Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091415\
Data File : BG018661.D
Acq On : 14 Sep 2015 14:32
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 15 01:19:34 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG091115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 11 10:44:16 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	111	0.00
2		1,4-Dioxane	0.484	0.485	-0.2	127	0.00
3		Pyridine	1.500	1.455	3.0	116	-0.01
4		n-Nitrosodimethylamine	0.522	0.506	3.1	114	0.00
5	S	2-Fluorophenol	1.158	1.175	-1.5	121	0.00
6		Aniline	2.288	2.281	0.3	117	0.00
7	S	Phenol-d6	1.660	1.689	-1.7	119	0.00
8		2-Chlorophenol	1.337	1.362	-1.9	122	0.00
9		Benzaldehyde	0.894	0.890	0.4	113	0.00
10	C	Phenol	1.754	1.765	-0.6	119	0.00
11		bis(2-Chloroethyl)ether	1.359	1.345	1.0	119	0.00
12		1,3-Dichlorobenzene	1.555	1.539	1.0	120	0.00
13	C	1,4-Dichlorobenzene	1.560	1.565	-0.3	119	0.00
14		1,2-Dichlorobenzene	1.489	1.521	-2.1	121	0.00
15		Benzyl Alcohol	1.197	1.209	-1.0	117	0.00
16		2,2'-oxybis(1-Chloropropane	1.543	1.460	5.4	115	0.00
17		2-Methylphenol	1.219	1.230	-0.9	120	0.00
18		Hexachloroethane	0.558	0.545	2.3	117	0.00
19	P	n-Nitroso-di-n-propylamine	1.169	1.147	1.9	118	0.00
20		3+4-Methylphenols	1.711	1.725	-0.8	121	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
22		Acetophenone	0.507	0.492	3.0	117	0.00
23	S	Nitrobenzene-d5	0.357	0.348	2.5	118	0.00
24		Nitrobenzene	0.380	0.372	2.1	118	0.00
25		Isophorone	0.769	0.724	5.9	115	0.00
26	C	2-Nitrophenol	0.177	0.181	-2.3	117	0.00
27		2,4-Dimethylphenol	0.322	0.313	2.8	118	0.00
28		bis(2-Chloroethoxy)methane	0.441	0.428	2.9	117	0.00
29	C	2,4-Dichlorophenol	0.325	0.331	-1.8	120	0.00
30		1,2,4-Trichlorobenzene	0.360	0.361	-0.3	121	0.00
31		Naphthalene	1.071	1.059	1.1	120	0.00
32		Benzoic acid	0.237	0.243	-2.5	119	0.00
33		4-Chloroaniline	0.485	0.480	1.0	121	0.00
34	C	Hexachlorobutadiene	0.212	0.213	-0.5	125	0.00
35		Caprolactam	0.140	0.136	2.9	113	0.04
36	C	4-Chloro-3-methylphenol	0.363	0.361	0.6	121	0.00
37		2-Methylnaphthalene	0.784	0.773	1.4	120	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.634	0.641	-1.1	122	0.00
40	P	Hexachlorocyclopentadiene	0.319	0.338	-6.0	122	0.00
41	S	2,4,6-Tribromophenol	0.269	0.281	-4.5	125	0.00
42	C	2,4,6-Trichlorophenol	0.439	0.468	-6.6	123	0.00
43		2,4,5-Trichlorophenol	0.482	0.479	0.6	119	0.00
44	S	2-Fluorobiphenyl	1.441	1.402	2.7	117	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.594	-0.3	121	0.00
46		2-Chloronaphthalene	1.225	1.232	-0.6	122	0.00
47		2-Nitroaniline	0.369	0.365	1.1	114	0.00
48		Acenaphthylene	2.166	2.161	0.2	119	0.00
49		Dimethylphthalate	1.656	1.631	1.5	120	0.00
50		2,6-Dinitrotoluene	0.360	0.371	-3.1	120	0.00
51	C	Acenaphthene	1.260	1.315	-4.4	126	0.00
52		3-Nitroaniline	0.430	0.444	-3.3	120	0.00
53	P	2,4-Dinitrophenol	0.152	0.156	-2.6	115	0.00
54		Dibenzofuran	1.958	1.941	0.9	120	0.00
55	P	4-Nitrophenol	0.332	0.343	-3.3	116	0.00
56		2,4-Dinitrotoluene	0.510	0.543	-6.5	121	0.00
57		Fluorene	1.686	1.644	2.5	118	0.00
58		2,3,4,6-Tetrachlorophenol	0.452	0.452	0.0	123	0.00
59		Diethylphthalate	1.727	1.677	2.9	117	0.00
60		4-Chlorophenyl-phenylether	0.810	0.816	-0.7	121	0.00
61		4-Nitroaniline	0.464	0.471	-1.5	118	0.01
62		Azobenzene	1.527	1.482	2.9	115	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
64		4,6-Dinitro-2-methylphenol	0.100	0.103	-3.0	123	0.00
65	c	n-Nitrosodiphenylamine	0.579	0.556	4.0	118	0.00
66		4-Bromophenyl-phenylether	0.203	0.204	-0.5	122	0.00
67		Hexachlorobenzene	0.221	0.227	-2.7	125	0.00
68		Atrazine	0.230	0.219	4.8	114	0.00
69	C	Pentachlorophenol	0.136	0.146	-7.4	122	0.00
70		Phenanthrene	1.052	1.013	3.7	117	0.00
71		Anthracene	1.065	1.027	3.6	118	0.00
72		Carbazole	1.031	1.000	3.0	116	0.00
73		Di-n-butylphthalate	1.216	1.201	1.2	116	0.00
74	C	Fluoranthene	1.303	1.262	3.1	117	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
76		Benzidine	0.601	0.614	-2.2	118	0.00
77		Pyrene	1.229	1.212	1.4	117	0.00
78	S	Terphenyl-d14	0.762	0.735	3.5	115	0.00
79		Butylbenzylphthalate	0.500	0.512	-2.4	118	0.00
80		Benzo(a)anthracene	1.151	1.139	1.0	117	0.00
81		3,3'-Dichlorobenzidine	0.437	0.455	-4.1	121	0.00
82		Chrysene	1.084	1.056	2.6	116	0.00
83		Bis(2-ethylhexyl)phthalate	0.721	0.735	-1.9	118	0.00
84	c	Di-n-octyl phthalate	1.219	1.257	-3.1	120	0.00
85		Indeno(1,2,3-cd)pyrene	1.377	1.406	-2.1	120	0.01
86	I	Perylene-d12	1.000	1.000	0.0	114	0.00
87		Benzo(b)fluoranthene	1.160	1.154	0.5	120	0.00

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88	Benzo(k)fluoranthene	1.162	1.130	2.8	121	0.01
89 C	Benzo(a)pyrene	1.104	1.104	0.0	121	0.00
90	Dibenzo(a,h)anthracene	1.087	1.100	-1.2	122	0.00
91	Benzo(g,h,i)perylene	1.107	1.099	0.7	121	0.01

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