Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011119\
Data File : BG039105.D
Acq On : 12 Jan 2019 5:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 12 06:29:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 09 14:23:53 2019
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	149	0.00
2		1,4-Dioxane	0.413	0.466	-12.8	170#	0.00
3		Pyridine	1.122	1.390	-23.9	179#	0.00
4		n-Nitrosodimethylamine	0.505	0.546	-8.1	154#	0.00
5	S	2-Fluorophenol	1.054	1.160	-10.1	158#	0.00
6		Aniline	1.822	1.942	-6.6	152#	0.00
7	S	Phenol-d6	1.517	1.719	-13.3	163#	0.00
8		2-Chlorophenol	1.243	1.407	-13.2	162#	0.00
9		Benzaldehyde	0.875	0.904	-3.3	163#	0.00
10	C	Phenol	1.521	1.739	-14.3	164#	0.00
11		bis(2-Chloroethyl)ether	1.163	1.266	-8.9	157#	0.00
12		1,3-Dichlorobenzene	1.489	1.550	-4.1	154#	0.00
13	C	1,4-Dichlorobenzene	1.508	1.593	-5.6	155#	0.00
14		1,2-Dichlorobenzene	1.438	1.605	-11.6	164#	0.00
15		Benzyl Alcohol	1.143	1.331	-16.4	164#	0.00
16		2,2'-oxybis(1-Chloropropane	2.229	2.464	-10.5	162#	0.00
17		2-Methylphenol	1.056	1.208	-14.4	163#	0.00
18		Hexachloroethane	0.512	0.560	-9.4	161#	0.00
19	P	n-Nitroso-di-n-propylamine	0.996	1.209	-21.4	179#	0.00
20		3+4-Methylphenols	1.506	1.823	-21.0	173#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	154#	0.00
22		Acetophenone	0.522	0.585	-12.1	171#	0.00
23	S	Nitrobenzene-d5	0.366	0.390	-6.6	163#	0.00
24		Nitrobenzene	0.369	0.387	-4.9	162#	0.00
25		Isophorone	0.630	0.720	-14.3	175#	0.00
26	C	2-Nitrophenol	0.200	0.221	-10.5	167#	0.00
27		2,4-Dimethylphenol	0.281	0.310	-10.3	166#	0.00
28		bis(2-Chloroethoxy)methane	0.378	0.438	-15.9	175#	0.00
29	C	2,4-Dichlorophenol	0.349	0.371	-6.3	158#	0.00
30		1,2,4-Trichlorobenzene	0.420	0.446	-6.2	164#	0.00
31		Naphthalene	1.003	1.068	-6.5	166#	0.00
32		Benzoic acid	0.149	0.127	14.8	134	0.00
33		4-Chloroaniline	0.449	0.488	-8.7	163#	0.00
34	C	Hexachlorobutadiene	0.289	0.288	0.3	152#	0.00
35		Caprolactam	0.140	0.142	-1.4	156#	0.00
36	C	4-Chloro-3-methylphenol	0.374	0.406	-8.6	166#	0.00
37		2-Methylnaphthalene	0.752	0.756	-0.5	156#	-0.01
38		1-Methylnaphthalene	0.722	0.728	-0.8	156#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	145	0.00
40		1,2,4,5-Tetrachlorobenzene	0.751	0.787	-4.8	151#	-0.01
41	P	Hexachlorocyclopentadiene	0.369	0.339	8.1	129	0.00
42	S	2,4,6-Tribromophenol	0.335	0.293	12.5	125	-0.01
43	C	2,4,6-Trichlorophenol	0.473	0.499	-5.5	149	0.00
44		2,4,5-Trichlorophenol	0.500	0.521	-4.2	145	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.461	1.577	-7.9	157#	0.00
46		1,1'-Biphenyl	1.585	1.692	-6.8	155#	-0.01
47		2-Chloronaphthalene	1.282	1.365	-6.5	154#	0.00
48		2-Nitroaniline	0.359	0.383	-6.7	149	0.00
49		Acenaphthylene	1.928	2.112	-9.5	159#	-0.01
50		Dimethylphthalate	1.690	1.856	-9.8	161#	0.00
51		2,6-Dinitrotoluene	0.378	0.408	-7.9	155#	0.00
52	C	Acenaphthene	1.158	1.250	-7.9	159#	0.00
53		3-Nitroaniline	0.383	0.391	-2.1	147	0.00
54	P	2,4-Dinitrophenol	0.195	0.303	-55.4#	211#	0.00
55		Dibenzofuran	2.045	2.122	-3.8	153#	0.00
56	P	4-Nitrophenol	0.276	0.243	12.0	123	0.00
57		2,4-Dinitrotoluene	0.542	0.534	1.5	144	0.00
58		Fluorene	1.540	1.514	1.7	144	-0.01
59		2,3,4,6-Tetrachlorophenol	0.472	0.458	3.0	141	0.00
60		Diethylphthalate	1.741	1.717	1.4	146	0.00
61		4-Chlorophenyl-phenylether	0.970	0.947	2.4	141	0.00
62		4-Nitroaniline	0.399	0.373	6.5	131	0.00
63		Azobenzene	1.362	1.350	0.9	147	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
65		4,6-Dinitro-2-methylphenol	0.148	0.137	7.4	112	0.00
66	c	n-Nitrosodiphenylamine	0.593	0.662	-11.6	137	-0.01
67		4-Bromophenyl-phenylether	0.257	0.284	-10.5	136	-0.01
68		Hexachlorobenzene	0.281	0.302	-7.5	135	0.00
69		Atrazine	0.252	0.257	-2.0	124	0.00
70	C	Pentachlorophenol	0.150	0.158	-5.3	124	0.00
71		Phenanthrene	1.057	1.124	-6.3	132	-0.01
72		Anthracene	1.045	1.159	-10.9	138	0.00
73		Carbazole	1.032	1.097	-6.3	133	0.00
74		Di-n-butylphthalate	1.148	1.170	-1.9	127	0.00
75	C	Fluoranthene	1.311	1.312	-0.1	125	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	119	-0.01
77		Benzidine	0.612	0.516	15.7	92	0.00
78		Pyrene	1.275	1.342	-5.3	123	0.00
79	S	Terphenyl-d14	1.081	1.087	-0.6	121	0.00
80		Butylbenzylphthalate	0.485	0.538	-10.9	130	-0.01
81		Benzo(a)anthracene	1.217	1.306	-7.3	125	-0.01
82		3,3'-Dichlorobenzidine	0.496	0.531	-7.1	125	-0.01
83		Chrysene	1.158	1.348	-16.4	137	-0.01
84		Bis(2-ethylhexyl)phthalate	0.673	0.793	-17.8	138	-0.01
85	c	Di-n-octyl phthalate	1.138	1.291	-13.4	132	-0.02
86		Indeno(1,2,3-cd)pyrene	1.384	1.495	-8.0	124	-0.03
87	I	Perylene-d12	1.000	1.000	0.0	120	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.103	1.162	-5.3	127	-0.02
89	Benzo(k)fluoranthene	1.090	1.173	-7.6	127	-0.02
90 C	Benzo(a)pyrene	1.066	1.147	-7.6	128	-0.02
91	Dibenzo(a,h)anthracene	1.060	1.112	-4.9	125	-0.04
92	Benzo(g,h,i)perylene	1.050	1.077	-2.6	123	-0.04

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