Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062516\
Data File : BG022578.D
Acq On : 25 Jun 2016 17:26
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 01:10:39 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	103	0.00
2		1,4-Dioxane	0.505	0.501	0.8	100	0.00
3		Pyridine	1.414	1.591	-12.5	113	-0.02
4		n-Nitrosodimethylamine	0.809	0.716	11.5	89	-0.01
5	S	2-Fluorophenol	1.247	1.195	4.2	99	0.00
6		Aniline	2.330	2.557	-9.7	111	0.00
7	S	Phenol-d6	1.758	1.776	-1.0	104	0.00
8		2-Chlorophenol	1.292	1.258	2.6	103	0.00
9		Benzaldehyde	0.619	0.682	-10.2	112	0.00
10	C	Phenol	1.858	1.878	-1.1	103	0.00
11		bis(2-Chloroethyl)ether	1.529	1.460	4.5	93	0.00
12		1,3-Dichlorobenzene	1.572	1.488	5.3	97	0.00
13	C	1,4-Dichlorobenzene	1.576	1.522	3.4	102	0.00
14		1,2-Dichlorobenzene	1.498	1.423	5.0	100	0.00
15		Benzyl Alcohol	1.626	1.765	-8.5	104	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.621	4.1	98	0.00
17		2-Methylphenol	1.224	1.240	-1.3	107	0.00
18		Hexachloroethane	0.652	0.615	5.7	95	0.00
19	P	n-Nitroso-di-n-propylamine	1.464	1.472	-0.5	103	0.00
20		3+4-Methylphenols	1.687	1.734	-2.8	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.578	0.565	2.2	104	0.00
23	S	Nitrobenzene-d5	0.473	0.460	2.7	103	0.00
24		Nitrobenzene	0.522	0.500	4.2	104	0.00
25		Isophorone	0.917	0.897	2.2	106	0.00
26	C	2-Nitrophenol	0.188	0.191	-1.6	111	0.00
27		2,4-Dimethylphenol	0.359	0.325	9.5	102	0.00
28		bis(2-Chloroethoxy)methane	0.501	0.480	4.2	106	0.00
29	C	2,4-Dichlorophenol	0.350	0.361	-3.1	109	0.00
30		1,2,4-Trichlorobenzene	0.416	0.392	5.8	103	0.00
31		Naphthalene	1.005	0.959	4.6	103	0.00
32		Benzoic acid	0.216	0.251	-16.2	120	0.06
33		4-Chloroaniline	0.433	0.459	-6.0	107	0.00
34	C	Hexachlorobutadiene	0.323	0.311	3.7	102	0.00
35		Caprolactam	0.110	0.120	-9.1	117	0.03
36	C	4-Chloro-3-methylphenol	0.430	0.446	-3.7	112	0.00
37		2-Methylnaphthalene	0.780	0.770	1.3	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.704	6.8	109	0.00
40	P	Hexachlorocyclopentadiene	0.603	0.550	8.8	101	0.00
41	S	2,4,6-Tribromophenol	0.315	0.327	-3.8	119	0.00
42	C	2,4,6-Trichlorophenol	0.484	0.488	-0.8	112	0.00
43		2,4,5-Trichlorophenol	0.507	0.526	-3.7	120	0.00
44	S	2-Fluorobiphenyl	1.261	1.238	1.8	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.524	1.441	5.4	108	0.00
46		2-Chloronaphthalene	1.254	1.195	4.7	110	0.00
47		2-Nitroaniline	0.484	0.496	-2.5	115	0.00
48		Acenaphthylene	1.818	1.740	4.3	109	0.00
49		Dimethylphthalate	1.659	1.590	4.2	111	0.00
50		2,6-Dinitrotoluene	0.361	0.363	-0.6	114	0.00
51	C	Acenaphthene	1.339	1.329	0.7	112	0.00
52		3-Nitroaniline	0.337	0.354	-5.0	118	0.00
53	P	2,4-Dinitrophenol	0.222	0.241	-8.6	120	0.00
54		Dibenzofuran	1.940	1.863	4.0	112	0.00
55	P	4-Nitrophenol	0.285	0.306	-7.4	124	0.01
56		2,4-Dinitrotoluene	0.499	0.519	-4.0	118	0.00
57		Fluorene	1.548	1.535	0.8	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.552	-5.1	121	0.00
59		Diethylphthalate	1.706	1.648	3.4	113	0.00
60		4-Chlorophenyl-phenylether	0.922	0.901	2.3	114	0.00
61		4-Nitroaniline	0.348	0.365	-4.9	119	0.02
62		Azobenzene	1.847	1.769	4.2	109	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.137	-2.2	115	0.00
65	c	n-Nitrosodiphenylamine	0.582	0.563	3.3	114	0.00
66		4-Bromophenyl-phenylether	0.259	0.246	5.0	114	0.00
67		Hexachlorobenzene	0.274	0.264	3.6	117	0.00
68		Atrazine	0.246	0.248	-0.8	119	0.01
69	C	Pentachlorophenol	0.189	0.193	-2.1	122	0.00
70		Phenanthrene	1.020	0.975	4.4	114	0.00
71		Anthracene	1.050	0.998	5.0	114	0.00
72		Carbazole	0.890	0.875	1.7	116	0.01
73		Di-n-butylphthalate	1.036	1.022	1.4	112	0.00
74	C	Fluoranthene	1.175	1.164	0.9	115	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	120	0.01
76		Benzidine	0.213	0.303	-42.3#	176#	-0.02
77		Pyrene	1.065	1.044	2.0	111	0.00
78	S	Terphenyl-d14	0.755	0.653	13.5	110	0.00
79		Butylbenzylphthalate	0.461	0.464	-0.7	116	0.00
80		Benzo(a)anthracene	1.107	1.116	-0.8	115	0.00
81		3,3'-Dichlorobenzidine	0.457	0.445	2.6	113	0.01
82		Chrysene	1.040	1.036	0.4	114	0.01
83		Bis(2-ethylhexyl)phthalate	0.649	0.619	4.6	112	0.00
84	c	Di-n-octyl phthalate	1.070	1.024	4.3	112	0.00
85		Indeno(1,2,3-cd)pyrene	1.367	1.344	1.7	117	0.02
86	I	Perylene-d12	1.000	1.000	0.0	122	0.01
87		Benzo(b)fluoranthene	1.203	1.141	5.2	114	0.02

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88	Benzo(k)fluoranthene	1.137	1.089	4.2	117	0.01
89 C	Benzo(a)pyrene	1.172	1.111	5.2	115	0.02
90	Dibenzo(a,h)anthracene	1.104	1.060	4.0	119	0.03
91	Benzo(g,h,i)perylene	1.123	1.066	5.1	118	0.02

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