Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024072.D
Acq On : 22 Sep 2016 15:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 23 01:25:52 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	152#	0.00
2		1,4-Dioxane	0.470	0.444	5.5	131	0.00
3		Pyridine	1.412	1.571	-11.3	153#	-0.02
4		n-Nitrosodimethylamine	0.744	0.785	-5.5	161#	0.00
5	S	2-Fluorophenol	1.139	1.203	-5.6	149	0.00
6		Aniline	2.328	2.608	-12.0	165#	0.00
7	S	Phenol-d6	1.754	1.934	-10.3	159#	0.00
8		2-Chlorophenol	1.320	1.337	-1.3	147	0.00
9		Benzaldehyde	1.246	1.216	2.4	140	0.00
10	C	Phenol	1.845	2.010	-8.9	154#	0.00
11		bis(2-Chloroethyl)ether	1.451	1.509	-4.0	163#	0.00
12		1,3-Dichlorobenzene	1.545	1.568	-1.5	154#	0.00
13	C	1,4-Dichlorobenzene	1.636	1.559	4.7	150	0.00
14		1,2-Dichlorobenzene	1.536	1.489	3.1	148	0.00
15		Benzyl Alcohol	1.546	1.741	-12.6	159#	0.00
16		2,2'-oxybis(1-Chloropropane	1.945	2.080	-6.9	167#	0.00
17		2-Methylphenol	1.243	1.364	-9.7	159#	0.00
18		Hexachloroethane	0.630	0.630	0.0	150#	0.00
19	P	n-Nitroso-di-n-propylamine	1.549	1.737	-12.1	169#	0.00
20		3+4-Methylphenols	1.732	1.909	-10.2	154#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
22		Acetophenone	0.518	0.576	-11.2	153#	0.00
23	S	Nitrobenzene-d5	0.448	0.514	-14.7	162#	0.00
24		Nitrobenzene	0.482	0.549	-13.9	162#	0.00
25		Isophorone	0.869	1.020	-17.4	164#	0.00
26	C	2-Nitrophenol	0.183	0.194	-6.0	148	0.00
27		2,4-Dimethylphenol	0.311	0.329	-5.8	137	0.00
28		bis(2-Chloroethoxy)methane	0.439	0.504	-14.8	163#	0.00
29	C	2,4-Dichlorophenol	0.330	0.356	-7.9	145	0.00
30		1,2,4-Trichlorobenzene	0.362	0.379	-4.7	150	0.00
31		Naphthalene	1.031	1.031	0.0	144	0.00
32		Benzoic acid	0.253	0.275	-8.7	148	0.00
33		4-Chloroaniline	0.430	0.454	-5.6	142	0.00
34	C	Hexachlorobutadiene	0.272	0.295	-8.5	148	0.00
35		Caprolactam	0.116	0.130	-12.1	148	0.00
36	C	4-Chloro-3-methylphenol	0.441	0.495	-12.2	152#	0.00
37		2-Methylnaphthalene	0.784	0.782	0.3	136	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	133	0.00
39		1,2,4,5-Tetrachlorobenzene	0.664	0.706	-6.3	139	0.00
40	P	Hexachlorocyclopentadiene	0.335	0.310	7.5	125	0.00
41	S	2,4,6-Tribromophenol	0.360	0.337	6.4	118	0.00
42	C	2,4,6-Trichlorophenol	0.443	0.452	-2.0	132	0.00
43		2,4,5-Trichlorophenol	0.448	0.463	-3.3	131	0.00
44	S	2-Fluorobiphenyl	1.395	1.386	0.6	134	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024072.D
Acq On : 22 Sep 2016 15:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 23 01:25:52 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	1.598	0.7	133	0.00
46		2-Chloronaphthalene	1.181	1.169	1.0	134	0.00
47		2-Nitroaniline	0.465	0.559	-20.2	162#	0.00
48		Acenaphthylene	1.969	1.977	-0.4	133	0.00
49		Dimethylphthalate	1.715	1.750	-2.0	137	0.00
50		2,6-Dinitrotoluene	0.362	0.379	-4.7	139	0.00
51	C	Acenaphthene	1.217	1.185	2.6	132	0.00
52		3-Nitroaniline	0.339	0.383	-13.0	142	0.00
53	P	2,4-Dinitrophenol	0.226	0.238	-5.3	132	0.00
54		Dibenzofuran	1.900	1.870	1.6	131	0.00
55	P	4-Nitrophenol	0.269	0.281	-4.5	144	0.00
56		2,4-Dinitrotoluene	0.532	0.562	-5.6	139	0.00
57		Fluorene	1.647	1.602	2.7	131	0.00
58		2,3,4,6-Tetrachlorophenol	0.455	0.455	0.0	131	0.00
59		Diethylphthalate	1.831	1.820	0.6	132	0.00
60		4-Chlorophenyl-phenylether	0.887	0.877	1.1	129	0.00
61		4-Nitroaniline	0.342	0.377	-10.2	144	0.00
62		Azobenzene	1.773	1.826	-3.0	139	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.148	-6.5	135	0.00
65	c	n-Nitrosodiphenylamine	0.568	0.575	-1.2	130	0.00
66		4-Bromophenyl-phenylether	0.243	0.239	1.6	124	0.00
67		Hexachlorobenzene	0.285	0.272	4.6	122	0.00
68		Atrazine	0.244	0.256	-4.9	131	0.00
69	C	Pentachlorophenol	0.152	0.141	7.2	113	0.00
70		Phenanthrene	1.040	1.023	1.6	129	0.00
71		Anthracene	1.061	1.054	0.7	130	0.00
72		Carbazole	0.963	0.958	0.5	130	0.00
73		Di-n-butylphthalate	1.149	1.214	-5.7	139	0.00
74	C	Fluoranthene	1.253	1.248	0.4	127	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	122	0.00
76		Benzidine	0.619	0.650	-5.0	124	0.00
77		Pyrene	1.230	1.258	-2.3	125	0.00
78	S	Terphenyl-d14	0.878	0.885	-0.8	123	0.00
79		Butylbenzylphthalate	0.474	0.518	-9.3	141	0.00
80		Benzo(a)anthracene	1.120	1.159	-3.5	128	0.00
81		3,3'-Dichlorobenzidine	0.448	0.464	-3.6	124	0.00
82		Chrysene	1.066	1.108	-3.9	130	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.708	-9.1	144	0.00
84	c	Di-n-octyl phthalate	1.065	1.196	-12.3	146	0.00
85		Indeno(1,2,3-cd)pyrene	1.180	1.352	-14.6	147	0.05
86	I	Perylene-d12	1.000	1.000	0.0	133	0.00
87		Benzo(b)fluoranthene	1.136	1.166	-2.6	138	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
Data File : BG024072.D
Acq On : 22 Sep 2016 15:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 23 01:25:52 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	1.122	0.4	132	0.01
89 C	Benzo(a)pyrene	1.079	1.099	-1.9	138	0.01
90	Dibenzo(a,h)anthracene	1.061	1.089	-2.6	138	0.03
91	Benzo(g,h,i)perylene	1.020	1.089	-6.8	145	0.04

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