Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG082319\
Data File : BG042435.D
Acq On : 23 Aug 2019 16:13
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 23 18:02:38 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	212#	0.00
2		1,4-Dioxane	0.412	0.465	-12.9	237#	0.00
3		Pyridine	1.057	1.284	-21.5	248#	0.00
4		n-Nitrosodimethylamine	0.377	0.382	-1.3	218#	0.00
5	S	2-Fluorophenol	0.988	1.083	-9.6	230#	0.00
6		Aniline	1.779	1.954	-9.8	232#	0.00
7	S	Phenol-d6	1.334	1.439	-7.9	225#	0.00
8		2-Chlorophenol	1.197	1.259	-5.2	220#	0.00
9		Benzaldehyde	0.668	0.689	-3.1	260#	0.00
10	C	Phenol	1.356	1.479	-9.1	228#	0.00
11		bis(2-Chloroethyl)ether	1.065	1.186	-11.4	233#	0.00
12		1,3-Dichlorobenzene	1.522	1.548	-1.7	216#	0.00
13	C	1,4-Dichlorobenzene	1.542	1.526	1.0	208#	0.00
14		1,2-Dichlorobenzene	1.477	1.466	0.7	211#	0.00
15		Benzyl Alcohol	0.960	0.943	1.8	207#	0.00
16		2,2'-oxybis(1-Chloropropane	1.444	1.592	-10.2	233#	0.00
17		2-Methylphenol	0.999	1.057	-5.8	226#	0.00
18		Hexachloroethane	0.528	0.548	-3.8	220#	0.00
19	P	n-Nitroso-di-n-propylamine	0.864	0.889	-2.9	218#	0.00
20		3+4-Methylphenols	1.382	1.484	-7.4	229#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	215#	0.00
22		Acetophenone	0.428	0.438	-2.3	218#	0.00
23	S	Nitrobenzene-d5	0.289	0.295	-2.1	215#	0.00
24		Nitrobenzene	0.293	0.301	-2.7	219#	0.00
25		Isophorone	0.571	0.595	-4.2	219#	0.00
26	C	2-Nitrophenol	0.184	0.187	-1.6	219#	0.00
27		2,4-Dimethylphenol	0.253	0.262	-3.6	217#	0.00
28		bis(2-Chloroethoxy)methane	0.360	0.395	-9.7	232#	0.00
29	C	2,4-Dichlorophenol	0.331	0.322	2.7	204#	0.00
30		1,2,4-Trichlorobenzene	0.406	0.375	7.6	197#	0.00
31		Naphthalene	0.929	0.946	-1.8	214#	0.00
32		Benzoic acid	0.169	0.167	1.2	220#	0.00
33		4-Chloroaniline	0.429	0.441	-2.8	214#	0.00
34	C	Hexachlorobutadiene	0.273	0.227	16.8	179#	0.00
35		Caprolactam	0.110	0.115	-4.5	212#	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.309	1.6	205#	0.00
37		2-Methylnaphthalene	0.746	0.740	0.8	206#	0.00
38		1-Methylnaphthalene	0.708	0.694	2.0	205#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	194#	0.00
40		1,2,4,5-Tetrachlorobenzene	0.642	0.577	10.1	172#	0.00
41	P	Hexachlorocyclopentadiene	0.337	0.281	16.6	166#	0.00
42	S	2,4,6-Tribromophenol	0.284	0.227	20.1	149	0.00
43	C	2,4,6-Trichlorophenol	0.434	0.411	5.3	181#	0.00
44		2,4,5-Trichlorophenol	0.450	0.423	6.0	181#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.183	1.148	3.0	182#	0.00
46		1,1'-Biphenyl	1.293	1.342	-3.8	198#	0.00
47		2-Chloronaphthalene	1.115	1.134	-1.7	194#	0.00
48		2-Nitroaniline	0.269	0.294	-9.3	207#	0.00
49		Acenaphthylene	1.677	1.715	-2.3	191#	0.00
50		Dimethylphthalate	1.443	1.421	1.5	182#	0.00
51		2,6-Dinitrotoluene	0.334	0.336	-0.6	187#	0.00
52	C	Acenaphthene	1.018	1.048	-2.9	193#	0.00
53		3-Nitroaniline	0.335	0.360	-7.5	200#	0.00
54	P	2,4-Dinitrophenol	0.198	0.174	12.1	166#	0.00
55		Dibenzofuran	1.686	1.651	2.1	181#	0.00
56	P	4-Nitrophenol	0.238	0.259	-8.8	202#	0.00
57		2,4-Dinitrotoluene	0.479	0.473	1.3	182#	0.00
58		Fluorene	1.378	1.341	2.7	180#	0.00
59		2,3,4,6-Tetrachlorophenol	0.445	0.399	10.3	169#	0.00
60		Diethylphthalate	1.411	1.371	2.8	178#	0.00
61		4-Chlorophenyl-phenylether	0.831	0.751	9.6	168#	0.00
62		4-Nitroaniline	0.358	0.379	-5.9	193#	0.00
63		Azobenzene	0.967	1.036	-7.1	201#	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	171#	0.00
65		4,6-Dinitro-2-methylphenol	0.125	0.122	2.4	160#	0.00
66	c	n-Nitrosodiphenylamine	0.550	0.595	-8.2	179#	0.00
67		4-Bromophenyl-phenylether	0.254	0.239	5.9	158#	0.00
68		Hexachlorobenzene	0.276	0.252	8.7	152#	0.00
69		Atrazine	0.195	0.176	9.7	156#	0.00
70	C	Pentachlorophenol	0.143	0.129	9.8	148	0.00
71		Phenanthrene	0.993	1.012	-1.9	164#	0.00
72		Anthracene	0.992	1.019	-2.7	165#	0.00
73		Carbazole	0.884	0.953	-7.8	174#	0.00
74		Di-n-butylphthalate	1.045	1.109	-6.1	168#	0.00
75	C	Fluoranthene	1.212	0.995	17.9	130	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	115	0.00
77		Benzidine	0.573	0.659	-15.0	151#	0.00
78		Pyrene	1.226	1.429	-16.6	127	0.00
79	S	Terphenyl-d14	0.925	0.972	-5.1	117	0.00
80		Butylbenzylphthalate	0.482	0.553	-14.7	127	0.00
81		Benzo(a)anthracene	1.217	1.269	-4.3	115	0.00
82		3,3'-Dichlorobenzidine	0.489	0.490	-0.2	113	0.00
83		Chrysene	1.173	1.236	-5.4	115	0.00
84		Bis(2-ethylhexyl)phthalate	0.670	0.746	-11.3	123	0.00
85	c	Di-n-octyl phthalate	1.152	1.254	-8.9	120	0.00
86		Indeno(1,2,3-cd)pyrene	1.595	1.586	0.6	112	0.00
87	I	Perylene-d12	1.000	1.000	0.0	113	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.100	1.123	-2.1	109	0.00
89	Benzo(k)fluoranthene	1.057	1.093	-3.4	113	0.00
90 C	Benzo(a)pyrene	1.043	1.086	-4.1	112	0.00
91	Dibenzo(a,h)anthracene	1.056	1.079	-2.2	112	0.00
92	Benzo(g,h,i)perylene	1.057	1.068	-1.0	112	-0.01

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