Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM052617\
Data File : BM010249.D
Acq On : 26 May 2017 10:23
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 26 23:52:57 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 20 04:25:49 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	48#	-0.04
2		1,4-Dioxane	0.509	0.538	-5.7	52	-0.02
3		Pyridine	1.362	1.430	-5.0	49#	-0.04
4		n-Nitrosodimethylamine	0.493	0.677	-37.3#	70	-0.02
5	S	2-Fluorophenol	1.108	1.080	2.5	47#	-0.04
6		Aniline	1.766	1.671	5.4	45#	-0.04
7	S	Phenol-d6	1.425	1.416	0.6	47#	-0.04
8		2-Chlorophenol	1.205	1.161	3.7	47#	-0.04
9		Benzaldehyde	0.887	0.883	0.5	46#	-0.04
10	C	Phenol	1.502	1.487	1.0	48#	-0.04
11		bis(2-Chloroethyl)ether	1.119	1.073	4.1	45#	-0.04
12		1,3-Dichlorobenzene	1.531	1.502	1.9	47#	-0.04
13	C	1,4-Dichlorobenzene	1.573	1.541	2.0	47#	-0.04
14		1,2-Dichlorobenzene	1.510	1.472	2.5	47#	-0.04
15		Benzyl Alcohol	1.078	0.958	11.1	41#	-0.04
16		2,2'-oxybis(1-Chloropropane	1.101	0.946	14.1	40#	-0.04
17		2-Methylphenol	1.069	1.035	3.2	45#	-0.04
18		Hexachloroethane	0.510	0.532	-4.3	50	-0.04
19	P	n-Nitroso-di-n-propylamine	1.006	1.116	-10.9	52	-0.05
20		3+4-Methylphenols	1.443	1.450	-0.5	47#	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	45#	-0.04
22		Acetophenone	0.513	0.543	-5.8	48#	-0.04
23	S	Nitrobenzene-d5	0.371	0.456	-22.9	54	-0.04
24		Nitrobenzene	0.382	0.471	-23.3	55	-0.04
25		Isophorone	0.641	0.730	-13.9	50	-0.05
26	C	2-Nitrophenol	0.161	0.179	-11.2	49#	-0.04
27		2,4-Dimethylphenol	0.291	0.297	-2.1	46#	-0.04
28		bis(2-Chloroethoxy)methane	0.407	0.412	-1.2	45#	-0.04
29	C	2,4-Dichlorophenol	0.335	0.374	-11.6	50#	-0.04
30		1,2,4-Trichlorobenzene	0.433	0.478	-10.4	51	-0.04
31		Naphthalene	1.070	1.072	-0.2	45#	-0.04
32		Benzoic acid	0.198	0.195	1.5	44#	-0.05
33		4-Chloroaniline	0.412	0.376	8.7	40#	-0.04
34	C	Hexachlorobutadiene	0.289	0.340	-17.6	54	-0.04
35		Caprolactam	0.096	0.101	-5.2	45#	-0.06
36	C	4-Chloro-3-methylphenol	0.362	0.394	-8.8	48#	-0.04
37		2-Methylnaphthalene	0.770	0.820	-6.5	48#	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	49#	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.769	0.815	-6.0	53	-0.04
40	P	Hexachlorocyclopentadiene	0.444	0.471	-6.1	52	-0.04
41	S	2,4,6-Tribromophenol	0.304	0.326	-7.2	52	-0.03
42	C	2,4,6-Trichlorophenol	0.453	0.514	-13.5	54	-0.04
43		2,4,5-Trichlorophenol	0.501	0.555	-10.8	52	-0.04
44	S	2-Fluorobiphenyl	1.556	1.640	-5.4	52	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM052617\
Data File : BM010249.D
Acq On : 26 May 2017 10:23
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 26 23:52:57 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 20 04:25:49 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.645	1.663	-1.1	50	-0.04
46		2-Chloronaphthalene	1.332	1.346	-1.1	50	-0.04
47		2-Nitroaniline	0.342	0.404	-18.1	58	-0.04
48		Acenaphthylene	1.969	1.986	-0.9	49#	-0.04
49		Dimethylphthalate	1.784	1.851	-3.8	51	-0.04
50		2,6-Dinitrotoluene	0.339	0.356	-5.0	49#	-0.03
51	C	Acenaphthene	1.224	1.234	-0.8	49#	-0.04
52		3-Nitroaniline	0.315	0.308	2.2	48#	-0.03
53	P	2,4-Dinitrophenol	0.195	0.232	-19.0	60	-0.03
54		Dibenzofuran	1.925	1.969	-2.3	51	-0.04
55	P	4-Nitrophenol	0.253	0.228	9.9	44#	-0.04
56		2,4-Dinitrotoluene	0.479	0.514	-7.3	53	-0.03
57		Fluorene	1.674	1.748	-4.4	52	-0.04
58		2,3,4,6-Tetrachlorophenol	0.426	0.481	-12.9	55	-0.04
59		Diethylphthalate	1.687	1.767	-4.7	52	-0.04
60		4-Chlorophenyl-phenylether	0.940	1.001	-6.5	54	-0.04
61		4-Nitroaniline	0.320	0.294	8.1	45#	-0.04
62		Azobenzene	1.240	1.521	-22.7	59	-0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	52	-0.03
64		4,6-Dinitro-2-methylphenol	0.122	0.140	-14.8	58	-0.04
65	c	n-Nitrosodiphenylamine	0.599	0.618	-3.2	52	-0.04
66		4-Bromophenyl-phenylether	0.256	0.272	-6.3	55	-0.04
67		Hexachlorobenzene	0.311	0.313	-0.6	52	-0.04
68		Atrazine	0.227	0.234	-3.1	51	-0.04
69	C	Pentachlorophenol	0.171	0.195	-14.0	56	-0.03
70		Phenanthrene	1.119	1.154	-3.1	53	-0.04
71		Anthracene	1.110	1.166	-5.0	53	-0.04
72		Carbazole	0.983	1.005	-2.2	52	-0.04
73		Di-n-butylphthalate	1.166	1.207	-3.5	52	-0.04
74	C	Fluoranthene	1.453	1.508	-3.8	54	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	55	-0.03
76		Benzidine	0.369	0.297	19.5	40#	-0.02
77		Pyrene	1.072	1.078	-0.6	55	-0.03
78	S	Terphenyl-d14	0.880	0.923	-4.9	56	-0.03
79		Butylbenzylphthalate	0.397	0.380	4.3	51	-0.04
80		Benzo(a)anthracene	1.102	1.126	-2.2	56	-0.04
81		3,3'-Dichlorobenzidine	0.443	0.426	3.8	52	-0.03
82		Chrysene	1.045	1.077	-3.1	57	-0.03
83		Bis(2-ethylhexyl)phthalate	0.594	0.580	2.4	53	-0.04
84	c	Di-n-octyl phthalate	1.063	1.050	1.2	54	-0.05
85		Indeno(1,2,3-cd)pyrene	1.533	1.591	-3.8	58	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	55	-0.05
87		Benzo(b)fluoranthene	1.170	1.177	-0.6	54	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM052617\
Data File : BM010249.D
Acq On : 26 May 2017 10:23
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 26 23:52:57 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 20 04:25:49 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.147	1.179	-2.8	57	-0.05
89 C	Benzo(a)pyrene	1.115	1.133	-1.6	56	-0.05
90	Dibenzo(a,h)anthracene	1.247	1.263	-1.3	56	-0.08
91	Benzo(g,h,i)perylene	1.222	1.278	-4.6	58	-0.08

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