Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061815\
Data File : BG017724.D
Acq On : 18 Jun 2015 14:57
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061815

Quant Time: Jun 19 06:40:41 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 06:05:07 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	101	0.00
2		1,4-Dioxane	40.000	39.195	2.0	101	0.00
3		Pyridine	40.000	41.362	-3.4	103	0.00
4		n-Nitrosodimethylamine	40.000	42.264	-5.7	103	0.00
5	S	2-Fluorophenol	80.000	80.975	-1.2	103	0.00
6		Aniline	40.000	40.920	-2.3	103	0.00
7	S	Phenol-d6	80.000	82.436	-3.0	103	0.00
8		2-Chlorophenol	40.000	40.744	-1.9	101	0.00
9		Benzaldehyde	40.000	42.018	-5.0	106	0.00
10	C	Phenol	40.000	40.983	-2.5	102	0.00
11		bis(2-Chloroethyl)ether	40.000	40.384	-1.0	101	0.00
12		1,3-Dichlorobenzene	40.000	40.194	-0.5	103	0.00
13	C	1,4-Dichlorobenzene	40.000	40.173	-0.4	103	0.00
14		1,2-Dichlorobenzene	40.000	40.510	-1.3	104	0.00
15		Benzyl Alcohol	40.000	41.585	-4.0	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.202	-3.0	106	0.00
17		2-Methylphenol	40.000	40.692	-1.7	103	0.00
18		Hexachloroethane	40.000	40.852	-2.1	104	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.779	0.6	103	0.00
20		3+4-Methylphenols	40.000	41.631	-4.1	103	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	39.927	0.2	101	0.00
23	S	Nitrobenzene-d5	80.000	83.430	-4.3	105	0.00
24		Nitrobenzene	40.000	41.467	-3.7	104	0.00
25		Isophorone	40.000	40.957	-2.4	103	0.00
26	C	2-Nitrophenol	40.000	42.071	-5.2	105	0.00
27		2,4-Dimethylphenol	40.000	40.633	-1.6	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.165	-0.4	102	0.00
29	C	2,4-Dichlorophenol	40.000	40.504	-1.3	101	0.00
30		1,2,4-Trichlorobenzene	40.000	39.908	0.2	102	0.00
31		Naphthalene	40.000	40.464	-1.2	102	0.00
32		Benzoic acid	40.000	43.745	-9.4	120	0.00
33		4-Chloroaniline	40.000	40.652	-1.6	101	0.00
34	C	Hexachlorobutadiene	40.000	40.762	-1.9	103	0.00
35		Caprolactam	40.000	39.948	0.1	102	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.147	-0.4	101	0.00
37		2-Methylnaphthalene	40.000	40.372	-0.9	101	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.424	-1.1	104	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.519	-3.8	102	0.00
41	S	2,4,6-Tribromophenol	80.000	84.896	-6.1	104	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.189	-3.0	101	0.00
43		2,4,5-Trichlorophenol	40.000	42.428	-6.1	106	0.00
44	S	2-Fluorobiphenyl	80.000	80.304	-0.4	100	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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	40.197	-0.5	101	0.00
46		2-Chloronaphthalene	40.000	40.686	-1.7	102	0.00
47		2-Nitroaniline	40.000	43.176	-7.9	106	0.00
48		Acenaphthylene	40.000	40.430	-1.1	101	0.00
49		Dimethylphthalate	40.000	40.181	-0.5	100	0.00
50		2,6-Dinitrotoluene	40.000	42.778	-6.9	103	0.00
51	C	Acenaphthene	40.000	39.934	0.2	100	0.00
52		3-Nitroaniline	40.000	42.553	-6.4	101	0.00
53	P	2,4-Dinitrophenol	40.000	40.815	-2.0	108	0.00
54		Dibenzofuran	40.000	40.588	-1.5	101	0.00
55	P	4-Nitrophenol	40.000	44.337	-10.8	104	0.00
56		2,4-Dinitrotoluene	40.000	40.797	-2.0	104	0.00
57		Fluorene	40.000	40.490	-1.2	101	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.387	-6.0	103	0.00
59		Diethylphthalate	40.000	39.669	0.8	100	0.00
60		4-Chlorophenyl-phenylether	40.000	40.181	-0.5	101	0.00
61		4-Nitroaniline	40.000	43.706	-9.3	103	0.00
62		Azobenzene	40.000	40.045	-0.1	100	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.306	-0.8	103	0.00
65	c	n-Nitrosodiphenylamine	40.000	41.313	-3.3	100	0.00
66		4-Bromophenyl-phenylether	40.000	41.385	-3.5	100	0.00
67		Hexachlorobenzene	40.000	41.108	-2.8	100	0.00
68		Atrazine	40.000	41.876	-4.7	101	0.00
69	C	Pentachlorophenol	40.000	44.999	-12.5	104	0.00
70		Phenanthrene	40.000	41.015	-2.5	100	0.00
71		Anthracene	40.000	41.045	-2.6	99	0.00
72		Carbazole	40.000	40.873	-2.2	100	0.00
73		Di-n-butylphthalate	40.000	40.896	-2.2	100	0.00
74	C	Fluoranthene	40.000	41.309	-3.3	101	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	101	0.00
76		Benzidine	40.000	40.565	-1.4	97	0.00
77		Pyrene	40.000	40.457	-1.1	101	0.00
78	S	Terphenyl-d14	80.000	80.635	-0.8	101	0.00
79		Butylbenzylphthalate	40.000	39.950	0.1	99	0.00
80		Benzo(a)anthracene	40.000	40.440	-1.1	101	0.00
81		3,3'-Dichlorobenzidine	40.000	39.814	0.5	97	0.00
82		Chrysene	40.000	40.271	-0.7	101	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.099	-0.2	100	0.00
84	c	Di-n-octyl phthalate	40.000	40.322	-0.8	100	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	41.182	-3.0	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	101	0.00
87		Benzo(b)fluoranthene	40.000	40.125	-0.3	97	0.00

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88	Benzo(k)fluoranthene	40.000	41.893	-4.7	103	0.00
89 C	Benzo(a)pyrene	40.000	41.000	-2.5	100	0.00
90	Dibenzo(a,h)anthracene	40.000	41.099	-2.7	101	0.00
91	Benzo(g,h,i)perylene	40.000	40.967	-2.4	101	-0.01

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