Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092718\
Data File : BG037035.D
Acq On : 28 Sep 2018 4:59
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 28 07:38:28 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 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	116	0.00
2		1,4-Dioxane	0.477	0.579	-21.4	131	0.00
3		Pyridine	1.378	1.540	-11.8	122	0.00
4		n-Nitrosodimethylamine	0.612	0.731	-19.4	134	0.00
5	S	2-Fluorophenol	1.043	1.198	-14.9	128	0.00
6		Aniline	2.094	2.092	0.1	113	-0.01
7	S	Phenol-d6	1.613	1.693	-5.0	117	-0.01
8		2-Chlorophenol	1.211	1.274	-5.2	117	0.00
9		Benzaldehyde	1.066	1.067	-0.1	113	0.00
10	C	Phenol	1.665	1.783	-7.1	120	0.00
11		bis(2-Chloroethyl)ether	1.540	1.484	3.6	115	-0.01
12		1,3-Dichlorobenzene	1.485	1.516	-2.1	115	-0.01
13	C	1,4-Dichlorobenzene	1.519	1.558	-2.6	117	-0.01
14		1,2-Dichlorobenzene	1.472	1.468	0.3	116	0.00
15		Benzyl Alcohol	1.309	1.286	1.8	111	-0.01
16		2,2'-oxybis(1-Chloropropane	2.957	3.012	-1.9	115	0.00
17		2-Methylphenol	1.160	1.153	0.6	114	0.00
18		Hexachloroethane	0.559	0.593	-6.1	120	0.00
19	P	n-Nitroso-di-n-propylamine	1.342	1.271	5.3	108	0.00
20		3+4-Methylphenols	1.614	1.618	-0.2	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.470	0.481	-2.3	109	0.00
23	S	Nitrobenzene-d5	0.373	0.409	-9.7	113	0.00
24		Nitrobenzene	0.399	0.413	-3.5	108	0.00
25		Isophorone	0.682	0.720	-5.6	111	-0.01
26	C	2-Nitrophenol	0.159	0.181	-13.8	115	0.00
27		2,4-Dimethylphenol	0.248	0.248	0.0	103	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.427	-1.4	105	0.00
29	C	2,4-Dichlorophenol	0.287	0.316	-10.1	111	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.363	-1.1	105	0.00
31		Naphthalene	0.952	0.962	-1.1	107	-0.01
32		Benzoic acid	0.175	0.155	11.4	92	0.00
33		4-Chloroaniline	0.433	0.437	-0.9	106	-0.01
34	C	Hexachlorobutadiene	0.248	0.252	-1.6	106	-0.01
35		Caprolactam	0.118	0.128	-8.5	111	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.384	-12.0	111	-0.01
37		2-Methylnaphthalene	0.725	0.718	1.0	105	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	0.698	0.690	1.1	100	0.00
40	P	Hexachlorocyclopentadiene	0.359	0.327	8.9	88	-0.01
41	S	2,4,6-Tribromophenol	0.239	0.252	-5.4	104	-0.01
42	C	2,4,6-Trichlorophenol	0.387	0.418	-8.0	106	-0.01
43		2,4,5-Trichlorophenol	0.413	0.462	-11.9	108	0.00
44	S	2-Fluorobiphenyl	1.343	1.339	0.3	101	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.532	1.526	0.4	102	-0.01
46		2-Chloronaphthalene	1.205	1.216	-0.9	103	0.00
47		2-Nitroaniline	0.417	0.467	-12.0	109	0.00
48		Acenaphthylene	1.888	1.932	-2.3	104	0.00
49		Dimethylphthalate	1.610	1.604	0.4	101	0.00
50		2,6-Dinitrotoluene	0.351	0.364	-3.7	104	0.00
51	C	Acenaphthene	1.141	1.136	0.4	102	0.00
52		3-Nitroaniline	0.370	0.389	-5.1	104	0.00
53	P	2,4-Dinitrophenol	0.136	0.114	16.2	82	0.00
54		Dibenzofuran	1.930	1.918	0.6	101	0.00
55	P	4-Nitrophenol	0.236	0.260	-10.2	107	-0.01
56		2,4-Dinitrotoluene	0.490	0.520	-6.1	106	0.00
57		Fluorene	1.442	1.458	-1.1	103	-0.01
58		2,3,4,6-Tetrachlorophenol	0.419	0.456	-8.8	106	0.00
59		Diethylphthalate	1.624	1.645	-1.3	103	0.00
60		4-Chlorophenyl-phenylether	0.913	0.908	0.5	101	-0.01
61		4-Nitroaniline	0.389	0.406	-4.4	103	-0.01
62		Azobenzene	1.560	1.688	-8.2	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.117	-11.4	112	0.00
65	c	n-Nitrosodiphenylamine	0.552	0.569	-3.1	103	0.00
66		4-Bromophenyl-phenylether	0.227	0.224	1.3	98	0.00
67		Hexachlorobenzene	0.234	0.233	0.4	99	0.00
68		Atrazine	0.224	0.224	0.0	98	-0.01
69	C	Pentachlorophenol	0.148	0.151	-2.0	99	-0.01
70		Phenanthrene	0.964	0.981	-1.8	103	-0.01
71		Anthracene	0.953	0.989	-3.8	104	0.00
72		Carbazole	0.929	0.972	-4.6	105	0.00
73		Di-n-butylphthalate	1.032	1.104	-7.0	107	-0.01
74	C	Fluoranthene	1.160	1.189	-2.5	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	107	-0.01
76		Benzidine	0.605	0.561	7.3	100	0.00
77		Pyrene	1.200	1.178	1.8	104	-0.01
78	S	Terphenyl-d14	0.761	0.727	4.5	102	0.00
79		Butylbenzylphthalate	0.478	0.507	-6.1	113	-0.01
80		Benzo(a)anthracene	1.167	1.149	1.5	107	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.449	-1.1	105	0.00
82		Chrysene	1.128	1.093	3.1	105	-0.01
83		Bis(2-ethylhexyl)phthalate	0.668	0.697	-4.3	113	-0.01
84	c	Di-n-octyl phthalate	1.100	1.180	-7.3	114	-0.02
85		Indeno(1,2,3-cd)pyrene	1.211	1.267	-4.6	109	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	108	-0.02
87		Benzo(b)fluoranthene	1.066	1.028	3.6	102	-0.02

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88	Benzo(k)fluoranthene	1.069	1.088	-1.8	112	-0.02
89 C	Benzo(a)pyrene	1.030	1.026	0.4	107	-0.02
90	Dibenzo(a,h)anthracene	0.966	0.990	-2.5	110	-0.02
91	Benzo(g,h,i)perylene	0.969	0.993	-2.5	108	-0.02

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