Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101618\
Data File : BG037374.D
Acq On : 17 Oct 2018 00:39
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 17 02:16:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 11 13:59:58 2018
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	141	-0.01
2		1,4-Dioxane	0.638	0.602	5.6	138	0.00
3		Pyridine	1.515	1.527	-0.8	136	0.00
4		n-Nitrosodimethylamine	0.588	0.561	4.6	135	0.00
5	S	2-Fluorophenol	1.136	1.200	-5.6	146	0.00
6		Aniline	2.258	2.192	2.9	135	0.00
7	S	Phenol-d6	1.725	1.803	-4.5	143	0.00
8		2-Chlorophenol	1.330	1.369	-2.9	139	0.00
9		Benzaldehyde	1.188	1.145	3.6	134	0.00
10	C	Phenol	1.817	1.892	-4.1	142	0.00
11		bis(2-Chloroethyl)ether	1.484	1.445	2.6	137	-0.01
12		1,3-Dichlorobenzene	1.564	1.549	1.0	141	-0.01
13	C	1,4-Dichlorobenzene	1.583	1.560	1.5	139	-0.01
14		1,2-Dichlorobenzene	1.534	1.473	4.0	137	0.00
15		Benzyl Alcohol	1.349	1.365	-1.2	137	0.00
16		2,2'-oxybis(1-Chloropropane	2.934	2.623	10.6	125	-0.02
17		2-Methylphenol	1.217	1.242	-2.1	139	0.00
18		Hexachloroethane	0.540	0.541	-0.2	139	-0.01
19	P	n-Nitroso-di-n-propylamine	1.306	1.248	4.4	133	0.00
20		3+4-Methylphenols	1.701	1.803	-6.0	142	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	137	-0.01
22		Acetophenone	0.504	0.492	2.4	133	0.00
23	S	Nitrobenzene-d5	0.304	0.356	-17.1	151#	0.00
24		Nitrobenzene	0.326	0.374	-14.7	148	-0.01
25		Isophorone	0.694	0.669	3.6	129	0.00
26	C	2-Nitrophenol	0.116	0.171	-47.4#	203#	0.00
27		2,4-Dimethylphenol	0.290	0.297	-2.4	138	0.00
28		bis(2-Chloroethoxy)methane	0.434	0.431	0.7	133	-0.01
29	C	2,4-Dichlorophenol	0.295	0.334	-13.2	144	0.00
30		1,2,4-Trichlorobenzene	0.359	0.355	1.1	137	-0.01
31		Naphthalene	0.972	0.970	0.2	137	-0.01
32		Benzoic acid	0.160	0.212	-32.5#	177#	0.00
33		4-Chloroaniline	0.456	0.457	-0.2	133	0.00
34	C	Hexachlorobutadiene	0.223	0.212	4.9	132	-0.02
35		Caprolactam	0.120	0.143	-19.2	150	0.00
36	C	4-Chloro-3-methylphenol	0.348	0.379	-8.9	139	0.00
37		2-Methylnaphthalene	0.709	0.713	-0.6	137	-0.01
38		1-Methylnaphthalene	0.693	0.691	0.3	135	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	135	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.648	0.637	1.7	135	0.00
41	P	Hexachlorocyclopentadiene	0.267	0.296	-10.9	149	-0.01
42	S	2,4,6-Tribromophenol	0.196	0.228	-16.3	148	-0.01
43	C	2,4,6-Trichlorophenol	0.376	0.444	-18.1	151#	-0.01
44		2,4,5-Trichlorophenol	0.404	0.473	-17.1	149	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.332	1.286	3.5	133	-0.01
46		1,1'-Biphenyl	1.645	1.634	0.7	135	-0.01
47		2-Chloronaphthalene	1.243	1.239	0.3	137	-0.01
48		2-Nitroaniline	0.300	0.389	-29.7#	165#	0.00
49		Acenaphthylene	1.901	1.885	0.8	135	0.00
50		Dimethylphthalate	1.572	1.559	0.8	131	-0.01
51		2,6-Dinitrotoluene	0.284	0.349	-22.9	160#	0.00
52	C	Acenaphthene	1.174	1.153	1.8	133	-0.01
53		3-Nitroaniline	0.329	0.387	-17.6	153#	0.00
54	P	2,4-Dinitrophenol	0.079	0.087	-10.1	154#	0.00
55		Dibenzofuran	1.905	1.875	1.6	134	-0.01
56	P	4-Nitrophenol	0.256	0.291	-13.7	147	0.00
57		2,4-Dinitrotoluene	0.356	0.479	-34.6#	175#	-0.01
58		Fluorene	1.441	1.407	2.4	134	-0.01
59		2,3,4,6-Tetrachlorophenol	0.358	0.410	-14.5	142	0.00
60		Diethylphthalate	1.629	1.606	1.4	129	-0.01
61		4-Chlorophenyl-phenylether	0.842	0.823	2.3	133	-0.01
62		4-Nitroaniline	0.345	0.399	-15.7	151#	0.00
63		Azobenzene	1.562	1.497	4.2	131	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	134	-0.01
65		4,6-Dinitro-2-methylphenol	0.056	0.084	-50.0#	192#	-0.01
66	c	n-Nitrosodiphenylamine	0.587	0.594	-1.2	133	-0.01
67		4-Bromophenyl-phenylether	0.216	0.219	-1.4	134	-0.01
68		Hexachlorobenzene	0.224	0.224	0.0	134	-0.01
69		Atrazine	0.218	0.226	-3.7	132	-0.01
70	C	Pentachlorophenol	0.144	0.152	-5.6	136	0.00
71		Phenanthrene	1.018	0.998	2.0	131	-0.01
72		Anthracene	0.996	0.986	1.0	132	-0.01
73		Carbazole	1.027	1.009	1.8	130	0.00
74		Di-n-butylphthalate	1.100	1.144	-4.0	129	-0.02
75	C	Fluoranthene	1.190	1.150	3.4	127	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	131	-0.02
77		Benzidine	0.765	0.635	17.0	98	-0.01
78		Pyrene	1.303	1.296	0.5	129	-0.01
79	S	Terphenyl-d14	0.952	0.910	4.4	124	-0.01
80		Butylbenzylphthalate	0.507	0.569	-12.2	141	-0.02
81		Benzo(a)anthracene	1.198	1.179	1.6	128	-0.01
82		3,3'-Dichlorobenzidine	0.462	0.468	-1.3	126	-0.02
83		Chrysene	1.143	1.131	1.0	130	-0.01
84		Bis(2-ethylhexyl)phthalate	0.728	0.797	-9.5	137	-0.03
85	c	Di-n-octyl phthalate	1.182	1.304	-10.3	138	-0.04
86		Indeno(1,2,3-cd)pyrene	1.266	1.289	-1.8	130	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	133	-0.02

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88	Benzo(b)fluoranthene	1.087	1.095	-0.7	132	-0.02
89	Benzo(k)fluoranthene	1.073	1.062	1.0	130	-0.02
90 C	Benzo(a)pyrene	1.043	1.048	-0.5	131	-0.02
91	Dibenzo(a,h)anthracene	0.998	0.985	1.3	129	-0.05
92	Benzo(g,h,i)perylene	0.996	0.995	0.1	130	-0.05

(#) = Out of Range SPCC's out = 0 CCC's out = 1