Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071417\
Data File : BG027926.D
Acq On : 14 Jul 2017 15:42
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071417

Quant Time: Jul 14 16:16:02 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 14 16:10:34 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	97	0.01
2		1,4-Dioxane	0.416	0.416	0.0	100	0.00
3		Pyridine	1.194	1.264	-5.9	102	0.00
4		n-Nitrosodimethylamine	0.671	0.684	-1.9	100	0.00
5	S	2-Fluorophenol	1.118	1.142	-2.1	101	0.00
6		Aniline	1.875	1.908	-1.8	98	0.00
7	S	Phenol-d6	1.611	1.654	-2.7	100	0.00
8		2-Chlorophenol	1.279	1.292	-1.0	101	0.00
9		Benzaldehyde	0.882	0.927	-5.1	102	0.00
10	C	Phenol	1.689	1.734	-2.7	101	0.00
11		bis(2-Chloroethyl)ether	1.238	1.245	-0.6	102	0.00
12		1,3-Dichlorobenzene	1.498	1.509	-0.7	102	0.00
13	C	1,4-Dichlorobenzene	1.549	1.562	-0.8	100	0.00
14		1,2-Dichlorobenzene	1.504	1.511	-0.5	100	0.00
15		Benzyl Alcohol	0.984	1.068	-8.5	102	0.00
16		2,2'-oxybis(1-Chloropropane	2.498	2.453	1.8	99	0.00
17		2-Methylphenol	1.111	1.126	-1.4	100	0.00
18		Hexachloroethane	0.504	0.512	-1.6	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.125	1.145	-1.8	97	0.00
20		3+4-Methylphenols	1.560	1.587	-1.7	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.01
22		Acetophenone	0.482	0.494	-2.5	100	0.00
23	S	Nitrobenzene-d5	0.358	0.373	-4.2	100	0.00
24		Nitrobenzene	0.376	0.383	-1.9	97	0.01
25		Isophorone	0.684	0.712	-4.1	98	0.01
26	C	2-Nitrophenol	0.181	0.190	-5.0	99	0.00
27		2,4-Dimethylphenol	0.317	0.327	-3.2	97	0.00
28		bis(2-Chloroethoxy)methane	0.400	0.410	-2.5	100	0.01
29	C	2,4-Dichlorophenol	0.343	0.354	-3.2	100	0.00
30		1,2,4-Trichlorobenzene	0.385	0.388	-0.8	98	0.00
31		Naphthalene	0.997	1.007	-1.0	98	0.00
32		Benzoic acid	0.154	0.178	-15.6	114	0.01
33		4-Chloroaniline	0.408	0.421	-3.2	98	0.00
34	C	Hexachlorobutadiene	0.269	0.276	-2.6	101	0.00
35		Caprolactam	0.105	0.113	-7.6	100	0.00
36	C	4-Chloro-3-methylphenol	0.370	0.386	-4.3	99	0.00
37		2-Methylnaphthalene	0.761	0.769	-1.1	98	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
39		1,2,4,5-Tetrachlorobenzene	0.791	0.806	-1.9	99	0.00
40	P	Hexachlorocyclopentadiene	0.369	0.396	-7.3	99	0.00
41	S	2,4,6-Tribromophenol	0.296	0.316	-6.8	100	0.00
42	C	2,4,6-Trichlorophenol	0.479	0.507	-5.8	99	0.00
43		2,4,5-Trichlorophenol	0.505	0.533	-5.5	98	0.00
44	S	2-Fluorobiphenyl	1.517	1.569	-3.4	99	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071417\
Data File : BG027926.D
Acq On : 14 Jul 2017 15:42
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071417

Quant Time: Jul 14 16:16:02 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 14 16:10:34 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	1.719	1.765	-2.7	98	0.01
46		2-Chloronaphthalene	1.251	1.270	-1.5	98	0.00
47		2-Nitroaniline	0.374	0.400	-7.0	100	0.00
48		Acenaphthylene	1.991	2.047	-2.8	98	0.00
49		Dimethylphthalate	1.670	1.730	-3.6	98	0.00
50		2,6-Dinitrotoluene	0.357	0.374	-4.8	98	0.00
51	C	Acenaphthene	1.249	1.257	-0.6	92	0.00
52		3-Nitroaniline	0.338	0.364	-7.7	98	0.00
53	P	2,4-Dinitrophenol	0.202	0.217	-7.4	98	0.00
54		Dibenzofuran	1.909	1.945	-1.9	99	0.00
55	P	4-Nitrophenol	0.253	0.274	-8.3	102	0.00
56		2,4-Dinitrotoluene	0.493	0.527	-6.9	100	0.00
57		Fluorene	1.706	1.760	-3.2	98	0.00
58		2,3,4,6-Tetrachlorophenol	0.448	0.479	-6.9	99	0.00
59		Diethylphthalate	1.619	1.675	-3.5	98	0.00
60		4-Chlorophenyl-phenylether	0.941	0.972	-3.3	99	0.00
61		4-Nitroaniline	0.361	0.384	-6.4	100	0.00
62		Azobenzene	1.269	1.315	-3.6	98	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.133	0.140	-5.3	102	0.00
65	c	n-Nitrosodiphenylamine	0.584	0.596	-2.1	98	0.00
66		4-Bromophenyl-phenylether	0.249	0.253	-1.6	97	0.00
67		Hexachlorobenzene	0.272	0.278	-2.2	99	0.00
68		Atrazine	0.224	0.238	-6.2	99	0.00
69	C	Pentachlorophenol	0.152	0.162	-6.6	103	0.00
70		Phenanthrene	1.046	1.071	-2.4	100	0.00
71		Anthracene	1.042	1.074	-3.1	100	0.00
72		Carbazole	0.938	0.975	-3.9	100	0.00
73		Di-n-butylphthalate	1.030	1.091	-5.9	100	0.00
74	C	Fluoranthene	1.256	1.311	-4.4	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
76		Benzidine	0.467	0.415	11.1	77	0.00
77		Pyrene	1.069	1.102	-3.1	99	0.00
78	S	Terphenyl-d14	0.885	0.920	-4.0	101	0.00
79		Butylbenzylphthalate	0.384	0.413	-7.6	102	0.00
80		Benzo(a)anthracene	1.099	1.136	-3.4	100	0.00
81		3,3'-Dichlorobenzidine	0.432	0.444	-2.8	97	0.00
82		Chrysene	1.052	1.072	-1.9	100	0.00
83		Bis(2-ethylhexyl)phthalate	0.559	0.591	-5.7	100	0.00
84	c	Di-n-octyl phthalate	0.949	1.009	-6.3	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.216	1.271	-4.5	100	0.00
86	I	Perylene-d12	1.000	1.000	0.0	97	0.00
87		Benzo(b)fluoranthene	1.169	1.196	-2.3	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071417\
Data File : BG027926.D
Acq On : 14 Jul 2017 15:42
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071417

Quant Time: Jul 14 16:16:02 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 14 16:10:34 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	1.168	1.209	-3.5	100	0.00
89 C	Benzo(a)pyrene	1.102	1.135	-3.0	100	0.01
90	Dibenzo(a,h)anthracene	1.113	1.146	-3.0	100	0.00
91	Benzo(g,h,i)perylene	1.073	1.106	-3.1	101	0.01

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