Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012120\
Data File : BG044151.D
Acq On : 22 Jan 2020 9:29
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 22 11:09:10 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 31 13:32:23 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	61	-0.04
2		1,4-Dioxane	0.475	0.497	-4.6	64	-0.04
3		Pyridine	1.403	1.457	-3.8	62	-0.04
4		n-Nitrosodimethylamine	0.625	0.596	4.6	59	-0.04
5	S	2-Fluorophenol	1.089	1.181	-8.4	64	-0.03
6		Aniline	2.000	2.012	-0.6	61	-0.04
7	S	Phenol-d6	1.523	1.662	-9.1	66	-0.02
8		2-Chlorophenol	1.279	1.325	-3.6	63	-0.04
9		Benzaldehyde	0.974	0.819	15.9	54	-0.03
10	C	Phenol	1.569	1.667	-6.2	65	-0.03
11		bis(2-Chloroethyl)ether	1.354	1.357	-0.2	61	-0.04
12		1,3-Dichlorobenzene	1.496	1.553	-3.8	63	-0.04
13	C	1,4-Dichlorobenzene	1.476	1.536	-4.1	63	-0.04
14		1,2-Dichlorobenzene	1.417	1.479	-4.4	64	-0.04
15		Benzyl Alcohol	1.202	1.194	0.7	61	-0.04
16		2,2'-oxybis(1-Chloropropane	2.095	1.958	6.5	58	-0.04
17		2-Methylphenol	1.135	1.178	-3.8	64	-0.03
18		Hexachloroethane	0.575	0.587	-2.1	63	-0.04
19	P	n-Nitroso-di-n-propylamine	1.134	1.113	1.9	60	-0.04
20		3+4-Methylphenols	1.584	1.636	-3.3	63	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	62	-0.04
22		Acetophenone	0.497	0.495	0.4	61	-0.04
23	S	Nitrobenzene-d5	0.374	0.361	3.5	62	-0.04
24		Nitrobenzene	0.370	0.361	2.4	61	-0.04
25		Isophorone	0.701	0.685	2.3	60	-0.04
26	C	2-Nitrophenol	0.175	0.188	-7.4	68	-0.04
27		2,4-Dimethylphenol	0.264	0.267	-1.1	64	-0.03
28		bis(2-Chloroethoxy)methane	0.468	0.458	2.1	61	-0.04
29	C	2,4-Dichlorophenol	0.315	0.322	-2.2	64	-0.03
30		1,2,4-Trichlorobenzene	0.353	0.356	-0.8	63	-0.04
31		Naphthalene	0.968	1.021	-5.5	64	-0.04
32		Benzoic acid	0.197	0.167	15.2	59	-0.02
33		4-Chloroaniline	0.462	0.460	0.4	61	-0.03
34	C	Hexachlorobutadiene	0.215	0.216	-0.5	63	-0.04
35		Caprolactam	0.117	0.118	-0.9	63	-0.03
36	C	4-Chloro-3-methylphenol	0.335	0.348	-3.9	65	-0.02
37		2-Methylnaphthalene	0.729	0.768	-5.3	65	-0.04
38		1-Methylnaphthalene	0.691	0.718	-3.9	64	-0.03
39	I	Acenaphthene-d10	1.000	1.000	0.0	62	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.586	0.605	-3.2	65	-0.04
41	P	Hexachlorocyclopentadiene	0.304	0.282	7.2	58	-0.04
42	S	2,4,6-Tribromophenol	0.209	0.233	-11.5	68	-0.03
43	C	2,4,6-Trichlorophenol	0.403	0.412	-2.2	63	-0.03
44		2,4,5-Trichlorophenol	0.416	0.427	-2.6	64	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.104	1.245	-12.8	66	-0.04
46		1,1'-Biphenyl	1.434	1.534	-7.0	65	-0.03
47		2-Chloronaphthalene	1.159	1.213	-4.7	64	-0.03
48		2-Nitroaniline	0.373	0.376	-0.8	63	-0.02
49		Acenaphthylene	1.665	1.813	-8.9	66	-0.03
50		Dimethylphthalate	1.420	1.496	-5.4	64	-0.03
51		2,6-Dinitrotoluene	0.321	0.333	-3.7	65	-0.03
52	C	Acenaphthene	1.168	1.172	-0.3	62	-0.03
53		3-Nitroaniline	0.364	0.375	-3.0	63	-0.03
54	P	2,4-Dinitrophenol	0.145	0.169	-16.6	73	-0.02
55		Dibenzofuran	1.608	1.758	-9.3	66	-0.02
56	P	4-Nitrophenol	0.279	0.283	-1.4	61	-0.01
57		2,4-Dinitrotoluene	0.437	0.473	-8.2	66	-0.03
58		Fluorene	1.274	1.389	-9.0	66	-0.03
59		2,3,4,6-Tetrachlorophenol	0.358	0.375	-4.7	64	-0.03
60		Diethylphthalate	1.420	1.532	-7.9	65	-0.04
61		4-Chlorophenyl-phenylether	0.692	0.727	-5.1	65	-0.03
62		4-Nitroaniline	0.360	0.384	-6.7	65	-0.03
63		Azobenzene	1.317	1.388	-5.4	64	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	64	-0.03
65		4,6-Dinitro-2-methylphenol	0.097	0.114	-17.5	77	-0.02
66	c	n-Nitrosodiphenylamine	0.589	0.601	-2.0	66	-0.03
67		4-Bromophenyl-phenylether	0.225	0.224	0.4	66	-0.03
68		Hexachlorobenzene	0.242	0.246	-1.7	67	-0.03
69		Atrazine	0.191	0.190	0.5	61	-0.02
70	C	Pentachlorophenol	0.134	0.119	11.2	56	-0.02
71		Phenanthrene	0.959	1.045	-9.0	68	-0.03
72		Anthracene	0.948	1.036	-9.3	68	-0.03
73		Carbazole	0.876	0.949	-8.3	67	-0.02
74		Di-n-butylphthalate	1.118	1.191	-6.5	67	-0.04
75	C	Fluoranthene	1.097	1.185	-8.0	69	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	64	-0.03
77		Benzidine	0.503	0.498	1.0	57	-0.02
78		Pyrene	1.305	1.401	-7.4	70	-0.03
79	S	Terphenyl-d14	0.828	0.953	-15.1	73	-0.03
80		Butylbenzylphthalate	0.622	0.668	-7.4	69	-0.03
81		Benzo(a)anthracene	1.240	1.331	-7.3	68	-0.04
82		3,3'-Dichlorobenzidine	0.490	0.504	-2.9	65	-0.04
83		Chrysene	1.178	1.287	-9.3	69	-0.04
84		Bis(2-ethylhexyl)phthalate	0.858	0.915	-6.6	68	-0.04
85	c	Di-n-octyl phthalate	1.442	1.594	-10.5	69	-0.05
86		Indeno(1,2,3-cd)pyrene	1.601	1.655	-3.4	68	-0.09
87	I	Perylene-d12	1.000	1.000	0.0	65	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.182	1.244	-5.2	69	-0.05
89	Benzo(k)fluoranthene	1.124	1.185	-5.4	67	-0.05
90 C	Benzo(a)pyrene	1.116	1.164	-4.3	67	-0.05
91	Dibenzo(a,h)anthracene	1.121	1.164	-3.8	68	-0.09
92	Benzo(g,h,i)perylene	1.113	1.161	-4.3	68	-0.09

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