Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062816\
Data File : BG022680.D
Acq On : 28 Jun 2016 17:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 28 18:51:43 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	96	-0.01
2		1,4-Dioxane	0.505	0.518	-2.6	96	-0.01
3		Pyridine	1.414	1.609	-13.8	106	-0.02
4		n-Nitrosodimethylamine	0.809	0.712	12.0	81	-0.02
5	S	2-Fluorophenol	1.247	1.246	0.1	96	0.00
6		Aniline	2.330	2.646	-13.6	106	0.00
7	S	Phenol-d6	1.758	1.860	-5.8	101	0.00
8		2-Chlorophenol	1.292	1.292	0.0	98	0.00
9		Benzaldehyde	0.619	0.774	-25.0#	117	0.00
10	C	Phenol	1.858	1.997	-7.5	101	0.00
11		bis(2-Chloroethyl)ether	1.529	1.507	1.4	89	0.00
12		1,3-Dichlorobenzene	1.572	1.522	3.2	92	0.00
13	C	1,4-Dichlorobenzene	1.576	1.535	2.6	95	-0.01
14		1,2-Dichlorobenzene	1.498	1.498	0.0	97	0.00
15		Benzyl Alcohol	1.626	1.794	-10.3	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.645	2.7	92	0.00
17		2-Methylphenol	1.224	1.271	-3.8	102	0.00
18		Hexachloroethane	0.652	0.619	5.1	89	-0.01
19	P	n-Nitroso-di-n-propylamine	1.464	1.480	-1.1	96	0.00
20		3+4-Methylphenols	1.687	1.781	-5.6	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.578	0.566	2.1	96	0.00
23	S	Nitrobenzene-d5	0.473	0.473	0.0	99	0.00
24		Nitrobenzene	0.522	0.512	1.9	99	0.00
25		Isophorone	0.917	0.894	2.5	98	-0.01
26	C	2-Nitrophenol	0.188	0.199	-5.9	107	0.00
27		2,4-Dimethylphenol	0.359	0.341	5.0	100	0.00
28		bis(2-Chloroethoxy)methane	0.501	0.498	0.6	102	0.00
29	C	2,4-Dichlorophenol	0.350	0.377	-7.7	106	0.00
30		1,2,4-Trichlorobenzene	0.416	0.410	1.4	99	0.00
31		Naphthalene	1.005	0.987	1.8	99	0.00
32		Benzoic acid	0.216	0.254	-17.6	113	0.06
33		4-Chloroaniline	0.433	0.484	-11.8	105	-0.01
34	C	Hexachlorobutadiene	0.323	0.313	3.1	95	-0.01
35		Caprolactam	0.110	0.125	-13.6	113	0.02
36	C	4-Chloro-3-methylphenol	0.430	0.464	-7.9	108	0.00
37		2-Methylnaphthalene	0.780	0.781	-0.1	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.714	5.4	106	0.00
40	P	Hexachlorocyclopentadiene	0.603	0.517	14.3	91	-0.01
41	S	2,4,6-Tribromophenol	0.315	0.336	-6.7	116	0.00
42	C	2,4,6-Trichlorophenol	0.484	0.494	-2.1	109	0.00
43		2,4,5-Trichlorophenol	0.507	0.541	-6.7	118	0.00
44	S	2-Fluorobiphenyl	1.261	1.244	1.3	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.524	1.447	5.1	104	0.00
46		2-Chloronaphthalene	1.254	1.231	1.8	108	0.00
47		2-Nitroaniline	0.484	0.482	0.4	107	0.00
48		Acenaphthylene	1.818	1.791	1.5	107	0.00
49		Dimethylphthalate	1.659	1.604	3.3	107	0.00
50		2,6-Dinitrotoluene	0.361	0.371	-2.8	111	0.00
51	C	Acenaphthene	1.339	1.182	11.7	95	0.00
52		3-Nitroaniline	0.337	0.360	-6.8	115	0.00
53	P	2,4-Dinitrophenol	0.222	0.245	-10.4	116	0.00
54		Dibenzofuran	1.940	1.912	1.4	110	0.00
55	P	4-Nitrophenol	0.285	0.297	-4.2	116	0.01
56		2,4-Dinitrotoluene	0.499	0.544	-9.0	118	0.00
57		Fluorene	1.548	1.547	0.1	110	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.548	-4.4	115	0.00
59		Diethylphthalate	1.706	1.644	3.6	108	0.00
60		4-Chlorophenyl-phenylether	0.922	0.910	1.3	110	0.00
61		4-Nitroaniline	0.348	0.357	-2.6	112	0.02
62		Azobenzene	1.847	1.754	5.0	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.146	-9.0	116	0.00
65	c	n-Nitrosodiphenylamine	0.582	0.567	2.6	109	0.00
66		4-Bromophenyl-phenylether	0.259	0.254	1.9	112	0.00
67		Hexachlorobenzene	0.274	0.277	-1.1	117	0.00
68		Atrazine	0.246	0.241	2.0	111	0.00
69	C	Pentachlorophenol	0.189	0.186	1.6	112	0.00
70		Phenanthrene	1.020	0.995	2.5	111	0.00
71		Anthracene	1.050	1.024	2.5	111	0.00
72		Carbazole	0.890	0.890	0.0	112	0.00
73		Di-n-butylphthalate	1.036	1.029	0.7	107	0.00
74	C	Fluoranthene	1.175	1.198	-2.0	112	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	120	0.00
76		Benzidine	0.213	0.320	-50.2#	184#	-0.02
77		Pyrene	1.065	1.046	1.8	110	0.00
78	S	Terphenyl-d14	0.755	0.649	14.0	109	0.00
79		Butylbenzylphthalate	0.461	0.443	3.9	110	0.00
80		Benzo(a)anthracene	1.107	1.106	0.1	113	0.00
81		3,3'-Dichlorobenzidine	0.457	0.437	4.4	110	0.00
82		Chrysene	1.040	1.042	-0.2	114	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.600	7.6	108	-0.01
84	c	Di-n-octyl phthalate	1.070	0.992	7.3	108	-0.02
85		Indeno(1,2,3-cd)pyrene	1.367	1.305	4.5	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	118	0.00
87		Benzo(b)fluoranthene	1.203	1.165	3.2	113	0.00

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88	Benzo(k)fluoranthene	1.137	1.115	1.9	116	0.00
89 C	Benzo(a)pyrene	1.172	1.142	2.6	114	0.00
90	Dibenzo(a,h)anthracene	1.104	1.063	3.7	116	0.00
91	Benzo(g,h,i)perylene	1.123	1.017	9.4	109	0.00

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