76	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	46.570	-16.4	79	-0.03
17		2-Methylphenol	40.000	43.757	-9.4	74	-0.03
18		Hexachloroethane	40.000	52.524	-31.3#	86	-0.04
19	P	n-Nitroso-di-n-propylamine	40.000	48.260	-20.6	81	-0.03
20		3+4-Methylphenols	40.000	44.493	-11.2	76	-0.03
21	I	Naphthalene-d8	20.000	20.000	0.0	75	-0.03
22		Acetophenone	40.000	41.270	-3.2	78	-0.03
23	S	Nitrobenzene-d5	80.000	90.930	-13.7	84	-0.03
24		Nitrobenzene	40.000	45.239	-13.1	85	-0.03
25		Isophorone	40.000	43.258	-8.1	79	-0.04
26	C	2-Nitrophenol	40.000	40.119	-0.3	73	-0.03
27		2,4-Dimethylphenol	40.000	40.807	-2.0	74	-0.03
28		bis(2-Chloroethoxy)methane	40.000	41.054	-2.6	76	-0.02
29	C	2,4-Dichlorophenol	40.000	40.771	-1.9	77	-0.03
30		1,2,4-Trichlorobenzene	40.000	38.842	2.9	72	-0.04
31		Naphthalene	40.000	40.252	-0.6	76	-0.04
32		Benzoic acid	40.000	41.031	-2.6	77	-0.05
33		4-Chloroaniline	40.000	40.512	-1.3	74	-0.03
34	C	Hexachlorobutadiene	40.000	40.733	-1.8	74	-0.04
35		Caprolactam	40.000	39.065	2.3	74	-0.03
36	C	4-Chloro-3-methylphenol	40.000	42.637	-6.6	78	-0.04
37		2-Methylnaphthalene	40.000	38.733	3.2	73	-0.03
38	I	Acenaphthene-d10	20.000	20.000	0.0	76	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	37.406	6.5	70	-0.04
40	P	Hexachlorocyclopentadiene	40.000	42.804	-7.0	90	-0.03
41	S	2,4,6-Tribromophenol	80.000	82.352	-2.9	72	-0.02
42	C	2,4,6-Trichlorophenol	40.000	39.622	0.9	74	-0.03
43		2,4,5-Trichlorophenol	40.000	38.955	2.6	71	-0.04
44	S	2-Fluorobiphenyl	80.000	79.674	0.4	74	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051617\
Data File : BM010055.D
Acq On : 16 May 2017 16:08
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 17 05:39:55 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 11 19:50:16 2017
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	40.615	-1.5	76	-0.04
46		2-Chloronaphthalene	40.000	40.813	-2.0	75	-0.04
47		2-Nitroaniline	40.000	52.430	-31.1#	91	-0.02
48		Acenaphthylene	40.000	41.130	-2.8	74	-0.03
49		Dimethylphthalate	40.000	42.658	-6.6	79	-0.02
50		2,6-Dinitrotoluene	40.000	41.892	-4.7	75	-0.02
51	C	Acenaphthene	40.000	40.873	-2.2	76	-0.03
52		3-Nitroaniline	40.000	42.902	-7.3	76	-0.02
53	P	2,4-Dinitrophenol	40.000	35.234	11.9	66	-0.02
54		Dibenzofuran	40.000	41.305	-3.3	77	-0.03
55	P	4-Nitrophenol	40.000	40.006	-0.0	73	-0.03
56		2,4-Dinitrotoluene	40.000	41.497	-3.7	78	-0.03
57		Fluorene	40.000	42.182	-5.5	77	-0.03
58		2,3,4,6-Tetrachlorophenol	40.000	38.320	4.2	70	-0.03
59		Diethylphthalate	40.000	45.269	-13.2	84	-0.03
60		4-Chlorophenyl-phenylether	40.000	40.107	-0.3	74	-0.03
61		4-Nitroaniline	40.000	43.635	-9.1	77	-0.02
62		Azobenzene	40.000	52.778	-31.9#	94	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	76	-0.03
64		4,6-Dinitro-2-methylphenol	40.000	36.395	9.0	70	-0.03
65	c	n-Nitrosodiphenylamine	40.000	40.250	-0.6	76	-0.03
66		4-Bromophenyl-phenylether	40.000	39.887	0.3	74	-0.03
67		Hexachlorobenzene	40.000	39.313	1.7	74	-0.04
68		Atrazine	40.000	43.866	-9.7	78	-0.02
69	C	Pentachlorophenol	40.000	39.548	1.1	80	-0.03
70		Phenanthrene	40.000	41.527	-3.8	79	-0.03
71		Anthracene	40.000	41.571	-3.9	78	-0.03
72		Carbazole	40.000	42.456	-6.1	79	-0.03
73		Di-n-butylphthalate	40.000	47.358	-18.4	87	-0.03
74	C	Fluoranthene	40.000	41.894	-4.7	78	-0.03
75	I	Chrysene-d12	20.000	20.000	0.0	78	-0.03
76		Benzidine	40.000	34.371	14.1	60	-0.02
77		Pyrene	40.000	41.057	-2.6	79	-0.03
78	S	Terphenyl-d14	80.000	85.939	-7.4	84	-0.02
79		Butylbenzylphthalate	40.000	48.562	-21.4	93	-0.02
80		Benzo(a)anthracene	40.000	40.515	-1.3	78	-0.03
81		3,3'-Dichlorobenzidine	40.000	38.779	3.1	72	-0.03
82		Chrysene	40.000	40.661	-1.7	78	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	49.250	-23.1	94	-0.02
84	c	Di-n-octyl phthalate	40.000	50.273	-25.7#	96	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	38.406	4.0	70	-0.08
86	I	Perylene-d12	20.000	20.000	0.0	71	-0.05
87		Benzo(b)fluoranthene	40.000	40.700	-1.8	74	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051617\
Data File : BM010055.D
Acq On : 16 May 2017 16:08
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 17 05:39:55 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 11 19:50:16 2017
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	42.969	-7.4	75	-0.04
89 C	Benzo(a)pyrene	40.000	41.292	-3.2	73	-0.04
90	Dibenzo(a,h)anthracene	40.000	39.140	2.1	69	-0.08
91	Benzo(g,h,i)perylene	40.000	40.067	-0.2	70	-0.08

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