Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032217\
Data File : BM009335.D
Acq On : 23 Mar 2017 01:21
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:06:42 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 16:43:02 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	98	0.00
2		1,4-Dioxane	0.603	0.609	-1.0	103	0.00
3		Pyridine	1.705	1.751	-2.7	97	0.00
4		n-Nitrosodimethylamine	0.908	0.921	-1.4	98	0.00
5	S	2-Fluorophenol	1.200	1.224	-2.0	99	0.00
6		Aniline	2.224	2.235	-0.5	94	0.00
7	S	Phenol-d6	1.636	1.672	-2.2	95	0.00
8		2-Chlorophenol	1.217	1.235	-1.5	94	0.00
9		Benzaldehyde	1.288	1.251	2.9	93	0.00
10	C	Phenol	1.773	1.784	-0.6	95	0.00
11		bis(2-Chloroethyl)ether	1.453	1.464	-0.8	96	0.00
12		1,3-Dichlorobenzene	1.457	1.506	-3.4	98	0.00
13	C	1,4-Dichlorobenzene	1.505	1.554	-3.3	99	0.00
14		1,2-Dichlorobenzene	1.461	1.477	-1.1	95	0.00
15		Benzyl Alcohol	1.430	1.435	-0.3	96	0.00
16		2,2'-oxybis(1-Chloropropane	2.526	2.462	2.5	94	0.00
17		2-Methylphenol	1.160	1.186	-2.2	95	0.00
18		Hexachloroethane	0.616	0.612	0.6	93	0.00
19	P	n-Nitroso-di-n-propylamine	1.319	1.316	0.2	95	0.00
20		3+4-Methylphenols	1.577	1.586	-0.6	94	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.564	0.566	-0.4	92	0.00
23	S	Nitrobenzene-d5	0.533	0.543	-1.9	94	0.00
24		Nitrobenzene	0.523	0.532	-1.7	94	0.00
25		Isophorone	0.814	0.825	-1.4	93	0.00
26	C	2-Nitrophenol	0.158	0.170	-7.6	97	0.00
27		2,4-Dimethylphenol	0.288	0.292	-1.4	94	0.00
28		bis(2-Chloroethoxy)methane	0.504	0.499	1.0	93	0.00
29	C	2,4-Dichlorophenol	0.301	0.316	-5.0	95	0.00
30		1,2,4-Trichlorobenzene	0.374	0.386	-3.2	95	0.00
31		Naphthalene	1.062	1.076	-1.3	94	0.00
32		Benzoic acid	0.196	0.198	-1.0	94	0.00
33		4-Chloroaniline	0.409	0.420	-2.7	94	0.00
34	C	Hexachlorobutadiene	0.256	0.260	-1.6	96	0.00
35		Caprolactam	0.092	0.095	-3.3	95	0.00
36	C	4-Chloro-3-methylphenol	0.341	0.347	-1.8	92	0.00
37		2-Methylnaphthalene	0.733	0.735	-0.3	93	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.773	-2.4	95	0.00
40	P	Hexachlorocyclopentadiene	0.433	0.450	-3.9	94	0.00
41	S	2,4,6-Tribromophenol	0.248	0.265	-6.9	97	0.00
42	C	2,4,6-Trichlorophenol	0.432	0.449	-3.9	93	0.00
43		2,4,5-Trichlorophenol	0.482	0.506	-5.0	95	0.00
44	S	2-Fluorobiphenyl	1.660	1.686	-1.6	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032217\
Data File : BM009335.D
Acq On : 23 Mar 2017 01:21
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:06:42 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 16:43:02 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	1.665	1.699	-2.0	94	0.00
46		2-Chloronaphthalene	1.293	1.295	-0.2	93	0.00
47		2-Nitroaniline	0.476	0.508	-6.7	96	0.00
48		Acenaphthylene	2.106	2.187	-3.8	95	0.00
49		Dimethylphthalate	1.536	1.554	-1.2	95	0.00
50		2,6-Dinitrotoluene	0.331	0.348	-5.1	94	0.00
51	C	Acenaphthene	1.271	1.296	-2.0	95	0.00
52		3-Nitroaniline	0.326	0.343	-5.2	94	0.00
53	P	2,4-Dinitrophenol	0.195	0.196	-0.5	97	0.00
54		Dibenzofuran	1.879	1.918	-2.1	96	0.00
55	P	4-Nitrophenol	0.269	0.270	-0.4	92	0.00
56		2,4-Dinitrotoluene	0.448	0.474	-5.8	98	0.00
57		Fluorene	1.689	1.686	0.2	93	0.00
58		2,3,4,6-Tetrachlorophenol	0.407	0.426	-4.7	97	0.00
59		Diethylphthalate	1.558	1.579	-1.3	95	0.00
60		4-Chlorophenyl-phenylether	0.882	0.890	-0.9	95	0.00
61		4-Nitroaniline	0.354	0.373	-5.4	100	0.00
62		Azobenzene	1.819	1.877	-3.2	95	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.126	0.126	0.0	95	0.00
65	c	n-Nitrosodiphenylamine	0.608	0.634	-4.3	95	0.00
66		4-Bromophenyl-phenylether	0.232	0.242	-4.3	99	0.00
67		Hexachlorobenzene	0.254	0.254	0.0	95	0.00
68		Atrazine	0.228	0.241	-5.7	99	0.00
69	C	Pentachlorophenol	0.155	0.160	-3.2	99	0.00
70		Phenanthrene	1.144	1.162	-1.6	97	0.00
71		Anthracene	1.158	1.196	-3.3	97	0.00
72		Carbazole	1.109	1.163	-4.9	100	0.00
73		Di-n-butylphthalate	1.133	1.180	-4.1	97	0.00
74	C	Fluoranthene	1.355	1.421	-4.9	102	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
76		Benzidine	0.596	0.602	-1.0	100	0.00
77		Pyrene	1.134	1.149	-1.3	101	0.00
78	S	Terphenyl-d14	0.957	0.969	-1.3	100	0.00
79		Butylbenzylphthalate	0.408	0.427	-4.7	102	0.00
80		Benzo(a)anthracene	1.168	1.176	-0.7	104	0.00
81		3,3'-Dichlorobenzidine	0.437	0.462	-5.7	107	0.00
82		Chrysene	1.115	1.126	-1.0	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.610	0.612	-0.3	99	0.00
84	c	Di-n-octyl phthalate	1.014	1.064	-4.9	104	0.00
85		Indeno(1,2,3-cd)pyrene	1.216	1.252	-3.0	108	0.00
86	I	Perylene-d12	1.000	1.000	0.0	107	0.00
87		Benzo(b)fluoranthene	1.268	1.276	-0.6	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032217\
Data File : BM009335.D
Acq On : 23 Mar 2017 01:21
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:06:42 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 16:43:02 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	1.219	1.272	-4.3	108	0.00
89 C	Benzo(a)pyrene	1.174	1.219	-3.8	108	0.00
90	Dibenzo(a,h)anthracene	1.150	1.188	-3.3	108	0.00
91	Benzo(g,h,i)perylene	1.121	1.161	-3.6	108	0.00

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