Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\
Data File : BG037207.D
Acq On : 6 Oct 2018 2:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 06 05:33:25 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	74	-0.01
2		1,4-Dioxane	0.477	0.001	99.8#	0#	-0.04
3		Pyridine	1.378	0.000	100.0#	0#	-3.93#
4		n-Nitrosodimethylamine	0.612	0.002	99.7#	0#	0.00
5	S	2-Fluorophenol	1.043	1.722	-65.1#	119	-0.02
6		Aniline	2.094	0.004	99.8#	0#	0.19
7	S	Phenol-d6	1.613	2.353	-45.9#	105	-0.02
8		2-Chlorophenol	1.211	0.000	100.0#	0#	-7.61#
9		Benzaldehyde	1.066	0.002	99.8#	0#	0.00
10	C	Phenol	1.665	0.008	99.5#	0#	0.33
11		bis(2-Chloroethyl)ether	1.540	0.004	99.7#	0#	0.09
12		1,3-Dichlorobenzene	1.485	0.000	100.0#	0#	-7.94#
13	C	1,4-Dichlorobenzene	1.519	0.000	100.0#	0#	-8.09#
14		1,2-Dichlorobenzene	1.472	0.000	100.0#	0#	-8.40#
15		Benzyl Alcohol	1.309	0.051	96.1#	3#	0.06
16		2,2'-oxybis(1-Chloropropane	2.957	0.000	100.0#	0#	-8.58#
17		2-Methylphenol	1.160	0.000	100.0#	0#	-8.49#
18		Hexachloroethane	0.559	0.000	100.0#	0#	-9.13#
19	P	n-Nitroso-di-n-propylamine	1.342	0.000#	100.0#	0#	-8.85#
20		3+4-Methylphenols	1.614	0.000	100.0#	0#	-8.82#
21	I	Naphthalene-d8	1.000	1.000	0.0	64	-0.02
22		Acetophenone	0.470	0.000	100.0#	0#	-8.87#
23	S	Nitrobenzene-d5	0.373	0.365	2.1	62	-0.02
24		Nitrobenzene	0.399	0.001	99.7#	0#	-0.05
25		Isophorone	0.682	0.002	99.7#	0#	-0.11
26	C	2-Nitrophenol	0.159	0.000	100.0#	0#	-9.96#
27		2,4-Dimethylphenol	0.248	0.000	100.0#	0#	-10.02#
28		bis(2-Chloroethoxy)methane	0.421	0.000	100.0#	0#	-10.26#
29	C	2,4-Dichlorophenol	0.287	0.000	100.0#	0#	-10.50#
30		1,2,4-Trichlorobenzene	0.359	0.000	100.0#	0#	-10.71#
31		Naphthalene	0.952	0.000	100.0#	0#	-10.91#
32		Benzoic acid	0.175	0.000	100.0#	0#	-10.16#
33		4-Chloroaniline	0.433	0.000	100.0#	0#	-11.01#
34	C	Hexachlorobutadiene	0.248	0.000	100.0#	0#	-11.19#
35		Caprolactam	0.118	0.000	100.0#	0#	-11.78#
36	C	4-Chloro-3-methylphenol	0.343	0.000	100.0#	0#	-12.13#
37		2-Methylnaphthalene	0.725	0.000	100.0#	0#	-12.51#
38	I	Acenaphthene-d10	1.000	1.000	0.0	64	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.698	0.000	100.0#	0#	-12.87#
40	P	Hexachlorocyclopentadiene	0.359	0.000#	100.0#	0#	-12.85#
41	S	2,4,6-Tribromophenol	0.239	0.347	-45.2#	90	-0.02
42	C	2,4,6-Trichlorophenol	0.387	0.000	100.0#	0#	-13.11#
43		2,4,5-Trichlorophenol	0.413	0.000	100.0#	0#	-13.18#
44	S	2-Fluorobiphenyl	1.343	1.234	8.1	58	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.532	0.001	99.9#	0#	0.00
46		2-Chloronaphthalene	1.205	0.000	100.0#	0#	-13.55#
47		2-Nitroaniline	0.417	0.005	98.8#	1#	-0.49
48		Acenaphthylene	1.888	0.000	100.0#	0#	-14.40#
49		Dimethylphthalate	1.610	0.000	100.0#	0#	-14.13#
50		2,6-Dinitrotoluene	0.351	0.000	100.0#	0#	-14.25#
51	C	Acenaphthene	1.141	0.001	99.9#	0#	-0.08
52		3-Nitroaniline	0.370	0.000	100.0#	0#	-14.58#
53	P	2,4-Dinitrophenol	0.136	0.000#	100.0#	0#	-14.79#
54		Dibenzofuran	1.930	0.000	100.0#	0#	-15.07#
55	P	4-Nitrophenol	0.236	0.000#	100.0#	0#	-14.89#
56		2,4-Dinitrotoluene	0.490	0.000	100.0#	0#	-15.03#
57		Fluorene	1.442	0.026	98.2#	1#	0.42
58		2,3,4,6-Tetrachlorophenol	0.419	0.000	100.0#	0#	-15.30#
59		Diethylphthalate	1.624	0.000	100.0#	0#	-15.49#
60		4-Chlorophenyl-phenylether	0.913	0.000	100.0#	0#	-15.71#
61		4-Nitroaniline	0.389	0.000	100.0#	0#	-15.75#
62		Azobenzene	1.560	0.004	99.7#	0#	0.15
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	-0.02
64		4,6-Dinitro-2-methylphenol	0.105	0.000	100.0#	0#	0.35
65	c	n-Nitrosodiphenylamine	0.552	0.010	98.2#	1#	0.22
66		4-Bromophenyl-phenylether	0.227	0.000	100.0#	0#	-16.61#
67		Hexachlorobenzene	0.234	0.000	100.0#	0#	-16.72#
68		Atrazine	0.224	0.000	100.0#	0#	-16.88#
69	C	Pentachlorophenol	0.148	0.000	100.0#	0#	-17.07#
70		Phenanthrene	0.964	0.000	100.0#	0#	-17.47#
71		Anthracene	0.953	0.000	100.0#	0#	-17.55#
72		Carbazole	0.929	0.000	100.0#	0#	-17.82#
73		Di-n-butylphthalate	1.032	0.000	100.0#	0#	0.00
74	C	Fluoranthene	1.160	0.000	100.0#	0#	-19.47#
75	I	Chrysene-d12	1.000	1.000	0.0	83	-0.02
76		Benzidine	0.605	0.028	95.4#	4#	0.36
77		Pyrene	1.200	0.006	99.5#	0#	0.18
78	S	Terphenyl-d14	0.761	0.833	-9.5	91	-0.02
79		Butylbenzylphthalate	0.478	0.000	100.0#	0#	-0.01
80		Benzo(a)anthracene	1.167	0.003	99.7#	0#	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.000	100.0#	0#	-21.58#
82		Chrysene	1.128	0.001	99.9#	0#	-0.02
83		Bis(2-ethylhexyl)phthalate	0.668	0.001	99.9#	0#	-0.02
84	c	Di-n-octyl phthalate	1.100	0.001	99.9#	0#	-0.04
85		Indeno(1,2,3-cd)pyrene	1.211	0.000	100.0#	0#	-28.55#
86	I	Perylene-d12	1.000	1.000	0.0	80	-0.05
87		Benzo(b)fluoranthene	1.066	0.000	100.0#	0#	0.00

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	Compound	AvgRF	CCRF	%Dev Area%	Dev(min)
88	Benzo(k)fluoranthene	1.069	0.000	100.0# 0#	-0.04
89 C	Benzo(a)pyrene	1.030	0.000	100.0# 0#	-0.05
90	Dibenzo(a,h)anthracene	0.966	0.000	100.0# 0#	-28.62#
91	Benzo(g,h,i)perylene	0.969	0.000	100.0# 0#	-29.70#

(#) = Out of Range SPCC's out = 4 CCC's out = 13