Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
Data File : BG016867.D
Acq On : 5 May 2015 16:11
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG050515

Quant Time: May 05 17:06:17 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 05 16:13:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	104	0.00
2		1,4-Dioxane	40.000	39.186	2.0	108	0.00
3		Pyridine	40.000	39.882	0.3	106	0.00
4		n-Nitrosodimethylamine	40.000	39.875	0.3	112	0.00
5	S	2-Fluorophenol	80.000	78.596	1.8	107	0.00
6		Aniline	40.000	39.678	0.8	108	0.00
7	S	Phenol-d6	80.000	78.414	2.0	107	0.00
8		2-Chlorophenol	40.000	38.805	3.0	107	0.00
9		Benzaldehyde	40.000	41.166	-2.9	108	0.00
10	C	Phenol	40.000	39.731	0.7	109	0.00
11		bis(2-Chloroethyl)ether	40.000	39.234	1.9	107	0.00
12		1,3-Dichlorobenzene	40.000	38.889	2.8	109	0.00
13	C	1,4-Dichlorobenzene	40.000	38.223	4.4	104	0.00
14		1,2-Dichlorobenzene	40.000	39.654	0.9	109	0.00
15		Benzyl Alcohol	40.000	39.720	0.7	109	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.402	4.0	107	0.00
17		2-Methylphenol	40.000	39.870	0.3	109	0.00
18		Hexachloroethane	40.000	38.496	3.8	106	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.573	1.1	110	0.00
20		3+4-Methylphenols	40.000	39.690	0.8	108	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	40.949	-2.4	109	0.00
23	S	Nitrobenzene-d5	80.000	82.534	-3.2	111	0.00
24		Nitrobenzene	40.000	40.382	-1.0	108	0.00
25		Isophorone	40.000	40.828	-2.1	110	0.00
26	C	2-Nitrophenol	40.000	41.774	-4.4	111	0.00
27		2,4-Dimethylphenol	40.000	41.508	-3.8	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.534	-1.3	112	0.00
29	C	2,4-Dichlorophenol	40.000	40.131	-0.3	107	0.00
30		1,2,4-Trichlorobenzene	40.000	40.708	-1.8	109	0.00
31		Naphthalene	40.000	39.791	0.5	108	0.00
32		Benzoic acid	40.000	44.288	-10.7	120	0.00
33		4-Chloroaniline	40.000	41.354	-3.4	109	0.00
34	C	Hexachlorobutadiene	40.000	40.211	-0.5	110	0.00
35		Caprolactam	40.000	41.489	-3.7	111	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.011	-2.5	110	0.00
37		2-Methylnaphthalene	40.000	40.241	-0.6	110	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.810	-2.0	108	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.711	-4.3	111	0.00
41	S	2,4,6-Tribromophenol	80.000	84.052	-5.1	112	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.977	-4.9	106	0.00
43		2,4,5-Trichlorophenol	40.000	41.714	-4.3	108	0.00
44	S	2-Fluorobiphenyl	80.000	82.665	-3.3	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
Data File : BG016867.D
Acq On : 5 May 2015 16:11
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG050515

Quant Time: May 05 17:06:17 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 05 16:13:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	41.612	-4.0	110	0.00
46		2-Chloronaphthalene	40.000	41.178	-2.9	109	0.00
47		2-Nitroaniline	40.000	43.942	-9.9	114	0.00
48		Acenaphthylene	40.000	41.082	-2.7	108	0.00
49		Dimethylphthalate	40.000	42.098	-5.2	112	0.00
50		2,6-Dinitrotoluene	40.000	43.138	-7.8	110	0.00
51	C	Acenaphthene	40.000	40.558	-1.4	108	0.00
52		3-Nitroaniline	40.000	43.111	-7.8	108	0.00
53	P	2,4-Dinitrophenol	40.000	43.313	-8.3	120	0.00
54		Dibenzofuran	40.000	41.750	-4.4	109	0.00
55	P	4-Nitrophenol	40.000	43.405	-8.5	113	0.00
56		2,4-Dinitrotoluene	40.000	45.585	-14.0	117	0.00
57		Fluorene	40.000	42.099	-5.2	111	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.703	-6.8	108	0.00
59		Diethylphthalate	40.000	41.702	-4.3	111	0.00
60		4-Chlorophenyl-phenylether	40.000	41.362	-3.4	112	0.00
61		4-Nitroaniline	40.000	43.138	-7.8	112	0.00
62		Azobenzene	40.000	41.506	-3.8	110	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.117	-7.8	115	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.987	-2.5	109	0.00
66		4-Bromophenyl-phenylether	40.000	41.627	-4.1	111	0.00
67		Hexachlorobenzene	40.000	41.353	-3.4	112	0.00
68		Atrazine	40.000	41.550	-3.9	108	0.00
69	C	Pentachlorophenol	40.000	42.755	-6.9	110	0.00
70		Phenanthrene	40.000	41.266	-3.2	109	0.00
71		Anthracene	40.000	41.398	-3.5	110	0.00
72		Carbazole	40.000	41.362	-3.4	108	0.00
73		Di-n-butylphthalate	40.000	41.197	-3.0	108	0.00
74	C	Fluoranthene	40.000	41.665	-4.2	108	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
76		Benzidine	40.000	41.152	-2.9	100	0.00
77		Pyrene	40.000	42.001	-5.0	108	0.00
78	S	Terphenyl-d14	80.000	84.301	-5.4	106	0.00
79		Butylbenzylphthalate	40.000	41.892	-4.7	106	0.00
80		Benzo(a)anthracene	40.000	41.719	-4.3	106	0.00
81		3,3'-Dichlorobenzidine	40.000	41.464	-3.7	105	0.00
82		Chrysene	40.000	41.809	-4.5	106	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	42.271	-5.7	106	0.00
84	c	Di-n-octyl phthalate	40.000	42.391	-6.0	106	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.506	-3.8	109	0.02
86	I	Perylene-d12	20.000	20.000	0.0	97	0.00
87		Benzo(b)fluoranthene	40.000	41.446	-3.6	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
Data File : BG016867.D
Acq On : 5 May 2015 16:11
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG050515

Quant Time: May 05 17:06:17 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 05 16:13:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	40.291	-0.7	105	0.00
89 C	Benzo(a)pyrene	40.000	40.576	-1.4	105	0.00
90	Dibenzo(a,h)anthracene	40.000	40.496	-1.2	107	0.01
91	Benzo(g,h,i)perylene	40.000	40.245	-0.6	109	0.03

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