Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062916\
Data File : BG022699.D
Acq On : 29 Jun 2016 18:40
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 30 04:34:47 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	98	-0.01
2		1,4-Dioxane	0.505	0.489	3.2	93	-0.02
3		Pyridine	1.414	1.677	-18.6	113	-0.02
4		n-Nitrosodimethylamine	0.809	0.739	8.7	87	-0.02
5	S	2-Fluorophenol	1.247	1.195	4.2	94	0.00
6		Aniline	2.330	2.660	-14.2	110	0.00
7	S	Phenol-d6	1.758	1.858	-5.7	103	0.00
8		2-Chlorophenol	1.292	1.325	-2.6	103	0.00
9		Benzaldehyde	0.619	0.807	-30.4#	126	0.00
10	C	Phenol	1.858	1.971	-6.1	102	0.00
11		bis(2-Chloroethyl)ether	1.529	1.516	0.9	92	0.00
12		1,3-Dichlorobenzene	1.572	1.543	1.8	96	0.00
13	C	1,4-Dichlorobenzene	1.576	1.570	0.4	99	-0.01
14		1,2-Dichlorobenzene	1.498	1.495	0.2	100	0.00
15		Benzyl Alcohol	1.626	1.913	-17.7	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.712	-1.2	99	-0.01
17		2-Methylphenol	1.224	1.333	-8.9	109	0.00
18		Hexachloroethane	0.652	0.637	2.3	93	-0.01
19	P	n-Nitroso-di-n-propylamine	1.464	1.669	-14.0	111	0.00
20		3+4-Methylphenols	1.687	1.876	-11.2	111	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.578	0.580	-0.3	109	0.00
23	S	Nitrobenzene-d5	0.473	0.462	2.3	107	0.00
24		Nitrobenzene	0.522	0.498	4.6	107	0.00
25		Isophorone	0.917	0.947	-3.3	115	0.00
26	C	2-Nitrophenol	0.188	0.194	-3.2	116	0.00
27		2,4-Dimethylphenol	0.359	0.347	3.3	112	0.00
28		bis(2-Chloroethoxy)methane	0.501	0.503	-0.4	114	0.00
29	C	2,4-Dichlorophenol	0.350	0.366	-4.6	114	0.00
30		1,2,4-Trichlorobenzene	0.416	0.405	2.6	109	0.00
31		Naphthalene	1.005	0.980	2.5	109	0.00
32		Benzoic acid	0.216	0.252	-16.7	124	0.08
33		4-Chloroaniline	0.433	0.496	-14.5	119	-0.01
34	C	Hexachlorobutadiene	0.323	0.318	1.5	107	-0.01
35		Caprolactam	0.110	0.144	-30.9#	144	0.02
36	C	4-Chloro-3-methylphenol	0.430	0.489	-13.7	126	0.00
37		2-Methylnaphthalene	0.780	0.824	-5.6	118	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.686	9.1	125	0.00
40	P	Hexachlorocyclopentadiene	0.603	0.533	11.6	115	-0.01
41	S	2,4,6-Tribromophenol	0.315	0.334	-6.0	142	0.00
42	C	2,4,6-Trichlorophenol	0.484	0.476	1.7	129	0.00
43		2,4,5-Trichlorophenol	0.507	0.500	1.4	134	0.00
44	S	2-Fluorobiphenyl	1.261	1.180	6.4	119	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.524	1.394	8.5	123	0.00
46		2-Chloronaphthalene	1.254	1.126	10.2	122	0.00
47		2-Nitroaniline	0.484	0.462	4.5	126	0.00
48		Acenaphthylene	1.818	1.732	4.7	127	0.00
49		Dimethylphthalate	1.659	1.597	3.7	131	0.00
50		2,6-Dinitrotoluene	0.361	0.358	0.8	132	0.00
51	C	Acenaphthene	1.339	1.174	12.3	116	0.00
52		3-Nitroaniline	0.337	0.343	-1.8	135	0.00
53	P	2,4-Dinitrophenol	0.222	0.238	-7.2	139	0.01
54		Dibenzofuran	1.940	1.847	4.8	130	0.00
55	P	4-Nitrophenol	0.285	0.291	-2.1	139	0.02
56		2,4-Dinitrotoluene	0.499	0.517	-3.6	138	0.00
57		Fluorene	1.548	1.511	2.4	132	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.546	-4.0	141	0.00
59		Diethylphthalate	1.706	1.660	2.7	134	0.00
60		4-Chlorophenyl-phenylether	0.922	0.926	-0.4	137	0.00
61		4-Nitroaniline	0.348	0.355	-2.0	136	0.02
62		Azobenzene	1.847	1.737	6.0	126	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.135	-0.7	132	0.01
65	c	n-Nitrosodiphenylamine	0.582	0.569	2.2	134	0.00
66		4-Bromophenyl-phenylether	0.259	0.260	-0.4	140	0.00
67		Hexachlorobenzene	0.274	0.274	0.0	141	0.00
68		Atrazine	0.246	0.248	-0.8	139	0.00
69	C	Pentachlorophenol	0.189	0.194	-2.6	143	0.00
70		Phenanthrene	1.020	0.963	5.6	132	0.00
71		Anthracene	1.050	0.987	6.0	131	0.00
72		Carbazole	0.890	0.852	4.3	132	0.00
73		Di-n-butylphthalate	1.036	0.940	9.3	120	0.00
74	C	Fluoranthene	1.175	1.110	5.5	127	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	140	0.00
76		Benzidine	0.213	0.398	-86.9#	268#	-0.02
77		Pyrene	1.065	1.017	4.5	126	0.00
78	S	Terphenyl-d14	0.755	0.614	18.7	120	0.00
79		Butylbenzylphthalate	0.461	0.434	5.9	126	-0.01
80		Benzo(a)anthracene	1.107	1.110	-0.3	133	0.00
81		3,3'-Dichlorobenzidine	0.457	0.462	-1.1	136	0.00
82		Chrysene	1.040	1.030	1.0	132	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.600	7.6	127	-0.02
84	c	Di-n-octyl phthalate	1.070	0.986	7.9	126	-0.04
85		Indeno(1,2,3-cd)pyrene	1.367	1.418	-3.7	144	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	143	-0.02
87		Benzo(b)fluoranthene	1.203	1.161	3.5	137	-0.01

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88	Benzo(k)fluoranthene	1.137	1.085	4.6	137	-0.01
89 C	Benzo(a)pyrene	1.172	1.122	4.3	136	-0.01
90	Dibenzo(a,h)anthracene	1.104	1.093	1.0	145	-0.02
91	Benzo(g,h,i)perylene	1.123	1.111	1.1	145	-0.02

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