Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022716\
Data File : BG021100.D
Acq On : 27 Feb 2016 5:19
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 29 00:11:00 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 26 16:39:44 2016
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	100	0.00
2		1,4-Dioxane	0.529	0.484	8.5	89	0.00
3		Pyridine	1.547	1.456	5.9	90	-0.01
4		n-Nitrosodimethylamine	0.784	0.705	10.1	84	-0.01
5	S	2-Fluorophenol	1.142	1.115	2.4	93	0.00
6		Aniline	2.103	2.229	-6.0	98	0.00
7	S	Phenol-d6	1.700	1.718	-1.1	95	-0.01
8		2-Chlorophenol	1.425	1.426	-0.1	95	-0.01
9		Benzaldehyde	0.921	0.942	-2.3	99	0.00
10	C	Phenol	1.741	1.796	-3.2	97	0.00
11		bis(2-Chloroethyl)ether	1.335	1.331	0.3	96	-0.01
12		1,3-Dichlorobenzene	1.516	1.520	-0.3	95	0.00
13	C	1,4-Dichlorobenzene	1.620	1.582	2.3	97	0.00
14		1,2-Dichlorobenzene	1.534	1.493	2.7	93	-0.01
15		Benzyl Alcohol	1.485	1.510	-1.7	95	-0.01
16		2,2'-oxybis(1-Chloropropane	1.530	1.710	-11.8	109	0.00
17		2-Methylphenol	1.229	1.245	-1.3	94	0.00
18		Hexachloroethane	0.613	0.604	1.5	93	-0.01
19	P	n-Nitroso-di-n-propylamine	1.311	1.364	-4.0	95	-0.01
20		3+4-Methylphenols	1.704	1.787	-4.9	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	96	-0.01
22		Acetophenone	0.552	0.555	-0.5	94	0.00
23	S	Nitrobenzene-d5	0.420	0.430	-2.4	96	-0.01
24		Nitrobenzene	0.434	0.460	-6.0	98	0.00
25		Isophorone	0.787	0.830	-5.5	98	0.00
26	C	2-Nitrophenol	0.183	0.183	0.0	93	0.00
27		2,4-Dimethylphenol	0.323	0.324	-0.3	94	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.404	-5.5	100	0.00
29	C	2,4-Dichlorophenol	0.317	0.313	1.3	95	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.341	5.0	89	0.00
31		Naphthalene	1.027	1.012	1.5	93	0.00
32		Benzoic acid	0.297	0.292	1.7	89	-0.03
33		4-Chloroaniline	0.429	0.437	-1.9	96	0.00
34	C	Hexachlorobutadiene	0.240	0.231	3.7	87	-0.01
35		Caprolactam	0.107	0.109	-1.9	94	-0.02
36	C	4-Chloro-3-methylphenol	0.388	0.403	-3.9	96	0.00
37		2-Methylnaphthalene	0.721	0.711	1.4	93	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	94	0.00
39		1,2,4,5-Tetrachlorobenzene	0.739	0.729	1.4	90	-0.01
40	P	Hexachlorocyclopentadiene	0.486	0.469	3.5	87	-0.01
41	S	2,4,6-Tribromophenol	0.262	0.252	3.8	86	0.00
42	C	2,4,6-Trichlorophenol	0.463	0.493	-6.5	96	0.06
43		2,4,5-Trichlorophenol	0.476	0.493	-3.6	91	0.00
44	S	2-Fluorobiphenyl	1.563	1.538	1.6	92	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.660	1.4	93	0.00
46		2-Chloronaphthalene	1.278	1.272	0.5	93	-0.01
47		2-Nitroaniline	0.414	0.464	-12.1	101	0.00
48		Acenaphthylene	2.022	2.056	-1.7	96	-0.01
49		Dimethylphthalate	1.724	1.697	1.6	93	0.00
50		2,6-Dinitrotoluene	0.355	0.373	-5.1	97	0.00
51	C	Acenaphthene	1.379	1.402	-1.7	93	0.00
52		3-Nitroaniline	0.345	0.363	-5.2	96	0.00
53	P	2,4-Dinitrophenol	0.247	0.256	-3.6	92	0.00
54		Dibenzofuran	1.747	1.765	-1.0	94	0.00
55	P	4-Nitrophenol	0.306	0.329	-7.5	101	0.00
56		2,4-Dinitrotoluene	0.509	0.527	-3.5	93	0.00
57		Fluorene	1.700	1.677	1.4	92	0.00
58		2,3,4,6-Tetrachlorophenol	0.414	0.419	-1.2	92	-0.01
59		Diethylphthalate	1.817	1.822	-0.3	94	-0.01
60		4-Chlorophenyl-phenylether	0.935	0.908	2.9	91	0.00
61		4-Nitroaniline	0.393	0.427	-8.7	102	0.00
62		Azobenzene	1.574	1.726	-9.7	103	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.147	0.143	2.7	91	0.00
65	c	n-Nitrosodiphenylamine	0.590	0.590	0.0	96	0.00
66		4-Bromophenyl-phenylether	0.235	0.223	5.1	89	0.00
67		Hexachlorobenzene	0.248	0.233	6.0	90	0.00
68		Atrazine	0.219	0.224	-2.3	95	0.00
69	C	Pentachlorophenol	0.168	0.155	7.7	86	0.00
70		Phenanthrene	1.076	1.068	0.7	95	0.00
71		Anthracene	1.084	1.096	-1.1	96	-0.01
72		Carbazole	0.936	0.953	-1.8	97	0.00
73		Di-n-butylphthalate	1.233	1.294	-4.9	98	-0.01
74	C	Fluoranthene	1.350	1.337	1.0	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	92	-0.01
76		Benzidine	0.526	0.560	-6.5	92	0.00
77		Pyrene	1.107	1.183	-6.9	96	0.00
78	S	Terphenyl-d14	0.856	0.881	-2.9	92	0.00
79		Butylbenzylphthalate	0.462	0.499	-8.0	94	0.00
80		Benzo(a)anthracene	1.151	1.166	-1.3	91	-0.01
81		3,3'-Dichlorobenzidine	0.450	0.454	-0.9	87	-0.01
82		Chrysene	1.108	1.112	-0.4	91	-0.01
83		Bis(2-ethylhexyl)phthalate	0.714	0.747	-4.6	91	-0.01
84	c	Di-n-octyl phthalate	1.196	1.250	-4.5	91	-0.02
85		Indeno(1,2,3-cd)pyrene	1.336	1.297	2.9	87	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.02
87		Benzo(b)fluoranthene	1.250	1.261	-0.9	89	-0.01

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88	Benzo(k)fluoranthene	1.232	1.272	-3.2	92	-0.01
89 C	Benzo(a)pyrene	1.200	1.199	0.1	88	-0.02
90	Dibenzo(a,h)anthracene	1.210	1.203	0.6	88	-0.03
91	Benzo(g,h,i)perylene	1.165	1.153	1.0	88	-0.02

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