Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024065.D
Acq On : 21 Sep 2016 22:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 96 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 22 11:54:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 19:01:00 2016
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	372#	0.00
2		1,4-Dioxane	0.470	0.437	7.0	316#	0.00
3		Pyridine	1.412	1.371	2.9	326#	0.00
4		n-Nitrosodimethylamine	0.744	0.737	0.9	369#	0.00
5	S	2-Fluorophenol	1.139	1.177	-3.3	357#	0.00
6		Aniline	2.328	2.300	1.2	355#	0.00
7	S	Phenol-d6	1.754	1.735	1.1	348#	0.01
8		2-Chlorophenol	1.320	1.266	4.1	341#	0.00
9		Benzaldehyde	1.246	1.229	1.4	346#	0.00
10	C	Phenol	1.845	1.832	0.7	344#	0.01
11		bis(2-Chloroethyl)ether	1.451	1.427	1.7	378#	0.00
12		1,3-Dichlorobenzene	1.545	1.536	0.6	370#	0.00
13	C	1,4-Dichlorobenzene	1.636	1.550	5.3	363#	0.00
14		1,2-Dichlorobenzene	1.536	1.452	5.5	353#	0.00
15		Benzyl Alcohol	1.546	1.590	-2.8	354#	0.00
16		2,2'-oxybis(1-Chloropropane	1.945	1.886	3.0	371#	0.01
17		2-Methylphenol	1.243	1.203	3.2	344#	0.00
18		Hexachloroethane	0.630	0.643	-2.1	374#	0.00
19	P	n-Nitroso-di-n-propylamine	1.549	1.434	7.4	341#	0.00
20		3+4-Methylphenols	1.732	1.671	3.5	330#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	294#	0.00
22		Acetophenone	0.518	0.563	-8.7	320#	0.00
23	S	Nitrobenzene-d5	0.448	0.509	-13.6	342#	0.00
24		Nitrobenzene	0.482	0.550	-14.1	347#	0.00
25		Isophorone	0.869	0.911	-4.8	313#	0.00
26	C	2-Nitrophenol	0.183	0.199	-8.7	326#	0.00
27		2,4-Dimethylphenol	0.311	0.323	-3.9	287#	0.00
28		bis(2-Chloroethoxy)methane	0.439	0.466	-6.2	323#	0.00
29	C	2,4-Dichlorophenol	0.330	0.353	-7.0	308#	0.00
30		1,2,4-Trichlorobenzene	0.362	0.391	-8.0	330#	0.00
31		Naphthalene	1.031	1.004	2.6	301#	0.00
32		Benzoic acid	0.253	0.241	4.7	279#	0.06
33		4-Chloroaniline	0.430	0.419	2.6	281#	0.00
34	C	Hexachlorobutadiene	0.272	0.314	-15.4	337#	0.00
35		Caprolactam	0.116	0.103	11.2	252#	0.03
36	C	4-Chloro-3-methylphenol	0.441	0.428	2.9	282#	0.01
37		2-Methylnaphthalene	0.784	0.725	7.5	270#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	242#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.664	0.770	-16.0	276#	0.00
40	P	Hexachlorocyclopentadiene	0.335	0.293	12.5	214#	0.00
41	S	2,4,6-Tribromophenol	0.360	0.313	13.1	199#	0.00
42	C	2,4,6-Trichlorophenol	0.443	0.488	-10.2	259#	0.00
43		2,4,5-Trichlorophenol	0.448	0.489	-9.2	251#	0.00
44	S	2-Fluorobiphenyl	1.395	1.383	0.9	243#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024065.D
Acq On : 21 Sep 2016 22:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 96 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 22 11:54:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 19:01:00 2016
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	1.609	1.643	-2.1	248#	0.00
46		2-Chloronaphthalene	1.181	1.229	-4.1	256#	0.00
47		2-Nitroaniline	0.465	0.554	-19.1	291#	0.00
48		Acenaphthylene	1.969	1.899	3.6	232#	0.00
49		Dimethylphthalate	1.715	1.601	6.6	228#	0.00
50		2,6-Dinitrotoluene	0.362	0.353	2.5	235#	0.00
51	C	Acenaphthene	1.217	1.185	2.6	240#	0.00
52		3-Nitroaniline	0.339	0.353	-4.1	238#	0.00
53	P	2,4-Dinitrophenol	0.226	0.157	30.5#	158#	0.00
54		Dibenzofuran	1.900	1.754	7.7	223#	0.00
55	P	4-Nitrophenol	0.269	0.363	-34.9#	336#	0.00
56		2,4-Dinitrotoluene	0.532	0.507	4.7	227#	0.00
57		Fluorene	1.647	1.464	11.1	218#	0.00
58		2,3,4,6-Tetrachlorophenol	0.455	0.416	8.6	217#	0.00
59		Diethylphthalate	1.831	1.613	11.9	213#	0.00
60		4-Chlorophenyl-phenylether	0.887	0.825	7.0	220#	0.00
61		4-Nitroaniline	0.342	0.335	2.0	232#	0.01
62		Azobenzene	1.773	1.675	5.5	231#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	201#	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.080	42.4#	113	0.01
65	c	n-Nitrosodiphenylamine	0.568	0.592	-4.2	209#	0.00
66		4-Bromophenyl-phenylether	0.243	0.258	-6.2	209#	0.00
67		Hexachlorobenzene	0.285	0.285	0.0	199#	0.00
68		Atrazine	0.244	0.251	-2.9	201#	0.00
69	C	Pentachlorophenol	0.152	0.223	-46.7#	279#	0.00
70		Phenanthrene	1.040	0.999	3.9	197#	0.00
71		Anthracene	1.061	1.006	5.2	193#	0.00
72		Carbazole	0.963	0.908	5.7	192#	0.00
73		Di-n-butylphthalate	1.149	1.090	5.1	195#	0.00
74	C	Fluoranthene	1.253	1.115	11.0	177#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	189#	0.00
76		Benzidine	0.619	0.686	-10.8	204#	0.00
77		Pyrene	1.230	1.107	10.0	171#	0.00
78	S	Terphenyl-d14	0.878	0.788	10.3	171#	0.00
79		Butylbenzylphthalate	0.474	0.488	-3.0	207#	0.00
80		Benzo(a)anthracene	1.120	1.116	0.4	192#	0.00
81		3,3'-Dichlorobenzidine	0.448	0.462	-3.1	192#	0.00
82		Chrysene	1.066	1.049	1.6	191#	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.676	-4.2	213#	0.00
84	c	Di-n-octyl phthalate	1.065	1.135	-6.6	216#	0.00
85		Indeno(1,2,3-cd)pyrene	1.180	1.073	9.1	182#	0.04
86	I	Perylene-d12	1.000	1.000	0.0	197#	0.00
87		Benzo(b)fluoranthene	1.136	1.135	0.1	200#	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024065.D
Acq On : 21 Sep 2016 22:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 96 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 22 11:54:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 19:01:00 2016
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	1.127	1.184	-5.1	208#	0.02
89 C	Benzo(a)pyrene	1.079	1.083	-0.4	202#	0.01
90	Dibenzo(a,h)anthracene	1.061	0.910	14.2	172#	0.04
91	Benzo(g,h,i)perylene	1.020	0.791	22.5	156#	0.05

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