Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021720\
Data File : BG044450.D
Acq On : 17 Feb 2020 17:01
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 18 06:04:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 10 16:11:50 2020
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	72	-0.01
2		1,4-Dioxane	0.509	0.539	-5.9	73	0.00
3		Pyridine	1.357	1.523	-12.2	72	0.00
4		n-Nitrosodimethylamine	0.567	0.562	0.9	70	0.00
5	S	2-Fluorophenol	1.133	1.214	-7.1	74	-0.01
6		Aniline	1.889	1.940	-2.7	70	-0.02
7	S	Phenol-d6	1.522	1.592	-4.6	72	-0.01
8		2-Chlorophenol	1.245	1.290	-3.6	71	-0.01
9		Benzaldehyde	0.867	0.846	2.4	71	-0.01
10	C	Phenol	1.506	1.564	-3.9	72	0.00
11		bis(2-Chloroethyl)ether	1.288	1.303	-1.2	70	-0.01
12		1,3-Dichlorobenzene	1.487	1.553	-4.4	74	-0.01
13	C	1,4-Dichlorobenzene	1.473	1.556	-5.6	73	-0.01
14		1,2-Dichlorobenzene	1.392	1.456	-4.6	72	-0.01
15		Benzyl Alcohol	1.077	1.094	-1.6	70	-0.01
16		2,2'-oxybis(1-Chloropropane	1.743	1.781	-2.2	71	-0.01
17		2-Methylphenol	1.075	1.097	-2.0	71	-0.01
18		Hexachloroethane	0.553	0.562	-1.6	72	-0.01
19	P	n-Nitroso-di-n-propylamine	1.014	1.012	0.2	69	-0.01
20		3+4-Methylphenols	1.481	1.510	-2.0	70	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	67	-0.01
22		Acetophenone	0.487	0.506	-3.9	70	-0.01
23	S	Nitrobenzene-d5	0.354	0.370	-4.5	70	-0.01
24		Nitrobenzene	0.346	0.359	-3.8	70	-0.01
25		Isophorone	0.660	0.682	-3.3	69	-0.02
26	C	2-Nitrophenol	0.179	0.193	-7.8	71	-0.01
27		2,4-Dimethylphenol	0.253	0.268	-5.9	70	-0.01
28		bis(2-Chloroethoxy)methane	0.440	0.452	-2.7	68	-0.02
29	C	2,4-Dichlorophenol	0.306	0.322	-5.2	70	-0.01
30		1,2,4-Trichlorobenzene	0.351	0.370	-5.4	71	-0.02
31		Naphthalene	0.970	1.019	-5.1	71	-0.01
32		Benzoic acid	0.135	0.110	18.5	56	-0.02
33		4-Chloroaniline	0.441	0.456	-3.4	70	-0.01
34	C	Hexachlorobutadiene	0.216	0.229	-6.0	71	-0.01
35		Caprolactam	0.112	0.115	-2.7	70	-0.02
36	C	4-Chloro-3-methylphenol	0.321	0.327	-1.9	69	-0.01
37		2-Methylnaphthalene	0.720	0.747	-3.8	70	-0.01
38		1-Methylnaphthalene	0.679	0.708	-4.3	70	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	66	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.598	0.648	-8.4	71	-0.02
41	P	Hexachlorocyclopentadiene	0.287	0.236	17.8	52	-0.01
42	S	2,4,6-Tribromophenol	0.235	0.248	-5.5	70	-0.01
43	C	2,4,6-Trichlorophenol	0.387	0.417	-7.8	70	-0.01
44		2,4,5-Trichlorophenol	0.395	0.426	-7.8	70	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.194	1.324	-10.9	72	-0.01
46		1,1'-Biphenyl	1.443	1.555	-7.8	71	-0.01
47		2-Chloronaphthalene	1.148	1.228	-7.0	71	-0.01
48		2-Nitroaniline	0.339	0.350	-3.2	68	-0.01
49		Acenaphthylene	1.684	1.816	-7.8	71	-0.01
50		Dimethylphthalate	1.438	1.512	-5.1	70	-0.01
51		2,6-Dinitrotoluene	0.321	0.332	-3.4	69	-0.02
52	C	Acenaphthene	1.091	1.151	-5.5	70	-0.01
53		3-Nitroaniline	0.355	0.364	-2.5	68	-0.01
54	P	2,4-Dinitrophenol	0.159	0.156	1.9	62	-0.01
55		Dibenzofuran	1.652	1.767	-7.0	70	-0.01
56	P	4-Nitrophenol	0.260	0.260	0.0	65	-0.01
57		2,4-Dinitrotoluene	0.449	0.468	-4.2	70	-0.01
58		Fluorene	1.301	1.398	-7.5	71	-0.01
59		2,3,4,6-Tetrachlorophenol	0.357	0.370	-3.6	68	-0.01
60		Diethylphthalate	1.433	1.490	-4.0	69	-0.01
61		4-Chlorophenyl-phenylether	0.706	0.745	-5.5	70	-0.01
62		4-Nitroaniline	0.354	0.368	-4.0	70	-0.01
63		Azobenzene	1.256	1.289	-2.6	68	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	68	-0.01
65		4,6-Dinitro-2-methylphenol	0.110	0.116	-5.5	70	-0.01
66	c	n-Nitrosodiphenylamine	0.560	0.584	-4.3	70	-0.02
67		4-Bromophenyl-phenylether	0.218	0.227	-4.1	70	-0.01
68		Hexachlorobenzene	0.244	0.256	-4.9	71	-0.01
69		Atrazine	0.192	0.187	2.6	64	-0.02
70	C	Pentachlorophenol	0.113	0.111	1.8	63	-0.01
71		Phenanthrene	0.968	1.039	-7.3	73	-0.01
72		Anthracene	0.957	1.026	-7.2	73	-0.02
73		Carbazole	0.883	0.947	-7.2	72	-0.01
74		Di-n-butylphthalate	1.091	1.173	-7.5	71	-0.02
75	C	Fluoranthene	1.120	1.207	-7.8	72	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	73	-0.02
77		Benzidine	0.555	0.560	-0.9	65	-0.02
78		Pyrene	1.284	1.309	-1.9	73	-0.02
79	S	Terphenyl-d14	0.972	0.943	3.0	75	-0.01
80		Butylbenzylphthalate	0.585	0.592	-1.2	72	-0.02
81		Benzo(a)anthracene	1.233	1.273	-3.2	74	-0.01
82		3,3'-Dichlorobenzidine	0.478	0.498	-4.2	72	-0.02
83		Chrysene	1.191	1.225	-2.9	73	-0.01
84		Bis(2-ethylhexyl)phthalate	0.806	0.817	-1.4	72	-0.02
85	c	Di-n-octyl phthalate	1.394	1.452	-4.2	74	-0.02
86		Indeno(1,2,3-cd)pyrene	1.554	1.568	-0.9	74	-0.04
87	I	Perylene-d12	1.000	1.000	0.0	70	-0.02

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88	Benzo(b)fluoranthene	1.152	1.235	-7.2	75	-0.02
89	Benzo(k)fluoranthene	1.123	1.186	-5.6	72	-0.02
90 C	Benzo(a)pyrene	1.090	1.157	-6.1	73	-0.02
91	Dibenzo(a,h)anthracene	1.078	1.152	-6.9	75	-0.04
92	Benzo(g,h,i)perylene	1.080	1.142	-5.7	74	-0.04

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