Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017926.D
Acq On : 17 Jul 2015 15:35
Operator : TP/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071715

Quant Time: Jul 18 02:18:23 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 2015
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	107	0.00
2		1,4-Dioxane	0.504	0.473	6.2	107	0.00
3		Pyridine	1.347	1.335	0.9	106	0.00
4		n-Nitrosodimethylamine	0.511	0.480	6.1	105	0.00
5	S	2-Fluorophenol	1.146	1.127	1.7	107	0.00
6		Aniline	2.003	1.952	2.5	108	0.00
7	S	Phenol-d6	1.502	1.459	2.9	107	0.00
8		2-Chlorophenol	1.352	1.325	2.0	109	0.00
9		Benzaldehyde	0.920	0.895	2.7	108	0.00
10	C	Phenol	1.603	1.565	2.4	109	0.00
11		bis(2-Chloroethyl)ether	1.254	1.209	3.6	108	0.00
12		1,3-Dichlorobenzene	1.484	1.428	3.8	107	0.00
13	C	1,4-Dichlorobenzene	1.521	1.453	4.5	107	0.00
14		1,2-Dichlorobenzene	1.454	1.386	4.7	108	0.00
15		Benzyl Alcohol	1.111	1.103	0.7	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.511	1.423	5.8	108	0.00
17		2-Methylphenol	1.118	1.091	2.4	109	0.00
18		Hexachloroethane	0.575	0.549	4.5	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.097	1.067	2.7	108	0.00
20		3+4-Methylphenols	1.513	1.507	0.4	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.432	0.417	3.5	109	0.00
23	S	Nitrobenzene-d5	0.309	0.303	1.9	109	0.00
24		Nitrobenzene	0.329	0.322	2.1	107	0.00
25		Isophorone	0.648	0.637	1.7	108	0.00
26	C	2-Nitrophenol	0.157	0.158	-0.6	114	0.00
27		2,4-Dimethylphenol	0.286	0.281	1.7	110	0.00
28		bis(2-Chloroethoxy)methane	0.361	0.351	2.8	107	0.00
29	C	2,4-Dichlorophenol	0.254	0.259	-2.0	109	0.00
30		1,2,4-Trichlorobenzene	0.302	0.291	3.6	110	0.00
31		Naphthalene	0.979	0.932	4.8	107	0.00
32		Benzoic acid	0.110	0.114	-3.6	130	0.00
33		4-Chloroaniline	0.436	0.423	3.0	106	0.00
34	C	Hexachlorobutadiene	0.170	0.165	2.9	107	0.00
35		Caprolactam	0.129	0.129	0.0	112	0.00
36	C	4-Chloro-3-methylphenol	0.289	0.289	0.0	109	0.00
37		2-Methylnaphthalene	0.703	0.666	5.3	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.536	0.521	2.8	106	0.00
40	P	Hexachlorocyclopentadiene	0.292	0.292	0.0	111	0.00
41	S	2,4,6-Tribromophenol	0.203	0.197	3.0	108	0.00
42	C	2,4,6-Trichlorophenol	0.349	0.344	1.4	109	0.00
43		2,4,5-Trichlorophenol	0.363	0.365	-0.6	108	0.00
44	S	2-Fluorobiphenyl	1.154	1.127	2.3	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017926.D
Acq On : 17 Jul 2015 15:35
Operator : TP/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071715

Quant Time: Jul 18 02:18:23 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 2015
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	1.400	1.370	2.1	106	0.00
46		2-Chloronaphthalene	1.135	1.097	3.3	106	0.00
47		2-Nitroaniline	0.305	0.303	0.7	109	0.00
48		Acenaphthylene	1.892	1.852	2.1	105	0.00
49		Dimethylphthalate	1.389	1.341	3.5	104	0.00
50		2,6-Dinitrotoluene	0.315	0.311	1.3	108	0.00
51	C	Acenaphthene	1.146	1.090	4.9	105	0.00
52		3-Nitroaniline	0.354	0.365	-3.1	108	0.00
53	P	2,4-Dinitrophenol	0.121	0.127	-5.0	119	0.00
54		Dibenzofuran	1.654	1.594	3.6	106	0.00
55	P	4-Nitrophenol	0.280	0.279	0.4	108	0.00
56		2,4-Dinitrotoluene	0.422	0.415	1.7	107	0.00
57		Fluorene	1.421	1.348	5.1	104	0.00
58		2,3,4,6-Tetrachlorophenol	0.313	0.304	2.9	106	0.00
59		Diethylphthalate	1.439	1.387	3.6	105	0.00
60		4-Chlorophenyl-phenylether	0.697	0.665	4.6	104	0.00
61		4-Nitroaniline	0.415	0.392	5.5	104	0.00
62		Azobenzene	1.330	1.270	4.5	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	0.103	0.107	-3.9	115	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.589	2.0	104	0.00
66		4-Bromophenyl-phenylether	0.200	0.193	3.5	104	0.00
67		Hexachlorobenzene	0.222	0.212	4.5	105	0.00
68		Atrazine	0.198	0.198	0.0	105	0.00
69	C	Pentachlorophenol	0.115	0.115	0.0	108	0.00
70		Phenanthrene	1.032	0.969	6.1	103	0.00
71		Anthracene	1.004	0.969	3.5	102	0.00
72		Carbazole	0.948	0.918	3.2	104	0.00
73		Di-n-butylphthalate	1.162	1.153	0.8	104	0.00
74	C	Fluoranthene	1.129	1.073	5.0	102	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76		Benzidine	0.568	0.566	0.4	106	0.00
77		Pyrene	1.115	1.058	5.1	103	0.00
78	S	Terphenyl-d14	0.794	0.759	4.4	104	0.00
79		Butylbenzylphthalate	0.500	0.508	-1.6	105	0.00
80		Benzo(a)anthracene	1.056	1.005	4.8	102	0.00
81		3,3'-Dichlorobenzidine	0.383	0.371	3.1	102	0.00
82		Chrysene	1.006	0.962	4.4	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.688	0.695	-1.0	104	0.00
84	c	Di-n-octyl phthalate	1.084	1.111	-2.5	104	0.00
85		Indeno(1,2,3-cd)pyrene	1.217	1.147	5.8	102	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Benzo(b)fluoranthene	1.094	1.052	3.8	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071715\
Data File : BG017926.D
Acq On : 17 Jul 2015 15:35
Operator : TP/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG071715

Quant Time: Jul 18 02:18:23 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 2015
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	1.093	1.060	3.0	102	0.00
89 C	Benzo(a)pyrene	1.020	0.985	3.4	102	0.00
90	Dibenzo(a,h)anthracene	1.057	1.025	3.0	104	0.00
91	Benzo(g,h,i)perylene	1.019	0.970	4.8	102	0.00

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