Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG100518\
Data File : BG037175.D
Acq On : 5 Oct 2018 5:26
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 05 08:47:03 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	99	-0.01
2		1,4-Dioxane	0.477	0.578	-21.2	112	-0.01
3		Pyridine	1.378	1.586	-15.1	109	-0.02
4		n-Nitrosodimethylamine	0.612	0.715	-16.8	113	-0.01
5	S	2-Fluorophenol	1.043	1.142	-9.5	105	-0.02
6		Aniline	2.094	2.133	-1.9	99	-0.02
7	S	Phenol-d6	1.613	1.768	-9.6	105	-0.02
8		2-Chlorophenol	1.211	1.313	-8.4	104	-0.02
9		Benzaldehyde	1.066	1.084	-1.7	99	-0.02
10	C	Phenol	1.665	1.880	-12.9	109	-0.02
11		bis(2-Chloroethyl)ether	1.540	1.514	1.7	101	-0.02
12		1,3-Dichlorobenzene	1.485	1.501	-1.1	98	-0.02
13	C	1,4-Dichlorobenzene	1.519	1.521	-0.1	99	-0.02
14		1,2-Dichlorobenzene	1.472	1.461	0.7	99	-0.02
15		Benzyl Alcohol	1.309	1.325	-1.2	98	-0.02
16		2,2'-oxybis(1-Chloropropane	2.957	3.122	-5.6	102	-0.02
17		2-Methylphenol	1.160	1.225	-5.6	104	-0.02
18		Hexachloroethane	0.559	0.593	-6.1	103	-0.02
19	P	n-Nitroso-di-n-propylamine	1.342	1.358	-1.2	99	-0.02
20		3+4-Methylphenols	1.614	1.755	-8.7	105	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.02
22		Acetophenone	0.470	0.491	-4.5	103	-0.02
23	S	Nitrobenzene-d5	0.373	0.399	-7.0	102	-0.02
24		Nitrobenzene	0.399	0.408	-2.3	99	-0.01
25		Isophorone	0.682	0.725	-6.3	103	-0.02
26	C	2-Nitrophenol	0.159	0.178	-11.9	105	-0.01
27		2,4-Dimethylphenol	0.248	0.259	-4.4	100	-0.02
28		bis(2-Chloroethoxy)methane	0.421	0.429	-1.9	98	-0.02
29	C	2,4-Dichlorophenol	0.287	0.318	-10.8	103	-0.02
30		1,2,4-Trichlorobenzene	0.359	0.350	2.5	94	-0.02
31		Naphthalene	0.952	0.953	-0.1	98	-0.02
32		Benzoic acid	0.175	0.163	6.9	90	-0.01
33		4-Chloroaniline	0.433	0.437	-0.9	98	-0.02
34	C	Hexachlorobutadiene	0.248	0.239	3.6	93	-0.02
35		Caprolactam	0.118	0.125	-5.9	100	-0.01
36	C	4-Chloro-3-methylphenol	0.343	0.404	-17.8	109	-0.02
37		2-Methylnaphthalene	0.725	0.716	1.2	97	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.698	0.667	4.4	93	-0.02
40	P	Hexachlorocyclopentadiene	0.359	0.291	18.9	75	-0.02
41	S	2,4,6-Tribromophenol	0.239	0.239	0.0	95	-0.02
42	C	2,4,6-Trichlorophenol	0.387	0.419	-8.3	102	-0.02
43		2,4,5-Trichlorophenol	0.413	0.460	-11.4	103	-0.01
44	S	2-Fluorobiphenyl	1.343	1.327	1.2	96	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.532	1.525	0.5	98	-0.02
46		2-Chloronaphthalene	1.205	1.192	1.1	97	-0.02
47		2-Nitroaniline	0.417	0.464	-11.3	104	-0.02
48		Acenaphthylene	1.888	1.908	-1.1	99	-0.02
49		Dimethylphthalate	1.610	1.602	0.5	97	-0.02
50		2,6-Dinitrotoluene	0.351	0.359	-2.3	99	-0.01
51	C	Acenaphthene	1.141	1.130	1.0	97	-0.02
52		3-Nitroaniline	0.370	0.379	-2.4	97	-0.01
53	P	2,4-Dinitrophenol	0.136	0.143	-5.1	99	-0.01
54		Dibenzofuran	1.930	1.935	-0.3	97	-0.02
55	P	4-Nitrophenol	0.236	0.292	-23.7	115	-0.02
56		2,4-Dinitrotoluene	0.490	0.502	-2.4	98	-0.01
57		Fluorene	1.442	1.440	0.1	98	-0.02
58		2,3,4,6-Tetrachlorophenol	0.419	0.444	-6.0	99	-0.01
59		Diethylphthalate	1.624	1.642	-1.1	99	-0.01
60		4-Chlorophenyl-phenylether	0.913	0.874	4.3	93	-0.02
61		4-Nitroaniline	0.389	0.383	1.5	94	-0.02
62		Azobenzene	1.560	1.662	-6.5	104	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	97	-0.02
64		4,6-Dinitro-2-methylphenol	0.105	0.108	-2.9	99	-0.01
65	c	n-Nitrosodiphenylamine	0.552	0.566	-2.5	98	-0.01
66		4-Bromophenyl-phenylether	0.227	0.219	3.5	91	-0.02
67		Hexachlorobenzene	0.234	0.226	3.4	92	-0.02
68		Atrazine	0.224	0.213	4.9	89	-0.02
69	C	Pentachlorophenol	0.148	0.139	6.1	87	-0.02
70		Phenanthrene	0.964	0.962	0.2	96	-0.02
71		Anthracene	0.953	0.972	-2.0	98	-0.02
72		Carbazole	0.929	0.948	-2.0	97	-0.02
73		Di-n-butylphthalate	1.032	1.077	-4.4	99	-0.02
74	C	Fluoranthene	1.160	1.149	0.9	95	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	99	-0.02
76		Benzidine	0.605	0.728	-20.3	120	-0.02
77		Pyrene	1.200	1.163	3.1	95	-0.02
78	S	Terphenyl-d14	0.761	0.727	4.5	95	-0.02
79		Butylbenzylphthalate	0.478	0.503	-5.2	104	-0.02
80		Benzo(a)anthracene	1.167	1.132	3.0	98	-0.02
81		3,3'-Dichlorobenzidine	0.444	0.444	0.0	96	-0.02
82		Chrysene	1.128	1.103	2.2	98	-0.02
83		Bis(2-ethylhexyl)phthalate	0.668	0.697	-4.3	105	-0.02
84	c	Di-n-octyl phthalate	1.100	1.172	-6.5	105	-0.03
85		Indeno(1,2,3-cd)pyrene	1.211	1.229	-1.5	98	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	100	-0.03
87		Benzo(b)fluoranthene	1.066	1.022	4.1	94	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.069	1.081	-1.1	102	-0.03
89 C	Benzo(a)pyrene	1.030	1.028	0.2	99	-0.03
90	Dibenzo(a,h)anthracene	0.966	0.971	-0.5	99	-0.06
91	Benzo(g,h,i)perylene	0.969	0.979	-1.0	98	-0.06

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