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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
2		1,4-Dioxane	0.533	0.541	-1.5	107	0.00
3		Pyridine	1.462	1.403	4.0	99	0.00
4		n-Nitrosodimethylamine	0.628	0.631	-0.5	105	0.00
5	S	2-Fluorophenol	2.326	2.296	1.3	103	0.00
6		Aniline	2.190	2.196	-0.3	104	0.00
7	S	Phenol-d6	3.239	3.269	-0.9	106	0.00
8		2-Chlorophenol	1.199	1.152	3.9	102	0.00
9		Benzaldehyde	1.091	1.064	2.5	102	0.00
10	C	Phenol	1.763	1.759	0.2	105	0.00
11		bis(2-Chloroethyl)ether	1.299	1.266	2.5	105	0.00
12		1,3-Dichlorobenzene	1.437	1.399	2.6	105	0.00
13	C	1,4-Dichlorobenzene	1.466	1.426	2.7	106	0.00
14		1,2-Dichlorobenzene	1.392	1.315	5.5	102	0.00
15		Benzyl Alcohol	1.582	1.568	0.9	102	0.00
16		2,2'-oxybis(1-Chloropropane	0.830	0.821	1.1	105	0.00
17		2-Methylphenol	1.203	1.178	2.1	101	0.00
18		Hexachloroethane	0.677	0.646	4.6	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.264	1.218	3.6	104	0.00
20		3+4-Methylphenols	1.604	1.549	3.4	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.586	0.568	3.1	101	0.00
23	S	Nitrobenzene-d5	1.019	1.026	-0.7	103	0.00
24		Nitrobenzene	0.513	0.521	-1.6	102	0.00
25		Isophorone	0.904	0.926	-2.4	104	0.00
26	C	2-Nitrophenol	0.197	0.201	-2.0	103	0.00
27		2,4-Dimethylphenol	0.329	0.316	4.0	101	0.00
28		bis(2-Chloroethoxy)methane	0.446	0.436	2.2	104	0.00
29	C	2,4-Dichlorophenol	0.333	0.339	-1.8	100	0.00
30		1,2,4-Trichlorobenzene	0.436	0.433	0.7	104	0.00
31		Naphthalene	1.059	1.068	-0.8	104	0.00
32		Benzoic acid	0.152	0.161	-5.9	110	0.00
33		4-Chloroaniline	0.431	0.425	1.4	102	0.00
34	C	Hexachlorobutadiene	0.405	0.408	-0.7	108	0.00
35		Caprolactam	0.146	0.147	-0.7	103	0.00
36	C	4-Chloro-3-methylphenol	0.433	0.454	-4.8	108	0.00
37		2-Methylnaphthalene	0.793	0.792	0.1	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	0.744	0.739	0.7	103	0.00
40	P	Hexachlorocyclopentadiene	0.495	0.540	-9.1	104	0.00
41	S	2,4,6-Tribromophenol	0.760	0.769	-1.2	99	0.00
42	C	2,4,6-Trichlorophenol	0.457	0.475	-3.9	104	0.00
43		2,4,5-Trichlorophenol	0.500	0.517	-3.4	105	0.00
44	S	2-Fluorobiphenyl	2.800	2.844	-1.6	106	0.00

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG010515

Quant Time: Jan 06 06:06:27 2015
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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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.428	1.430	-0.1	102	0.00
46		2-Chloronaphthalene	1.138	1.129	0.8	102	0.00
47		2-Nitroaniline	0.372	0.382	-2.7	98	0.00
48		Acenaphthylene	1.946	1.938	0.4	103	0.00
49		Dimethylphthalate	1.476	1.457	1.3	102	0.00
50		2,6-Dinitrotoluene	0.318	0.323	-1.6	99	0.00
51	C	Acenaphthene	1.143	1.139	0.3	102	0.00
52		3-Nitroaniline	0.303	0.309	-2.0	102	0.00
53	P	2,4-Dinitrophenol	0.177	0.187	-5.6	104	0.00
54		Dibenzofuran	1.854	1.829	1.3	100	0.00
55	P	4-Nitrophenol	0.249	0.255	-2.4	105	0.00
56		2,4-Dinitrotoluene	0.458	0.467	-2.0	102	0.00
57		Fluorene	1.511	1.507	0.3	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.564	0.560	0.7	103	0.00
59		Diethylphthalate	1.630	1.585	2.8	100	0.00
60		4-Chlorophenyl-phenylether	0.987	0.981	0.6	102	0.00
61		4-Nitroaniline	0.356	0.349	2.0	97	0.00
62		Azobenzene	1.756	1.743	0.7	102	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.125	0.150	-20.0	104	0.00
65	c	n-Nitrosodiphenylamine	0.575	0.589	-2.4	100	0.00
66		4-Bromophenyl-phenylether	0.274	0.285	-4.0	103	0.00
67		Hexachlorobenzene	0.311	0.318	-2.3	103	0.00
68		Atrazine	0.257	0.257	0.0	98	0.00
69	C	Pentachlorophenol	0.169	0.185	-9.5	96	0.00
70		Phenanthrene	1.064	1.068	-0.4	98	0.00
71		Anthracene	1.057	1.075	-1.7	98	0.00
72		Carbazole	1.014	1.017	-0.3	97	0.00
73		Di-n-butylphthalate	1.194	1.212	-1.5	97	0.00
74	C	Fluoranthene	1.610	1.630	-1.2	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.463	0.476	-2.8	96	0.00
77		Pyrene	1.069	1.055	1.3	94	0.00
78	S	Terphenyl-d14	1.790	1.805	-0.8	98	0.00
79		Butylbenzylphthalate	0.401	0.398	0.7	99	0.00
80		Benzo(a)anthracene	1.188	1.228	-3.4	101	0.00
81		3,3'-Dichlorobenzidine	0.436	0.461	-5.7	105	0.00
82		Chrysene	1.092	1.111	-1.7	99	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.592	-1.5	99	0.00
84	c	Di-n-octyl phthalate	0.949	0.990	-4.3	103	0.00
85		Indeno(1,2,3-cd)pyrene	1.290	1.333	-3.3	102	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.249	1.250	-0.1	100	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.202	1.249	-3.9	108	0.00
89 C	Benzo(a)pyrene	1.125	1.157	-2.8	103	0.00
90	Dibenzo(a,h)anthracene	1.105	1.130	-2.3	102	0.00
91	Benzo(g,h,i)perylene	1.099	1.105	-0.5	102	-0.01

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