Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG010515\
Data File : BG015845.D
Acq On : 5 Jan 2015 16:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG010515

Quant Time: Jan 06 06:06:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG010515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 06 03:51:36 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	102	0.00
2		1,4-Dioxane	40.000	40.559	-1.4	107	0.00
3		Pyridine	40.000	38.386	4.0	99	0.00
4		n-Nitrosodimethylamine	40.000	40.169	-0.4	105	0.00
5	S	2-Fluorophenol	40.000	39.498	1.3	103	0.00
6		Aniline	40.000	40.116	-0.3	104	0.00
7	S	Phenol-d6	40.000	40.371	-0.9	106	0.00
8		2-Chlorophenol	40.000	38.460	3.8	102	0.00
9		Benzaldehyde	40.000	39.008	2.5	102	0.00
10	C	Phenol	40.000	39.909	0.2	105	0.00
11		bis(2-Chloroethyl)ether	40.000	38.957	2.6	105	0.00
12		1,3-Dichlorobenzene	40.000	38.932	2.7	105	0.00
13	C	1,4-Dichlorobenzene	40.000	38.895	2.8	106	0.00
14		1,2-Dichlorobenzene	40.000	37.808	5.5	102	0.00
15		Benzyl Alcohol	40.000	39.657	0.9	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.566	1.1	105	0.00
17		2-Methylphenol	40.000	39.166	2.1	101	0.00
18		Hexachloroethane	40.000	38.189	4.5	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.547	3.6	104	0.00
20		3+4-Methylphenols	40.000	38.637	3.4	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	38.723	3.2	101	0.00
23	S	Nitrobenzene-d5	40.000	40.275	-0.7	103	0.00
24		Nitrobenzene	40.000	40.577	-1.4	102	0.00
25		Isophorone	40.000	40.968	-2.4	104	0.00
26	C	2-Nitrophenol	40.000	40.940	-2.3	103	0.00
27		2,4-Dimethylphenol	40.000	38.415	4.0	101	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.134	2.2	104	0.00
29	C	2,4-Dichlorophenol	40.000	40.785	-2.0	100	0.00
30		1,2,4-Trichlorobenzene	40.000	39.790	0.5	104	0.00
31		Naphthalene	40.000	40.313	-0.8	104	0.00
32		Benzoic acid	40.000	42.428	-6.1	110	0.00
33		4-Chloroaniline	40.000	39.391	1.5	102	0.00
34	C	Hexachlorobutadiene	40.000	40.334	-0.8	108	0.00
35		Caprolactam	40.000	40.246	-0.6	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.903	-4.8	108	0.00
37		2-Methylnaphthalene	40.000	39.958	0.1	102	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.735	0.7	103	0.00
40	P	Hexachlorocyclopentadiene	40.000	43.665	-9.2	104	0.00
41	S	2,4,6-Tribromophenol	40.000	40.459	-1.1	99	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.609	-4.0	104	0.00
43		2,4,5-Trichlorophenol	40.000	41.387	-3.5	105	0.00
44	S	2-Fluorobiphenyl	40.000	40.619	-1.5	106	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG010515\
Data File : BG015845.D
Acq On : 5 Jan 2015 16:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG010515

Quant Time: Jan 06 06:06:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG010515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 06 03:51:36 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	40.064	-0.2	102	0.00
46		2-Chloronaphthalene	40.000	39.681	0.8	102	0.00
47		2-Nitroaniline	40.000	41.113	-2.8	98	0.00
48		Acenaphthylene	40.000	39.822	0.4	103	0.00
49		Dimethylphthalate	40.000	39.490	1.3	102	0.00
50		2,6-Dinitrotoluene	40.000	40.580	-1.4	99	0.00
51	C	Acenaphthene	40.000	39.852	0.4	102	0.00
52		3-Nitroaniline	40.000	40.772	-1.9	102	0.00
53	P	2,4-Dinitrophenol	40.000	42.334	-5.8	104	0.00
54		Dibenzofuran	40.000	39.455	1.4	100	0.00
55	P	4-Nitrophenol	40.000	40.903	-2.3	105	0.00
56		2,4-Dinitrotoluene	40.000	40.788	-2.0	102	0.00
57		Fluorene	40.000	39.879	0.3	100	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.704	0.7	103	0.00
59		Diethylphthalate	40.000	38.908	2.7	100	0.00
60		4-Chlorophenyl-phenylether	40.000	39.766	0.6	102	0.00
61		4-Nitroaniline	40.000	39.258	1.9	97	0.00
62		Azobenzene	40.000	39.703	0.7	102	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.821	-9.6	104	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.963	-2.4	100	0.00
66		4-Bromophenyl-phenylether	40.000	41.592	-4.0	103	0.00
67		Hexachlorobenzene	40.000	40.918	-2.3	103	0.00
68		Atrazine	40.000	40.021	-0.1	98	0.00
69	C	Pentachlorophenol	40.000	40.329	-0.8	96	0.00
70		Phenanthrene	40.000	40.142	-0.4	98	0.00
71		Anthracene	40.000	40.680	-1.7	98	0.00
72		Carbazole	40.000	40.127	-0.3	97	0.00
73		Di-n-butylphthalate	40.000	40.593	-1.5	97	0.00
74	C	Fluoranthene	40.000	40.495	-1.2	98	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	98	0.00
76		Benzidine	40.000	41.133	-2.8	96	0.00
77		Pyrene	40.000	39.475	1.3	94	0.00
78	S	Terphenyl-d14	40.000	40.336	-0.8	98	0.00
79		Butylbenzylphthalate	40.000	39.719	0.7	99	0.00
80		Benzo(a)anthracene	40.000	41.352	-3.4	101	0.00
81		3,3'-Dichlorobenzidine	40.000	42.279	-5.7	105	0.00
82		Chrysene	40.000	40.716	-1.8	99	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.637	-1.6	99	0.00
84	c	Di-n-octyl phthalate	40.000	41.759	-4.4	103	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.342	-3.4	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	102	0.00
87		Benzo(b)fluoranthene	40.000	40.060	-0.2	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG010515\
Data File : BG015845.D
Acq On : 5 Jan 2015 16:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG010515

Quant Time: Jan 06 06:06:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG010515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 06 03:51:36 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	41.557	-3.9	108	0.00
89 C	Benzo(a)pyrene	40.000	41.145	-2.9	103	0.00
90	Dibenzo(a,h)anthracene	40.000	40.881	-2.2	102	0.00
91	Benzo(g,h,i)perylene	40.000	40.231	-0.6	102	-0.01

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