Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG010219\
Data File : BG038894.D
Acq On : 2 Jan 2019 16:20
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 03 01:13:39 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	-0.01
2		1,4-Dioxane	0.525	0.452	13.9	83	0.00
3		Pyridine	1.445	1.239	14.3	70	0.00
4		n-Nitrosodimethylamine	0.708	0.592	16.4	75	0.00
5	S	2-Fluorophenol	1.128	1.126	0.2	98	0.00
6		Aniline	1.976	1.964	0.6	96	-0.01
7	S	Phenol-d6	1.548	1.575	-1.7	100	0.00
8		2-Chlorophenol	1.305	1.309	-0.3	98	0.00
9		Benzaldehyde	1.004	0.898	10.6	95	-0.01
10	C	Phenol	1.645	1.658	-0.8	99	0.00
11		bis(2-Chloroethyl)ether	1.309	1.187	9.3	91	-0.01
12		1,3-Dichlorobenzene	1.543	1.516	1.7	97	-0.01
13	C	1,4-Dichlorobenzene	1.580	1.544	2.3	98	-0.01
14		1,2-Dichlorobenzene	1.490	1.501	-0.7	102	-0.02
15		Benzyl Alcohol	1.208	1.281	-6.0	99	0.00
16		2,2'-oxybis(1-Chloropropane	2.999	2.430	19.0	80	-0.01
17		2-Methylphenol	1.102	1.153	-4.6	100	0.00
18		Hexachloroethane	0.577	0.532	7.8	92	-0.02
19	P	n-Nitroso-di-n-propylamine	1.087	1.068	1.7	96	-0.01
20		3+4-Methylphenols	1.522	1.613	-6.0	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	-0.01
22		Acetophenone	0.500	0.488	2.4	100	-0.01
23	S	Nitrobenzene-d5	0.391	0.363	7.2	95	-0.01
24		Nitrobenzene	0.396	0.363	8.3	94	-0.01
25		Isophorone	0.703	0.636	9.5	80	-0.01
26	C	2-Nitrophenol	0.203	0.206	-1.5	105	-0.01
27		2,4-Dimethylphenol	0.291	0.277	4.8	98	0.00
28		bis(2-Chloroethoxy)methane	0.397	0.373	6.0	95	-0.02
29	C	2,4-Dichlorophenol	0.323	0.370	-14.6	113	0.00
30		1,2,4-Trichlorobenzene	0.390	0.405	-3.8	107	-0.01
31		Naphthalene	0.985	0.976	0.9	102	-0.02
32		Benzoic acid	0.154	0.041	73.4#	25#	-0.06
33		4-Chloroaniline	0.428	0.448	-4.7	105	-0.01
34	C	Hexachlorobutadiene	0.264	0.276	-4.5	106	-0.01
35		Caprolactam	0.134	0.137	-2.2	110	0.00
36	C	4-Chloro-3-methylphenol	0.369	0.375	-1.6	106	0.00
37		2-Methylnaphthalene	0.703	0.716	-1.8	105	-0.01
38		1-Methylnaphthalene	0.682	0.698	-2.3	106	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.748	0.728	2.7	110	-0.01
41	P	Hexachlorocyclopentadiene	0.411	0.539	-31.1#	146	-0.02
42	S	2,4,6-Tribromophenol	0.276	0.326	-18.1	132	-0.01
43	C	2,4,6-Trichlorophenol	0.460	0.523	-13.7	122	0.00
44		2,4,5-Trichlorophenol	0.487	0.533	-9.4	119	0.00

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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)
45	S	2-Fluorobiphenyl	1.458	1.416	2.9	111	-0.01
46		1,1'-Biphenyl	1.677	1.602	4.5	107	-0.01
47		2-Chloronaphthalene	1.295	1.241	4.2	109	-0.01
48		2-Nitroaniline	0.413	0.374	9.4	100	0.00
49		Acenaphthylene	1.936	1.878	3.0	109	-0.01
50		Dimethylphthalate	1.633	1.591	2.6	109	-0.01
51		2,6-Dinitrotoluene	0.370	0.358	3.2	109	-0.01
52	C	Acenaphthene	1.158	1.120	3.3	110	-0.01
53		3-Nitroaniline	0.377	0.370	1.9	109	0.00
54	P	2,4-Dinitrophenol	0.193	0.259	-34.2#	147	0.00
55		Dibenzofuran	1.985	1.937	2.4	112	0.00
56	P	4-Nitrophenol	0.286	0.253	11.5	95	0.00
57		2,4-Dinitrotoluene	0.521	0.501	3.8	106	0.00
58		Fluorene	1.470	1.441	2.0	110	-0.01
59		2,3,4,6-Tetrachlorophenol	0.438	0.489	-11.6	121	0.00
60		Diethylphthalate	1.793	1.682	6.2	107	0.00
61		4-Chlorophenyl-phenylether	0.917	0.927	-1.1	114	-0.01
62		4-Nitroaniline	0.388	0.376	3.1	104	0.00
63		Azobenzene	1.498	1.327	11.4	100	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
65		4,6-Dinitro-2-methylphenol	0.143	0.119	16.8	93	0.00
66	c	n-Nitrosodiphenylamine	0.588	0.585	0.5	113	-0.01
67		4-Bromophenyl-phenylether	0.244	0.251	-2.9	116	-0.01
68		Hexachlorobenzene	0.253	0.261	-3.2	115	0.00
69		Atrazine	0.218	0.207	5.0	105	-0.01
70	C	Pentachlorophenol	0.150	0.166	-10.7	122	0.00
71		Phenanthrene	1.037	1.009	2.7	111	-0.01
72		Anthracene	1.024	1.011	1.3	111	0.00
73		Carbazole	1.013	0.992	2.1	111	0.00
74		Di-n-butylphthalate	1.162	1.109	4.6	107	0.00
75	C	Fluoranthene	1.283	1.189	7.3	108	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	111	-0.01
77		Benzidine	0.591	0.500	15.4	78	-0.01
78		Pyrene	1.321	1.257	4.8	108	0.00
79	S	Terphenyl-d14	0.919	0.902	1.8	109	-0.01
80		Butylbenzylphthalate	0.516	0.482	6.6	101	-0.01
81		Benzo(a)anthracene	1.219	1.186	2.7	106	0.00
82		3,3'-Dichlorobenzidine	0.483	0.481	0.4	105	0.00
83		Chrysene	1.174	1.121	4.5	105	-0.01
84		Bis(2-ethylhexyl)phthalate	0.732	0.686	6.3	100	-0.01
85	c	Di-n-octyl phthalate	1.226	1.108	9.6	96	-0.02
86		Indeno(1,2,3-cd)pyrene	1.380	1.302	5.7	101	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	107	-0.02

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.098	1.071	2.5	102	-0.01
89	Benzo(k)fluoranthene	1.093	1.048	4.1	100	-0.01
90 C	Benzo(a)pyrene	1.059	1.034	2.4	102	-0.02
91	Dibenzo(a,h)anthracene	1.055	1.027	2.7	101	-0.02
92	Benzo(g,h,i)perylene	1.039	1.010	2.8	101	-0.02

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