Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG110417\
Data File : BG030717.D
Acq On : 4 Nov 2017 1:54
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 04 01:39:18 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 03 18:14:01 2017
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	0.486	0.476	2.1	99	0.00
3		Pyridine	1.415	1.473	-4.1	104	0.00
4		n-Nitrosodimethylamine	0.661	0.645	2.4	99	0.00
5	S	2-Fluorophenol	1.197	1.221	-2.0	102	0.00
6		Aniline	2.081	2.215	-6.4	105	0.00
7	S	Phenol-d6	1.680	1.762	-4.9	104	0.00
8		2-Chlorophenol	1.372	1.363	0.7	101	0.00
9		Benzaldehyde	0.845	1.004	-18.8	118	0.00
10	C	Phenol	1.812	1.775	2.0	97	0.00
11		bis(2-Chloroethyl)ether	1.320	1.400	-6.1	104	0.00
12		1,3-Dichlorobenzene	1.541	1.547	-0.4	102	0.00
13	C	1,4-Dichlorobenzene	1.573	1.578	-0.3	101	0.00
14		1,2-Dichlorobenzene	1.517	1.528	-0.7	102	0.00
15		Benzyl Alcohol	1.184	1.312	-10.8	111	0.00
16		2,2'-oxybis(1-Chloropropane	2.242	1.915	14.6	86	0.00
17		2-Methylphenol	1.204	1.215	-0.9	100	0.00
18		Hexachloroethane	0.536	0.538	-0.4	100	0.00
19	P	n-Nitroso-di-n-propylamine	1.143	1.263	-10.5	111	0.00
20		3+4-Methylphenols	1.696	1.786	-5.3	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.482	0.488	-1.2	108	0.00
23	S	Nitrobenzene-d5	0.315	0.333	-5.7	110	0.00
24		Nitrobenzene	0.354	0.340	4.0	100	0.00
25		Isophorone	0.657	0.630	4.1	100	0.00
26	C	2-Nitrophenol	0.169	0.170	-0.6	102	0.00
27		2,4-Dimethylphenol	0.306	0.266	13.1	92	0.00
28		bis(2-Chloroethoxy)methane	0.403	0.416	-3.2	109	0.00
29	C	2,4-Dichlorophenol	0.326	0.321	1.5	104	0.00
30		1,2,4-Trichlorobenzene	0.353	0.358	-1.4	108	0.00
31		Naphthalene	0.997	0.971	2.6	104	0.00
32		Benzoic acid	0.256	0.243	5.1	94	0.00
33		4-Chloroaniline	0.419	0.433	-3.3	109	0.00
34	C	Hexachlorobutadiene	0.219	0.230	-5.0	111	0.00
35		Caprolactam	0.108	0.112	-3.7	110	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.341	5.0	101	0.00
37		2-Methylnaphthalene	0.726	0.744	-2.5	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
39		1,2,4,5-Tetrachlorobenzene	0.730	0.742	-1.6	115	0.00
40	P	Hexachlorocyclopentadiene	0.247	0.291	-17.8	129	0.00
41	S	2,4,6-Tribromophenol	0.258	0.280	-8.5	119	0.00
42	C	2,4,6-Trichlorophenol	0.458	0.425	7.2	103	0.00
43		2,4,5-Trichlorophenol	0.471	0.481	-2.1	114	0.00
44	S	2-Fluorobiphenyl	1.364	1.358	0.4	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.656	3.7	108	0.00
46		2-Chloronaphthalene	1.254	1.198	4.5	107	0.00
47		2-Nitroaniline	0.344	0.350	-1.7	109	0.00
48		Acenaphthylene	1.962	1.956	0.3	112	0.00
49		Dimethylphthalate	1.560	1.582	-1.4	112	0.00
50		2,6-Dinitrotoluene	0.327	0.343	-4.9	111	0.00
51	C	Acenaphthene	1.277	1.235	3.3	107	0.00
52		3-Nitroaniline	0.353	0.371	-5.1	114	0.00
53	P	2,4-Dinitrophenol	0.153	0.146	4.6	106	0.00
54		Dibenzofuran	1.823	1.909	-4.7	118	0.00
55	P	4-Nitrophenol	0.291	0.276	5.2	102	0.00
56		2,4-Dinitrotoluene	0.443	0.485	-9.5	117	0.00
57		Fluorene	1.506	1.503	0.2	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.393	0.427	-8.7	119	0.00
59		Diethylphthalate	1.555	1.575	-1.3	113	0.00
60		4-Chlorophenyl-phenylether	0.803	0.868	-8.1	121	0.00
61		4-Nitroaniline	0.362	0.384	-6.1	117	0.00
62		Azobenzene	1.311	1.345	-2.6	116	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.117	-11.4	128	0.00
65	c	n-Nitrosodiphenylamine	0.599	0.604	-0.8	120	0.00
66		4-Bromophenyl-phenylether	0.229	0.231	-0.9	119	0.00
67		Hexachlorobenzene	0.260	0.254	2.3	117	0.00
68		Atrazine	0.214	0.226	-5.6	122	0.00
69	C	Pentachlorophenol	0.149	0.157	-5.4	119	0.00
70		Phenanthrene	1.033	1.024	0.9	118	0.00
71		Anthracene	1.037	1.025	1.2	117	0.00
72		Carbazole	0.997	0.959	3.8	114	0.00
73		Di-n-butylphthalate	1.146	1.135	1.0	116	0.00
74	C	Fluoranthene	1.208	1.251	-3.6	122	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	130	0.00
76		Benzidine	0.604	0.600	0.7	118	0.00
77		Pyrene	1.213	1.193	1.6	120	0.00
78	S	Terphenyl-d14	0.879	0.850	3.3	118	0.00
79		Butylbenzylphthalate	0.517	0.496	4.1	117	0.00
80		Benzo(a)anthracene	1.174	1.195	-1.8	124	0.00
81		3,3'-Dichlorobenzidine	0.485	0.488	-0.6	123	0.00
82		Chrysene	1.125	1.118	0.6	123	0.00
83		Bis(2-ethylhexyl)phthalate	0.742	0.697	6.1	114	0.00
84	c	Di-n-octyl phthalate	1.293	1.182	8.6	111	-0.01
85		Indeno(1,2,3-cd)pyrene	1.383	1.421	-2.7	126	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	130	-0.01
87		Benzo(b)fluoranthene	1.145	1.158	-1.1	123	0.00

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88	Benzo(k)fluoranthene	1.141	1.168	-2.4	126	-0.01
89 C	Benzo(a)pyrene	1.101	1.108	-0.6	123	0.00
90	Dibenzo(a,h)anthracene	1.114	1.159	-4.0	126	-0.01
91	Benzo(g,h,i)perylene	1.035	1.116	-7.8	129	-0.02

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