Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG051119\
Data File : BG040892.D
Acq On : 12 May 2019 4:04
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 13 04:32:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 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	128	-0.02
2		1,4-Dioxane	0.516	0.485	6.0	125	-0.02
3		Pyridine	1.496	1.428	4.5	121	-0.02
4		n-Nitrosodimethylamine	0.830	0.761	8.3	123	-0.02
5	S	2-Fluorophenol	1.135	1.122	1.1	131	-0.02
6		Aniline	2.315	2.173	6.1	123	-0.02
7	S	Phenol-d6	1.747	1.706	2.3	129	-0.02
8		2-Chlorophenol	1.385	1.352	2.4	127	-0.03
9		Benzaldehyde	1.083	1.037	4.2	130	-0.02
10	C	Phenol	1.878	1.838	2.1	129	-0.02
11		bis(2-Chloroethyl)ether	1.408	1.332	5.4	125	-0.03
12		1,3-Dichlorobenzene	1.512	1.520	-0.5	133	-0.02
13	C	1,4-Dichlorobenzene	1.533	1.498	2.3	129	-0.02
14		1,2-Dichlorobenzene	1.478	1.478	0.0	133	-0.03
15		Benzyl Alcohol	1.818	1.610	11.4	116	-0.03
16		2,2'-oxybis(1-Chloropropane	2.804	2.256	19.5	106	-0.02
17		2-Methylphenol	1.329	1.285	3.3	128	-0.02
18		Hexachloroethane	0.629	0.602	4.3	127	-0.03
19	P	n-Nitroso-di-n-propylamine	1.573	1.390	11.6	119	-0.03
20		3+4-Methylphenols	1.851	1.779	3.9	126	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	126	-0.03
22		Acetophenone	0.556	0.558	-0.4	126	-0.03
23	S	Nitrobenzene-d5	0.433	0.460	-6.2	135	-0.02
24		Nitrobenzene	0.462	0.484	-4.8	135	-0.02
25		Isophorone	0.964	0.903	6.3	119	-0.04
26	C	2-Nitrophenol	0.166	0.207	-24.7#	160#	-0.02
27		2,4-Dimethylphenol	0.311	0.310	0.3	127	-0.03
28		bis(2-Chloroethoxy)methane	0.484	0.457	5.6	119	-0.02
29	C	2,4-Dichlorophenol	0.353	0.376	-6.5	133	-0.03
30		1,2,4-Trichlorobenzene	0.392	0.428	-9.2	138	-0.02
31		Naphthalene	1.063	1.073	-0.9	126	-0.03
32		Benzoic acid	0.213	0.260	-22.1	160#	-0.02
33		4-Chloroaniline	0.471	0.470	0.2	128	-0.03
34	C	Hexachlorobutadiene	0.289	0.337	-16.6	148	-0.03
35		Caprolactam	0.139	0.139	0.0	125	-0.02
36	C	4-Chloro-3-methylphenol	0.445	0.466	-4.7	136	-0.02
37		2-Methylnaphthalene	0.794	0.809	-1.9	128	-0.03
38		1-Methylnaphthalene	0.777	0.784	-0.9	128	-0.03
39	I	Acenaphthene-d10	1.000	1.000	0.0	136	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.745	0.793	-6.4	144	-0.03
41	P	Hexachlorocyclopentadiene	0.440	0.429	2.5	133	-0.03
42	S	2,4,6-Tribromophenol	0.275	0.321	-16.7	161#	-0.02
43	C	2,4,6-Trichlorophenol	0.450	0.503	-11.8	156#	-0.03
44		2,4,5-Trichlorophenol	0.490	0.539	-10.0	151#	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.385	1.383	0.1	135	-0.03
46		1,1'-Biphenyl	1.553	1.507	3.0	132	-0.03
47		2-Chloronaphthalene	1.215	1.221	-0.5	137	-0.02
48		2-Nitroaniline	0.433	0.470	-8.5	152#	-0.02
49		Acenaphthylene	2.062	1.964	4.8	128	-0.02
50		Dimethylphthalate	1.673	1.670	0.2	135	-0.02
51		2,6-Dinitrotoluene	0.327	0.368	-12.5	154#	-0.02
52	C	Acenaphthene	1.201	1.124	6.4	130	-0.02
53		3-Nitroaniline	0.335	0.369	-10.1	151#	-0.02
54	P	2,4-Dinitrophenol	0.161	0.229	-42.2#	201#	-0.02
55		Dibenzofuran	1.967	1.949	0.9	136	-0.02
56	P	4-Nitrophenol	0.290	0.301	-3.8	145	-0.02
57		2,4-Dinitrotoluene	0.444	0.540	-21.6	169#	-0.02
58		Fluorene	1.590	1.557	2.1	134	-0.02
59		2,3,4,6-Tetrachlorophenol	0.494	0.560	-13.4	154#	-0.02
60		Diethylphthalate	1.766	1.681	4.8	130	-0.03
61		4-Chlorophenyl-phenylether	0.925	0.972	-5.1	144	-0.03
62		4-Nitroaniline	0.374	0.401	-7.2	147	-0.02
63		Azobenzene	1.818	1.593	12.4	119	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	144	-0.03
65		4,6-Dinitro-2-methylphenol	0.092	0.126	-37.0#	211#	-0.02
66	c	n-Nitrosodiphenylamine	0.588	0.541	8.0	134	-0.02
67		4-Bromophenyl-phenylether	0.249	0.253	-1.6	149	-0.03
68		Hexachlorobenzene	0.258	0.265	-2.7	151#	-0.02
69		Atrazine	0.233	0.199	14.6	121	-0.03
70	C	Pentachlorophenol	0.151	0.165	-9.3	158#	-0.02
71		Phenanthrene	1.077	1.015	5.8	136	-0.02
72		Anthracene	1.083	1.013	6.5	134	-0.03
73		Carbazole	0.987	0.919	6.9	134	-0.02
74		Di-n-butylphthalate	1.191	1.073	9.9	130	-0.04
75	C	Fluoranthene	1.342	1.315	2.0	141	-0.03
76	I	Chrysene-d12	1.000	1.000	0.0	135	-0.04
77		Benzidine	0.548	0.492	10.2	115	-0.02
78		Pyrene	1.254	1.312	-4.6	139	-0.02
79	S	Terphenyl-d14	0.903	0.937	-3.8	137	-0.03
80		Butylbenzylphthalate	0.489	0.498	-1.8	138	-0.04
81		Benzo(a)anthracene	1.264	1.332	-5.4	141	-0.03
82		3,3'-Dichlorobenzidine	0.451	0.486	-7.8	143	-0.04
83		Chrysene	1.202	1.261	-4.9	140	-0.03
84		Bis(2-ethylhexyl)phthalate	0.705	0.687	2.6	131	-0.05
85	c	Di-n-octyl phthalate	1.188	1.155	2.8	131	-0.08
86		Indeno(1,2,3-cd)pyrene	1.453	1.583	-8.9	148	-0.10
87	I	Perylene-d12	1.000	1.000	0.0	149	-0.07

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.222	1.191	2.5	145	-0.06
89	Benzo(k)fluoranthene	1.141	1.131	0.9	148	-0.05
90 C	Benzo(a)pyrene	1.157	1.141	1.4	146	-0.07
91	Dibenzo(a,h)anthracene	1.171	1.155	1.4	146	-0.11
92	Benzo(g,h,i)perylene	1.165	1.152	1.1	146	-0.12

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