Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
Data File : BG023035.D
Acq On : 12 Jul 2016 14:04
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 13 01:04:25 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 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	109	0.00
2		1,4-Dioxane	0.475	0.460	3.2	109	0.00
3		Pyridine	1.627	1.529	6.0	102	0.00
4		n-Nitrosodimethylamine	0.698	0.667	4.4	103	0.00
5	S	2-Fluorophenol	1.223	1.149	6.1	104	0.00
6		Aniline	2.654	2.578	2.9	104	0.00
7	S	Phenol-d6	1.915	1.906	0.5	106	0.00
8		2-Chlorophenol	1.301	1.297	0.3	108	0.00
9		Benzaldehyde	1.280	1.223	4.5	106	0.00
10	C	Phenol	1.971	2.010	-2.0	109	0.00
11		bis(2-Chloroethyl)ether	1.562	1.469	6.0	102	0.00
12		1,3-Dichlorobenzene	1.593	1.495	6.2	105	0.00
13	C	1,4-Dichlorobenzene	1.595	1.531	4.0	106	0.00
14		1,2-Dichlorobenzene	1.585	1.512	4.6	108	0.00
15		Benzyl Alcohol	1.754	1.785	-1.8	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.863	2.4	107	0.00
17		2-Methylphenol	1.283	1.261	1.7	106	0.00
18		Hexachloroethane	0.673	0.646	4.0	112	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.650	4.5	101	0.00
20		3+4-Methylphenols	1.789	1.840	-2.9	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.600	0.582	3.0	103	0.00
23	S	Nitrobenzene-d5	0.467	0.465	0.4	106	0.00
24		Nitrobenzene	0.496	0.502	-1.2	109	0.00
25		Isophorone	0.939	0.885	5.8	97	0.00
26	C	2-Nitrophenol	0.195	0.194	0.5	99	0.00
27		2,4-Dimethylphenol	0.337	0.330	2.1	104	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.498	5.0	99	0.00
29	C	2,4-Dichlorophenol	0.359	0.363	-1.1	105	0.00
30		1,2,4-Trichlorobenzene	0.411	0.404	1.7	104	0.00
31		Naphthalene	1.018	1.007	1.1	106	0.00
32		Benzoic acid	0.306	0.282	7.8	92	0.02
33		4-Chloroaniline	0.498	0.494	0.8	105	0.00
34	C	Hexachlorobutadiene	0.334	0.309	7.5	100	0.00
35		Caprolactam	0.148	0.129	12.8	92	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.475	3.3	100	0.00
37		2-Methylnaphthalene	0.805	0.811	-0.7	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.855	0.8	100	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.551	-11.1	111	0.00
41	S	2,4,6-Tribromophenol	0.359	0.297	17.3	84	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.485	2.0	96	0.00
43		2,4,5-Trichlorophenol	0.509	0.487	4.3	97	0.00
44	S	2-Fluorobiphenyl	1.270	1.244	2.0	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.527	-1.7	102	0.00
46		2-Chloronaphthalene	1.217	1.240	-1.9	102	0.00
47		2-Nitroaniline	0.519	0.529	-1.9	102	0.00
48		Acenaphthylene	1.826	1.798	1.5	99	0.00
49		Dimethylphthalate	1.634	1.493	8.6	93	0.00
50		2,6-Dinitrotoluene	0.367	0.354	3.5	97	0.00
51	C	Acenaphthene	1.193	1.189	0.3	101	0.00
52		3-Nitroaniline	0.366	0.359	1.9	97	0.00
53	P	2,4-Dinitrophenol	0.252	0.317	-25.8#	121	0.00
54		Dibenzofuran	1.786	1.720	3.7	98	0.00
55	P	4-Nitrophenol	0.307	0.270	12.1	88	0.02
56		2,4-Dinitrotoluene	0.535	0.498	6.9	94	0.00
57		Fluorene	1.537	1.456	5.3	97	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.447	12.2	88	0.02
59		Diethylphthalate	1.662	1.526	8.2	94	0.00
60		4-Chlorophenyl-phenylether	0.936	0.861	8.0	92	0.00
61		4-Nitroaniline	0.396	0.371	6.3	95	0.00
62		Azobenzene	1.590	1.587	0.2	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.136	8.7	81	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.584	-1.4	93	0.00
66		4-Bromophenyl-phenylether	0.260	0.251	3.5	88	0.00
67		Hexachlorobenzene	0.283	0.273	3.5	89	0.00
68		Atrazine	0.256	0.248	3.1	91	0.00
69	C	Pentachlorophenol	0.183	0.166	9.3	79	0.00
70		Phenanthrene	0.980	0.968	1.2	92	0.00
71		Anthracene	0.992	0.978	1.4	93	0.00
72		Carbazole	0.858	0.850	0.9	93	0.00
73		Di-n-butylphthalate	0.954	1.012	-6.1	96	0.00
74	C	Fluoranthene	1.203	1.154	4.1	94	0.02
75	I	Chrysene-d12	1.000	1.000	0.0	93	0.03
76		Benzidine	0.573	0.637	-11.2	100	0.02
77		Pyrene	1.124	1.078	4.1	95	0.00
78	S	Terphenyl-d14	0.646	0.650	-0.6	94	0.01
79		Butylbenzylphthalate	0.445	0.464	-4.3	98	0.02
80		Benzo(a)anthracene	1.119	1.115	0.4	95	0.03
81		3,3'-Dichlorobenzidine	0.491	0.475	3.3	89	0.03
82		Chrysene	1.050	1.042	0.8	95	0.03
83		Bis(2-ethylhexyl)phthalate	0.618	0.636	-2.9	98	0.03
84	c	Di-n-octyl phthalate	1.031	1.066	-3.4	97	0.05
85		Indeno(1,2,3-cd)pyrene	1.338	1.243	7.1	89	0.13
86	I	Perylene-d12	1.000	1.000	0.0	93	0.07
87		Benzo(b)fluoranthene	1.154	1.121	2.9	93	0.06

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88	Benzo(k)fluoranthene	1.095	1.108	-1.2	93	0.07
89 C	Benzo(a)pyrene	1.106	1.079	2.4	92	0.07
90	Dibenzo(a,h)anthracene	1.098	1.047	4.6	90	0.12
91	Benzo(g,h,i)perylene	1.099	1.030	6.3	88	0.14

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