Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110818\
Data File : BG037901.D
Acq On : 9 Nov 2018 2:23
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 09 14:43:40 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 14:06:42 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	98	-0.02
2		1,4-Dioxane	0.607	0.578	4.8	97	0.00
3		Pyridine	1.529	1.434	6.2	94	0.00
4		n-Nitrosodimethylamine	0.545	0.547	-0.4	101	0.00
5	S	2-Fluorophenol	1.196	1.177	1.6	98	-0.02
6		Aniline	2.207	2.181	1.2	98	-0.02
7	S	Phenol-d6	1.809	1.762	2.6	96	-0.02
8		2-Chlorophenol	1.399	1.387	0.9	98	-0.02
9		Benzaldehyde	1.132	0.968	14.5	85	-0.02
10	C	Phenol	1.909	1.874	1.8	97	0.00
11		bis(2-Chloroethyl)ether	1.450	1.411	2.7	98	-0.02
12		1,3-Dichlorobenzene	1.558	1.559	-0.1	101	-0.02
13	C	1,4-Dichlorobenzene	1.594	1.595	-0.1	101	-0.02
14		1,2-Dichlorobenzene	1.533	1.514	1.2	99	-0.02
15		Benzyl Alcohol	1.337	1.278	4.4	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.594	2.596	-0.1	100	-0.02
17		2-Methylphenol	1.262	1.237	2.0	96	0.00
18		Hexachloroethane	0.543	0.580	-6.8	106	-0.02
19	P	n-Nitroso-di-n-propylamine	1.246	1.191	4.4	93	-0.02
20		3+4-Methylphenols	1.804	1.767	2.1	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	-0.02
22		Acetophenone	0.502	0.473	5.8	93	-0.02
23	S	Nitrobenzene-d5	0.357	0.361	-1.1	100	0.00
24		Nitrobenzene	0.371	0.372	-0.3	98	-0.02
25		Isophorone	0.652	0.631	3.2	93	-0.02
26	C	2-Nitrophenol	0.167	0.190	-13.8	109	-0.02
27		2,4-Dimethylphenol	0.298	0.289	3.0	95	0.00
28		bis(2-Chloroethoxy)methane	0.427	0.420	1.6	96	-0.02
29	C	2,4-Dichlorophenol	0.332	0.338	-1.8	100	0.00
30		1,2,4-Trichlorobenzene	0.361	0.367	-1.7	102	-0.02
31		Naphthalene	0.982	0.970	1.2	99	-0.02
32		Benzoic acid	0.194	0.206	-6.2	102	0.00
33		4-Chloroaniline	0.462	0.452	2.2	96	-0.02
34	C	Hexachlorobutadiene	0.214	0.223	-4.2	102	-0.02
35		Caprolactam	0.133	0.143	-7.5	103	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.361	3.7	94	-0.02
37		2-Methylnaphthalene	0.716	0.713	0.4	98	-0.02
38		1-Methylnaphthalene	0.695	0.688	1.0	99	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.645	0.677	-5.0	101	-0.02
41	P	Hexachlorocyclopentadiene	0.308	0.342	-11.0	104	-0.02
42	S	2,4,6-Tribromophenol	0.218	0.242	-11.0	102	0.00
43	C	2,4,6-Trichlorophenol	0.431	0.453	-5.1	98	-0.02
44		2,4,5-Trichlorophenol	0.469	0.485	-3.4	97	0.00

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.326	1.323	0.2	98	-0.02
46		1,1'-Biphenyl	1.656	1.668	-0.7	98	-0.02
47		2-Chloronaphthalene	1.253	1.260	-0.6	98	-0.02
48		2-Nitroaniline	0.380	0.420	-10.5	105	0.00
49		Acenaphthylene	1.885	1.899	-0.7	97	0.00
50		Dimethylphthalate	1.547	1.589	-2.7	99	0.00
51		2,6-Dinitrotoluene	0.342	0.374	-9.4	102	0.00
52	C	Acenaphthene	1.161	1.150	0.9	97	-0.02
53		3-Nitroaniline	0.380	0.421	-10.8	103	0.00
54	P	2,4-Dinitrophenol	0.100	0.130	-30.0#	127	0.00
55		Dibenzofuran	1.923	1.893	1.6	96	-0.02
56	P	4-Nitrophenol	0.281	0.297	-5.7	98	0.00
57		2,4-Dinitrotoluene	0.449	0.535	-19.2	109	0.00
58		Fluorene	1.440	1.425	1.0	97	-0.02
59		2,3,4,6-Tetrachlorophenol	0.402	0.434	-8.0	103	-0.02
60		Diethylphthalate	1.588	1.710	-7.7	104	-0.02
61		4-Chlorophenyl-phenylether	0.840	0.840	0.0	97	-0.02
62		4-Nitroaniline	0.378	0.420	-11.1	103	0.00
63		Azobenzene	1.516	1.509	0.5	98	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	0.093	0.115	-23.7	124	0.00
66	c	n-Nitrosodiphenylamine	0.602	0.594	1.3	100	-0.02
67		4-Bromophenyl-phenylether	0.220	0.217	1.4	100	-0.02
68		Hexachlorobenzene	0.228	0.226	0.9	101	0.00
69		Atrazine	0.218	0.197	9.6	89	-0.02
70	C	Pentachlorophenol	0.152	0.156	-2.6	101	-0.02
71		Phenanthrene	1.016	1.002	1.4	101	-0.02
72		Anthracene	0.998	1.006	-0.8	103	0.00
73		Carbazole	0.998	1.027	-2.9	104	0.00
74		Di-n-butylphthalate	1.110	1.165	-5.0	104	-0.02
75	C	Fluoranthene	1.153	1.160	-0.6	102	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	106	-0.02
77		Benzidine	0.680	0.536	21.2	81	-0.02
78		Pyrene	1.307	1.283	1.8	104	0.00
79	S	Terphenyl-d14	0.936	0.911	2.7	103	-0.02
80		Butylbenzylphthalate	0.535	0.573	-7.1	108	-0.02
81		Benzo(a)anthracene	1.183	1.174	0.8	105	-0.02
82		3,3'-Dichlorobenzidine	0.459	0.448	2.4	99	-0.02
83		Chrysene	1.138	1.139	-0.1	106	-0.02
84		Bis(2-ethylhexyl)phthalate	0.750	0.812	-8.3	109	-0.03
85	c	Di-n-octyl phthalate	1.223	1.345	-10.0	111	-0.04
86		Indeno(1,2,3-cd)pyrene	1.220	1.049	14.0	89	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	104	-0.03

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88	Benzo(b)fluoranthene	1.096	1.103	-0.6	104	-0.03
89	Benzo(k)fluoranthene	1.074	1.080	-0.6	104	-0.02
90 C	Benzo(a)pyrene	1.042	1.052	-1.0	104	-0.03
91	Dibenzo(a,h)anthracene	0.961	0.828	13.8	87	-0.05
92	Benzo(g,h,i)perylene	0.972	0.827	14.9	88	-0.06

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