Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040319\
Data File : BG040345.D
Acq On : 4 Apr 2019 4:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 04 05:11:04 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 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	130	-0.03
2		1,4-Dioxane	0.566	0.599	-5.8	136	-0.03
3		Pyridine	1.523	1.654	-8.6	131	-0.03
4		n-Nitrosodimethylamine	0.705	0.744	-5.5	135	-0.02
5	S	2-Fluorophenol	1.203	1.300	-8.1	136	-0.02
6		Aniline	2.160	2.192	-1.5	127	-0.02
7	S	Phenol-d6	1.663	1.751	-5.3	133	-0.02
8		2-Chlorophenol	1.408	1.481	-5.2	132	-0.02
9		Benzaldehyde	1.140	1.090	4.4	116	-0.02
10	C	Phenol	1.805	1.914	-6.0	134	-0.03
11		bis(2-Chloroethyl)ether	1.445	1.455	-0.7	126	-0.02
12		1,3-Dichlorobenzene	1.531	1.609	-5.1	132	-0.03
13	C	1,4-Dichlorobenzene	1.525	1.618	-6.1	134	-0.02
14		1,2-Dichlorobenzene	1.458	1.518	-4.1	132	-0.02
15		Benzyl Alcohol	1.339	1.304	2.6	122	-0.02
16		2,2'-oxybis(1-Chloropropane	2.774	2.729	1.6	124	-0.01
17		2-Methylphenol	1.249	1.267	-1.4	127	-0.02
18		Hexachloroethane	0.580	0.598	-3.1	131	-0.03
19	P	n-Nitroso-di-n-propylamine	1.192	1.157	2.9	124	-0.03
20		3+4-Methylphenols	1.692	1.733	-2.4	128	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	125	-0.02
22		Acetophenone	0.499	0.488	2.2	123	-0.02
23	S	Nitrobenzene-d5	0.376	0.381	-1.3	125	-0.02
24		Nitrobenzene	0.390	0.394	-1.0	125	-0.03
25		Isophorone	0.774	0.762	1.6	123	-0.03
26	C	2-Nitrophenol	0.185	0.195	-5.4	131	-0.02
27		2,4-Dimethylphenol	0.293	0.283	3.4	120	-0.03
28		bis(2-Chloroethoxy)methane	0.458	0.462	-0.9	123	-0.03
29	C	2,4-Dichlorophenol	0.318	0.337	-6.0	129	-0.03
30		1,2,4-Trichlorobenzene	0.350	0.360	-2.9	129	-0.03
31		Naphthalene	1.025	1.046	-2.0	125	-0.03
32		Benzoic acid	0.177	0.172	2.8	130	-0.02
33		4-Chloroaniline	0.437	0.455	-4.1	127	-0.02
34	C	Hexachlorobutadiene	0.224	0.226	-0.9	126	-0.03
35		Caprolactam	0.126	0.138	-9.5	134	-0.01
36	C	4-Chloro-3-methylphenol	0.366	0.386	-5.5	130	-0.02
37		2-Methylnaphthalene	0.758	0.774	-2.1	127	-0.03
38		1-Methylnaphthalene	0.735	0.759	-3.3	128	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	130	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.636	0.642	-0.9	129	-0.02
41	P	Hexachlorocyclopentadiene	0.349	0.361	-3.4	136	-0.02
42	S	2,4,6-Tribromophenol	0.216	0.255	-18.1	151#	-0.02
43	C	2,4,6-Trichlorophenol	0.410	0.435	-6.1	136	-0.02
44		2,4,5-Trichlorophenol	0.447	0.476	-6.5	138	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.350	1.349	0.1	126	-0.02
46		1,1'-Biphenyl	1.580	1.593	-0.8	129	-0.02
47		2-Chloronaphthalene	1.212	1.218	-0.5	129	-0.03
48		2-Nitroaniline	0.409	0.426	-4.2	132	-0.03
49		Acenaphthylene	2.055	2.099	-2.1	129	-0.02
50		Dimethylphthalate	1.565	1.616	-3.3	132	-0.02
51		2,6-Dinitrotoluene	0.351	0.378	-7.7	139	-0.03
52	C	Acenaphthene	1.178	1.211	-2.8	132	-0.03
53		3-Nitroaniline	0.372	0.410	-10.2	141	-0.01
54	P	2,4-Dinitrophenol	0.187	0.291	-55.6#	212#	-0.02
55		Dibenzofuran	1.892	1.977	-4.5	134	-0.02
56	P	4-Nitrophenol	0.300	0.334	-11.3	144	-0.01
57		2,4-Dinitrotoluene	0.488	0.544	-11.5	142	-0.02
58		Fluorene	1.488	1.567	-5.3	135	-0.03
59		2,3,4,6-Tetrachlorophenol	0.405	0.471	-16.3	148	-0.02
60		Diethylphthalate	1.602	1.703	-6.3	136	-0.02
61		4-Chlorophenyl-phenylether	0.795	0.827	-4.0	132	-0.02
62		4-Nitroaniline	0.399	0.457	-14.5	147	-0.02
63		Azobenzene	1.511	1.573	-4.1	134	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	143	-0.02
65		4,6-Dinitro-2-methylphenol	0.112	0.127	-13.4	167#	-0.02
66	c	n-Nitrosodiphenylamine	0.585	0.585	0.0	139	-0.03
67		4-Bromophenyl-phenylether	0.225	0.227	-0.9	141	-0.03
68		Hexachlorobenzene	0.226	0.231	-2.2	143	-0.03
69		Atrazine	0.218	0.228	-4.6	145	-0.02
70	C	Pentachlorophenol	0.131	0.208	-58.8#	227#	-0.02
71		Phenanthrene	1.051	1.068	-1.6	141	-0.03
72		Anthracene	1.052	1.073	-2.0	141	-0.03
73		Carbazole	0.996	1.056	-6.0	147	-0.03
74		Di-n-butylphthalate	1.180	1.233	-4.5	144	-0.03
75	C	Fluoranthene	1.245	1.329	-6.7	147	-0.03
76	I	Chrysene-d12	1.000	1.000	0.0	157#	-0.03
77		Benzidine	0.722	0.937	-29.8#	196#	-0.02
78		Pyrene	1.315	1.264	3.9	146	-0.03
79	S	Terphenyl-d14	0.889	0.816	8.2	141	-0.03
80		Butylbenzylphthalate	0.563	0.567	-0.7	154#	-0.02
81		Benzo(a)anthracene	1.232	1.249	-1.4	155#	-0.03
82		3,3'-Dichlorobenzidine	0.469	0.494	-5.3	159#	-0.03
83		Chrysene	1.184	1.205	-1.8	156#	-0.03
84		Bis(2-ethylhexyl)phthalate	0.805	0.821	-2.0	158#	-0.03
85	c	Di-n-octyl phthalate	1.371	1.389	-1.3	154#	-0.04
86		Indeno(1,2,3-cd)pyrene	1.448	1.506	-4.0	161#	-0.07
87	I	Perylene-d12	1.000	1.000	0.0	152#	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.224	1.241	-1.4	151#	-0.05
89	Benzo(k)fluoranthene	1.126	1.160	-3.0	156#	-0.04
90 C	Benzo(a)pyrene	1.149	1.194	-3.9	155#	-0.05
91	Dibenzo(a,h)anthracene	1.067	1.147	-7.5	162#	-0.07
92	Benzo(g,h,i)perylene	1.097	1.177	-7.3	162#	-0.10

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