Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062816\
Data File : BG022669.D
Acq On : 28 Jun 2016 10:09
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 28 15:06:01 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	102	-0.01
2		1,4-Dioxane	0.505	0.505	0.0	100	-0.01
3		Pyridine	1.414	1.648	-16.5	116	-0.02
4		n-Nitrosodimethylamine	0.809	0.713	11.9	87	-0.02
5	S	2-Fluorophenol	1.247	1.220	2.2	100	0.00
6		Aniline	2.330	2.581	-10.8	111	0.00
7	S	Phenol-d6	1.758	1.876	-6.7	109	0.00
8		2-Chlorophenol	1.292	1.322	-2.3	107	0.00
9		Benzaldehyde	0.619	0.747	-20.7	121	0.00
10	C	Phenol	1.858	1.990	-7.1	108	0.00
11		bis(2-Chloroethyl)ether	1.529	1.485	2.9	94	0.00
12		1,3-Dichlorobenzene	1.572	1.550	1.4	100	0.00
13	C	1,4-Dichlorobenzene	1.576	1.583	-0.4	104	-0.01
14		1,2-Dichlorobenzene	1.498	1.475	1.5	103	0.00
15		Benzyl Alcohol	1.626	1.812	-11.4	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.654	2.2	99	0.00
17		2-Methylphenol	1.224	1.322	-8.0	113	0.00
18		Hexachloroethane	0.652	0.620	4.9	95	-0.01
19	P	n-Nitroso-di-n-propylamine	1.464	1.516	-3.6	105	0.00
20		3+4-Methylphenols	1.687	1.798	-6.6	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.578	0.565	2.2	103	0.00
23	S	Nitrobenzene-d5	0.473	0.465	1.7	104	0.00
24		Nitrobenzene	0.522	0.509	2.5	106	0.00
25		Isophorone	0.917	0.894	2.5	106	0.00
26	C	2-Nitrophenol	0.188	0.196	-4.3	114	0.00
27		2,4-Dimethylphenol	0.359	0.334	7.0	105	0.00
28		bis(2-Chloroethoxy)methane	0.501	0.499	0.4	110	0.00
29	C	2,4-Dichlorophenol	0.350	0.370	-5.7	112	0.00
30		1,2,4-Trichlorobenzene	0.416	0.410	1.4	107	0.00
31		Naphthalene	1.005	0.999	0.6	107	0.00
32		Benzoic acid	0.216	0.243	-12.5	116	0.06
33		4-Chloroaniline	0.433	0.478	-10.4	112	-0.01
34	C	Hexachlorobutadiene	0.323	0.310	4.0	102	-0.01
35		Caprolactam	0.110	0.127	-15.5	123	0.02
36	C	4-Chloro-3-methylphenol	0.430	0.463	-7.7	116	0.00
37		2-Methylnaphthalene	0.780	0.773	0.9	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.739	2.1	117	0.00
40	P	Hexachlorocyclopentadiene	0.603	0.522	13.4	98	-0.01
41	S	2,4,6-Tribromophenol	0.315	0.322	-2.2	119	0.00
42	C	2,4,6-Trichlorophenol	0.484	0.492	-1.7	115	0.00
43		2,4,5-Trichlorophenol	0.507	0.522	-3.0	121	0.00
44	S	2-Fluorobiphenyl	1.261	1.243	1.4	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062816\
Data File : BG022669.D
Acq On : 28 Jun 2016 10:09
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 28 15:06:01 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	1.524	1.448	5.0	111	0.00
46		2-Chloronaphthalene	1.254	1.211	3.4	113	0.00
47		2-Nitroaniline	0.484	0.492	-1.7	116	0.00
48		Acenaphthylene	1.818	1.757	3.4	112	0.00
49		Dimethylphthalate	1.659	1.589	4.2	113	0.00
50		2,6-Dinitrotoluene	0.361	0.375	-3.9	119	0.00
51	C	Acenaphthene	1.339	1.321	1.3	113	0.00
52		3-Nitroaniline	0.337	0.345	-2.4	117	0.00
53	P	2,4-Dinitrophenol	0.222	0.230	-3.6	116	0.00
54		Dibenzofuran	1.940	1.855	4.4	113	0.00
55	P	4-Nitrophenol	0.285	0.294	-3.2	122	0.01
56		2,4-Dinitrotoluene	0.499	0.541	-8.4	125	0.00
57		Fluorene	1.548	1.522	1.7	115	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.562	-7.0	125	0.00
59		Diethylphthalate	1.706	1.617	5.2	113	0.00
60		4-Chlorophenyl-phenylether	0.922	0.903	2.1	116	0.00
61		4-Nitroaniline	0.348	0.348	0.0	116	0.02
62		Azobenzene	1.847	1.733	6.2	109	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.140	-4.5	118	0.00
65	c	n-Nitrosodiphenylamine	0.582	0.568	2.4	116	0.00
66		4-Bromophenyl-phenylether	0.259	0.247	4.6	115	0.00
67		Hexachlorobenzene	0.274	0.272	0.7	122	0.00
68		Atrazine	0.246	0.245	0.4	119	0.00
69	C	Pentachlorophenol	0.189	0.182	3.7	117	0.00
70		Phenanthrene	1.020	0.966	5.3	115	0.00
71		Anthracene	1.050	0.997	5.0	115	0.00
72		Carbazole	0.890	0.852	4.3	114	0.01
73		Di-n-butylphthalate	1.036	0.998	3.7	111	0.00
74	C	Fluoranthene	1.175	1.149	2.2	114	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	119	0.00
76		Benzidine	0.213	0.337	-58.2#	193#	-0.02
77		Pyrene	1.065	1.070	-0.5	112	0.00
78	S	Terphenyl-d14	0.755	0.665	11.9	111	0.00
79		Butylbenzylphthalate	0.461	0.456	1.1	112	0.00
80		Benzo(a)anthracene	1.107	1.115	-0.7	114	0.00
81		3,3'-Dichlorobenzidine	0.457	0.437	4.4	110	0.01
82		Chrysene	1.040	1.054	-1.3	115	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.606	6.6	109	0.00
84	c	Di-n-octyl phthalate	1.070	1.010	5.6	110	-0.01
85		Indeno(1,2,3-cd)pyrene	1.367	1.319	3.5	114	0.01
86	I	Perylene-d12	1.000	1.000	0.0	117	0.00
87		Benzo(b)fluoranthene	1.203	1.163	3.3	112	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062816\
Data File : BG022669.D
Acq On : 28 Jun 2016 10:09
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 28 15:06:01 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	1.137	1.122	1.3	116	0.01
89 C	Benzo(a)pyrene	1.172	1.142	2.6	114	0.01
90	Dibenzo(a,h)anthracene	1.104	1.069	3.2	116	0.02
91	Benzo(g,h,i)perylene	1.123	1.019	9.3	109	0.02

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