Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG090717\
Data File : BG028599.D
Acq On : 7 Sep 2017 18:34
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG090717

Quant Time: Sep 07 19:22:30 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG090717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 07 18:39:50 2017
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	104	0.00
2		1,4-Dioxane	0.486	0.478	1.6	104	0.00
3		Pyridine	1.509	1.441	4.5	101	0.00
4		n-Nitrosodimethylamine	0.719	0.673	6.4	99	0.00
5	S	2-Fluorophenol	1.188	1.141	4.0	101	0.00
6		Aniline	2.206	2.145	2.8	99	0.00
7	S	Phenol-d6	1.723	1.686	2.1	100	0.00
8		2-Chlorophenol	1.341	1.289	3.9	101	0.00
9		Benzaldehyde	1.017	0.974	4.2	101	0.00
10	C	Phenol	1.866	1.826	2.1	101	0.00
11		bis(2-Chloroethyl)ether	1.372	1.302	5.1	100	0.00
12		1,3-Dichlorobenzene	1.570	1.481	5.7	101	0.00
13	C	1,4-Dichlorobenzene	1.548	1.511	2.4	101	0.00
14		1,2-Dichlorobenzene	1.521	1.455	4.3	100	0.00
15		Benzyl Alcohol	1.366	1.303	4.6	97	0.00
16		2,2'-oxybis(1-Chloropropane	2.829	2.645	6.5	98	0.00
17		2-Methylphenol	1.197	1.148	4.1	101	0.00
18		Hexachloroethane	0.608	0.556	8.6	96	0.00
19	P	n-Nitroso-di-n-propylamine	1.362	1.319	3.2	100	0.00
20		3+4-Methylphenols	1.657	1.632	1.5	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.556	0.555	0.2	100	0.00
23	S	Nitrobenzene-d5	0.436	0.432	0.9	99	0.00
24		Nitrobenzene	0.460	0.464	-0.9	100	0.00
25		Isophorone	0.802	0.797	0.6	98	0.00
26	C	2-Nitrophenol	0.191	0.192	-0.5	103	0.00
27		2,4-Dimethylphenol	0.345	0.339	1.7	97	0.00
28		bis(2-Chloroethoxy)methane	0.479	0.474	1.0	101	0.00
29	C	2,4-Dichlorophenol	0.338	0.342	-1.2	101	0.00
30		1,2,4-Trichlorobenzene	0.416	0.413	0.7	99	0.00
31		Naphthalene	1.086	1.068	1.7	99	0.00
32		Benzoic acid	0.217	0.230	-6.0	103	0.00
33		4-Chloroaniline	0.458	0.456	0.4	98	0.00
34	C	Hexachlorobutadiene	0.301	0.299	0.7	99	0.00
35		Caprolactam	0.128	0.129	-0.8	101	0.00
36	C	4-Chloro-3-methylphenol	0.431	0.433	-0.5	98	0.00
37		2-Methylnaphthalene	0.800	0.798	0.3	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	0.777	0.783	-0.8	100	0.00
40	P	Hexachlorocyclopentadiene	0.271	0.283	-4.4	100	0.00
41	S	2,4,6-Tribromophenol	0.304	0.311	-2.3	101	0.00
42	C	2,4,6-Trichlorophenol	0.496	0.499	-0.6	98	0.00
43		2,4,5-Trichlorophenol	0.496	0.500	-0.8	99	0.00
44	S	2-Fluorobiphenyl	1.529	1.523	0.4	99	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.679	1.666	0.8	100	0.00
46		2-Chloronaphthalene	1.275	1.292	-1.3	101	0.00
47		2-Nitroaniline	0.520	0.512	1.5	99	0.00
48		Acenaphthylene	1.956	1.939	0.9	100	0.00
49		Dimethylphthalate	1.648	1.620	1.7	99	0.00
50		2,6-Dinitrotoluene	0.361	0.365	-1.1	98	0.00
51	C	Acenaphthene	1.198	1.211	-1.1	100	0.00
52		3-Nitroaniline	0.365	0.363	0.5	98	0.00
53	P	2,4-Dinitrophenol	0.195	0.207	-6.2	98	0.00
54		Dibenzofuran	1.852	1.872	-1.1	101	0.00
55	P	4-Nitrophenol	0.260	0.264	-1.5	103	0.00
56		2,4-Dinitrotoluene	0.520	0.531	-2.1	104	0.00
57		Fluorene	1.693	1.703	-0.6	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.443	0.444	-0.2	101	0.00
59		Diethylphthalate	1.727	1.727	0.0	102	0.00
60		4-Chlorophenyl-phenylether	0.956	0.960	-0.4	101	0.00
61		4-Nitroaniline	0.395	0.399	-1.0	100	0.00
62		Azobenzene	1.542	1.509	2.1	99	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.142	0.146	-2.8	105	0.00
65	c	n-Nitrosodiphenylamine	0.591	0.592	-0.2	99	0.00
66		4-Bromophenyl-phenylether	0.260	0.263	-1.2	100	0.00
67		Hexachlorobenzene	0.286	0.289	-1.0	101	0.00
68		Atrazine	0.213	0.210	1.4	101	0.00
69	C	Pentachlorophenol	0.133	0.134	-0.8	103	0.00
70		Phenanthrene	1.064	1.081	-1.6	102	0.00
71		Anthracene	1.082	1.092	-0.9	102	0.00
72		Carbazole	0.968	0.965	0.3	100	0.00
73		Di-n-butylphthalate	1.175	1.153	1.9	99	0.00
74	C	Fluoranthene	1.358	1.364	-0.4	101	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
76		Benzidine	0.406	0.393	3.2	95	0.00
77		Pyrene	1.186	1.202	-1.3	100	0.00
78	S	Terphenyl-d14	0.928	0.945	-1.8	101	0.00
79		Butylbenzylphthalate	0.459	0.461	-0.4	102	0.00
80		Benzo(a)anthracene	1.197	1.200	-0.3	101	0.00
81		3,3'-Dichlorobenzidine	0.460	0.463	-0.7	100	0.00
82		Chrysene	1.129	1.118	1.0	99	0.00
83		Bis(2-ethylhexyl)phthalate	0.661	0.647	2.1	99	0.00
84	c	Di-n-octyl phthalate	1.133	1.122	1.0	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.338	1.347	-0.7	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.253	1.241	1.0	101	0.00

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88	Benzo(k)fluoranthene	1.210	1.220	-0.8	99	0.00
89 C	Benzo(a)pyrene	1.175	1.177	-0.2	101	0.00
90	Dibenzo(a,h)anthracene	1.219	1.230	-0.9	100	-0.01
91	Benzo(g,h,i)perylene	1.196	1.200	-0.3	101	-0.02

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