Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100118\
Data File : BG037095.D
Acq On : 2 Oct 2018 4:02
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 02 08:58:00 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	123	0.00
2		1,4-Dioxane	0.477	0.552	-15.7	133	0.00
3		Pyridine	1.378	1.504	-9.1	127	0.00
4		n-Nitrosodimethylamine	0.612	0.652	-6.5	128	0.00
5	S	2-Fluorophenol	1.043	1.152	-10.5	131	-0.01
6		Aniline	2.094	2.025	3.3	117	-0.01
7	S	Phenol-d6	1.613	1.657	-2.7	122	-0.01
8		2-Chlorophenol	1.211	1.268	-4.7	124	0.00
9		Benzaldehyde	1.066	1.038	2.6	117	-0.01
10	C	Phenol	1.665	1.772	-6.4	127	-0.01
11		bis(2-Chloroethyl)ether	1.540	1.464	4.9	120	-0.01
12		1,3-Dichlorobenzene	1.485	1.505	-1.3	121	-0.01
13	C	1,4-Dichlorobenzene	1.519	1.509	0.7	121	-0.01
14		1,2-Dichlorobenzene	1.472	1.445	1.8	122	0.00
15		Benzyl Alcohol	1.309	1.209	7.6	111	-0.01
16		2,2'-oxybis(1-Chloropropane	2.957	2.879	2.6	117	-0.01
17		2-Methylphenol	1.160	1.124	3.1	118	-0.01
18		Hexachloroethane	0.559	0.590	-5.5	127	-0.01
19	P	n-Nitroso-di-n-propylamine	1.342	1.229	8.4	111	-0.01
20		3+4-Methylphenols	1.614	1.574	2.5	117	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	113	-0.01
22		Acetophenone	0.470	0.467	0.6	115	-0.01
23	S	Nitrobenzene-d5	0.373	0.392	-5.1	117	0.00
24		Nitrobenzene	0.399	0.406	-1.8	114	0.00
25		Isophorone	0.682	0.686	-0.6	114	-0.01
26	C	2-Nitrophenol	0.159	0.178	-11.9	123	0.00
27		2,4-Dimethylphenol	0.248	0.241	2.8	108	-0.01
28		bis(2-Chloroethoxy)methane	0.421	0.416	1.2	110	-0.01
29	C	2,4-Dichlorophenol	0.287	0.295	-2.8	112	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.359	0.0	112	-0.01
31		Naphthalene	0.952	0.943	0.9	113	-0.02
32		Benzoic acid	0.175	0.166	5.1	107	-0.01
33		4-Chloroaniline	0.433	0.416	3.9	109	-0.01
34	C	Hexachlorobutadiene	0.248	0.249	-0.4	113	-0.02
35		Caprolactam	0.118	0.119	-0.8	112	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.367	-7.0	115	-0.01
37		2-Methylnaphthalene	0.725	0.703	3.0	112	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.698	0.683	2.1	107	-0.01
40	P	Hexachlorocyclopentadiene	0.359	0.300	16.4	87	-0.02
41	S	2,4,6-Tribromophenol	0.239	0.251	-5.0	112	-0.01
42	C	2,4,6-Trichlorophenol	0.387	0.394	-1.8	108	-0.01
43		2,4,5-Trichlorophenol	0.413	0.434	-5.1	109	0.00
44	S	2-Fluorobiphenyl	1.343	1.317	1.9	106	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.532	1.515	1.1	109	-0.01
46		2-Chloronaphthalene	1.205	1.191	1.2	109	-0.01
47		2-Nitroaniline	0.417	0.444	-6.5	112	-0.01
48		Acenaphthylene	1.888	1.906	-1.0	110	-0.01
49		Dimethylphthalate	1.610	1.573	2.3	106	-0.01
50		2,6-Dinitrotoluene	0.351	0.366	-4.3	113	0.00
51	C	Acenaphthene	1.141	1.132	0.8	109	-0.01
52		3-Nitroaniline	0.370	0.390	-5.4	112	0.00
53	P	2,4-Dinitrophenol	0.136	0.113	16.9	87	0.00
54		Dibenzofuran	1.930	1.915	0.8	108	-0.01
55	P	4-Nitrophenol	0.236	0.245	-3.8	109	-0.01
56		2,4-Dinitrotoluene	0.490	0.514	-4.9	113	0.00
57		Fluorene	1.442	1.443	-0.1	110	-0.01
58		2,3,4,6-Tetrachlorophenol	0.419	0.439	-4.8	109	-0.01
59		Diethylphthalate	1.624	1.613	0.7	109	0.00
60		4-Chlorophenyl-phenylether	0.913	0.886	3.0	105	-0.01
61		4-Nitroaniline	0.389	0.403	-3.6	110	-0.01
62		Azobenzene	1.560	1.639	-5.1	115	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.103	1.9	107	0.00
65	c	n-Nitrosodiphenylamine	0.552	0.553	-0.2	109	0.00
66		4-Bromophenyl-phenylether	0.227	0.221	2.6	105	-0.01
67		Hexachlorobenzene	0.234	0.226	3.4	105	-0.01
68		Atrazine	0.224	0.217	3.1	104	-0.01
69	C	Pentachlorophenol	0.148	0.141	4.7	101	-0.01
70		Phenanthrene	0.964	0.964	0.0	110	-0.01
71		Anthracene	0.953	0.971	-1.9	112	-0.01
72		Carbazole	0.929	0.965	-3.9	113	-0.01
73		Di-n-butylphthalate	1.032	1.084	-5.0	115	-0.01
74	C	Fluoranthene	1.160	1.170	-0.9	111	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	115	-0.01
76		Benzidine	0.605	0.655	-8.3	125	-0.01
77		Pyrene	1.200	1.170	2.5	111	-0.01
78	S	Terphenyl-d14	0.761	0.713	6.3	108	-0.01
79		Butylbenzylphthalate	0.478	0.500	-4.6	120	-0.01
80		Benzo(a)anthracene	1.167	1.141	2.2	115	-0.02
81		3,3'-Dichlorobenzidine	0.444	0.435	2.0	110	-0.01
82		Chrysene	1.128	1.108	1.8	115	-0.01
83		Bis(2-ethylhexyl)phthalate	0.668	0.693	-3.7	122	-0.02
84	c	Di-n-octyl phthalate	1.100	1.164	-5.8	121	-0.02
85		Indeno(1,2,3-cd)pyrene	1.211	1.237	-2.1	115	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	117	-0.03
87		Benzo(b)fluoranthene	1.066	1.028	3.6	110	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.069	1.088	-1.8	121	-0.02
89 C	Benzo(a)pyrene	1.030	1.028	0.2	116	-0.02
90	Dibenzo(a,h)anthracene	0.966	0.969	-0.3	116	-0.03
91	Benzo(g,h,i)perylene	0.969	0.976	-0.7	115	-0.03

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