Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121517\
Data File : BG031593.D
Acq On : 15 Dec 2017 18:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 15 23:00:46 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	137	0.00
2		1,4-Dioxane	0.538	0.485	9.9	126	-0.01
3		Pyridine	1.417	1.415	0.1	130	-0.01
4		n-Nitrosodimethylamine	0.668	0.643	3.7	125	-0.01
5	S	2-Fluorophenol	1.128	1.171	-3.8	135	-0.01
6		Aniline	2.063	2.082	-0.9	134	-0.01
7	S	Phenol-d6	1.566	1.614	-3.1	136	0.00
8		2-Chlorophenol	1.259	1.319	-4.8	138	0.00
9		Benzaldehyde	0.908	0.889	2.1	131	-0.01
10	C	Phenol	1.609	1.628	-1.2	132	0.00
11		bis(2-Chloroethyl)ether	1.313	1.325	-0.9	132	0.00
12		1,3-Dichlorobenzene	1.497	1.495	0.1	134	0.00
13	C	1,4-Dichlorobenzene	1.556	1.532	1.5	133	-0.01
14		1,2-Dichlorobenzene	1.482	1.472	0.7	136	-0.01
15		Benzyl Alcohol	1.225	1.213	1.0	133	0.00
16		2,2'-oxybis(1-Chloropropane	1.475	1.373	6.9	126	0.00
17		2-Methylphenol	1.097	1.139	-3.8	136	-0.01
18		Hexachloroethane	0.524	0.530	-1.1	136	-0.01
19	P	n-Nitroso-di-n-propylamine	1.233	1.169	5.2	127	-0.01
20		3+4-Methylphenols	1.634	1.618	1.0	134	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
22		Acetophenone	0.480	0.503	-4.8	130	-0.01
23	S	Nitrobenzene-d5	0.324	0.338	-4.3	128	0.00
24		Nitrobenzene	0.333	0.348	-4.5	128	0.00
25		Isophorone	0.613	0.648	-5.7	127	-0.01
26	C	2-Nitrophenol	0.164	0.194	-18.3	146	0.00
27		2,4-Dimethylphenol	0.250	0.276	-10.4	135	-0.01
28		bis(2-Chloroethoxy)methane	0.396	0.424	-7.1	133	0.00
29	C	2,4-Dichlorophenol	0.294	0.336	-14.3	136	0.00
30		1,2,4-Trichlorobenzene	0.376	0.389	-3.5	133	-0.01
31		Naphthalene	0.983	0.995	-1.2	130	-0.01
32		Benzoic acid	0.132	0.091	31.1#	89	0.00
33		4-Chloroaniline	0.427	0.447	-4.7	131	-0.01
34	C	Hexachlorobutadiene	0.249	0.260	-4.4	137	-0.01
35		Caprolactam	0.116	0.122	-5.2	130	0.00
36	C	4-Chloro-3-methylphenol	0.336	0.352	-4.8	132	0.00
37		2-Methylnaphthalene	0.761	0.778	-2.2	129	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
39		1,2,4,5-Tetrachlorobenzene	0.782	0.815	-4.2	128	0.00
40	P	Hexachlorocyclopentadiene	0.182	0.245	-34.6#	163#	-0.01
41	S	2,4,6-Tribromophenol	0.234	0.266	-13.7	136	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.457	-14.5	134	0.00
43		2,4,5-Trichlorophenol	0.430	0.490	-14.0	133	0.00
44	S	2-Fluorobiphenyl	1.363	1.383	-1.5	126	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121517\
Data File : BG031593.D
Acq On : 15 Dec 2017 18:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 15 23:00:46 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	1.643	1.694	-3.1	127	0.00
46		2-Chloronaphthalene	1.202	1.217	-1.2	125	0.00
47		2-Nitroaniline	0.338	0.350	-3.6	125	0.00
48		Acenaphthylene	1.935	1.973	-2.0	125	0.00
49		Dimethylphthalate	1.656	1.735	-4.8	128	0.00
50		2,6-Dinitrotoluene	0.354	0.376	-6.2	129	0.00
51	C	Acenaphthene	1.223	1.240	-1.4	126	0.00
52		3-Nitroaniline	0.361	0.384	-6.4	131	0.00
53	P	2,4-Dinitrophenol	0.116	0.130	-12.1	130	0.00
54		Dibenzofuran	1.952	1.897	2.8	121	0.00
55	P	4-Nitrophenol	0.222	0.239	-7.7	128	0.00
56		2,4-Dinitrotoluene	0.503	0.522	-3.8	126	0.00
57		Fluorene	1.564	1.542	1.4	122	0.00
58		2,3,4,6-Tetrachlorophenol	0.404	0.470	-16.3	135	0.00
59		Diethylphthalate	1.660	1.684	-1.4	124	0.00
60		4-Chlorophenyl-phenylether	0.957	0.931	2.7	120	0.00
61		4-Nitroaniline	0.379	0.401	-5.8	129	0.00
62		Azobenzene	1.375	1.276	7.2	113	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	123	0.00
64		4,6-Dinitro-2-methylphenol	0.112	0.107	4.5	111	0.00
65	c	n-Nitrosodiphenylamine	0.585	0.605	-3.4	123	0.00
66		4-Bromophenyl-phenylether	0.233	0.247	-6.0	128	0.00
67		Hexachlorobenzene	0.240	0.253	-5.4	129	0.00
68		Atrazine	0.214	0.230	-7.5	126	0.00
69	C	Pentachlorophenol	0.100	0.101	-1.0	116	0.00
70		Phenanthrene	1.045	1.050	-0.5	124	0.00
71		Anthracene	1.033	1.059	-2.5	124	0.00
72		Carbazole	0.935	0.994	-6.3	128	0.00
73		Di-n-butylphthalate	1.080	1.158	-7.2	127	0.00
74	C	Fluoranthene	1.307	1.349	-3.2	126	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	122	0.00
76		Benzidine	0.465	0.472	-1.5	114	0.00
77		Pyrene	1.243	1.285	-3.4	122	0.00
78	S	Terphenyl-d14	0.879	0.902	-2.6	123	0.00
79		Butylbenzylphthalate	0.436	0.495	-13.5	129	0.00
80		Benzo(a)anthracene	1.207	1.250	-3.6	123	0.00
81		3,3'-Dichlorobenzidine	0.420	0.448	-6.7	121	0.00
82		Chrysene	1.157	1.155	0.2	119	0.00
83		Bis(2-ethylhexyl)phthalate	0.627	0.666	-6.2	122	-0.01
84	c	Di-n-octyl phthalate	1.035	1.055	-1.9	115	-0.01
85		Indeno(1,2,3-cd)pyrene	1.237	1.123	9.2	106	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	111	0.00
87		Benzo(b)fluoranthene	1.196	1.255	-4.9	116	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG121517\
Data File : BG031593.D
Acq On : 15 Dec 2017 18:20
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 15 23:00:46 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 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	1.220	1.256	-3.0	113	0.00
89 C	Benzo(a)pyrene	1.134	1.160	-2.3	111	0.00
90	Dibenzo(a,h)anthracene	1.085	1.066	1.8	107	-0.02
91	Benzo(g,h,i)perylene	1.066	1.030	3.4	105	-0.01

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