Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG100419\
Data File : BG043045.D
Acq On : 5 Oct 2019 3:35
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 05 06:26:30 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 25 15:35:52 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	105	-0.02
2		1,4-Dioxane	0.467	0.002	99.6#	1#	-0.18
3		Pyridine	1.314	0.000	100.0#	0#	-0.01
4		n-Nitrosodimethylamine	0.632	0.000	100.0#	0#	0.00
5	S	2-Fluorophenol	1.121	1.303	-16.2	126	-0.02
6		Aniline	1.948	0.001	99.9#	0#	-0.01
7	S	Phenol-d6	1.614	1.224	24.2	80	-0.01
8		2-Chlorophenol	1.169	0.000	100.0#	0#	-0.17
9		Benzaldehyde	0.986	0.008	99.2#	1#	0.03
10	C	Phenol	1.481	0.347	76.6#	25#	-0.02
11		bis(2-Chloroethyl)ether	1.146	0.001	99.9#	0#	0.06
12		1,3-Dichlorobenzene	1.491	0.001	99.9#	0#	-0.17
13	C	1,4-Dichlorobenzene	1.486	0.001	99.9#	0#	-0.31
14		1,2-Dichlorobenzene	1.415	0.000	100.0#	0#	-8.43#
15		Benzyl Alcohol	1.213	0.056	95.4#	5#	0.06
16		2,2'-oxybis(1-Chloropropane	2.200	0.000	100.0#	0#	-0.08
17		2-Methylphenol	1.036	0.000	100.0#	0#	0.03
18		Hexachloroethane	0.560	0.000	100.0#	0#	-9.17#
19	P	n-Nitroso-di-n-propylamine	1.060	0.470	55.7#	46#	0.33
20		3+4-Methylphenols	1.420	0.000	100.0#	0#	-0.30
21	I	Naphthalene-d8	1.000	1.000	0.0	96	-0.02
22		Acetophenone	0.516	0.003	99.4#	0#	-0.02
23	S	Nitrobenzene-d5	0.411	0.446	-8.5	104	-0.01
24		Nitrobenzene	0.392	0.003	99.2#	1#	-0.06
25		Isophorone	0.723	0.001	99.9#	0#	0.00
26	C	2-Nitrophenol	0.167	0.000	100.0#	0#	-9.99#
27		2,4-Dimethylphenol	0.257	0.000	100.0#	0#	-10.05#
28		bis(2-Chloroethoxy)methane	0.410	0.000	100.0#	0#	0.32
29	C	2,4-Dichlorophenol	0.319	0.000	100.0#	0#	-10.53#
30		1,2,4-Trichlorobenzene	0.412	0.000	100.0#	0#	-10.74#
31		Naphthalene	0.985	0.000	100.0#	0#	0.00
32		Benzoic acid	0.194	0.000	100.0#	0#	-10.20#
33		4-Chloroaniline	0.427	0.000	100.0#	0#	-0.12
34	C	Hexachlorobutadiene	0.309	0.000	100.0#	0#	-0.11
35		Caprolactam	0.113	0.017	85.0#	14#	-0.05
36	C	4-Chloro-3-methylphenol	0.347	0.000	100.0#	0#	0.03
37		2-Methylnaphthalene	0.751	0.000	100.0#	0#	-12.53#
38		1-Methylnaphthalene	0.711	0.000	100.0#	0#	-12.74#
39	I	Acenaphthene-d10	1.000	1.000	0.0	91	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.752	0.000	100.0#	0#	0.30
41	P	Hexachlorocyclopentadiene	0.479	0.000#	100.0#	0#	-0.19
42	S	2,4,6-Tribromophenol	0.344	0.624	-81.4#	174#	-0.01
43	C	2,4,6-Trichlorophenol	0.440	0.000	100.0#	0#	-13.13#
44		2,4,5-Trichlorophenol	0.453	0.000	100.0#	0#	-13.21#

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.380	1.505	-9.1	102	-0.02
46		1,1'-Biphenyl	1.517	0.004	99.7#	0#	-0.01
47		2-Chloronaphthalene	1.137	0.000	100.0#	0#	0.29
48		2-Nitroaniline	0.378	0.000	100.0#	0#	-0.10
49		Acenaphthylene	1.740	0.000	100.0#	0#	-0.09
50		Dimethylphthalate	1.499	0.525	65.0#	33#	-0.02
51		2,6-Dinitrotoluene	0.319	0.030	90.6#	9#	-0.05
52	C	Acenaphthene	1.165	0.000	100.0#	0#	-0.01
53		3-Nitroaniline	0.330	0.000	100.0#	0#	-0.15
54	P	2,4-Dinitrophenol	0.198	0.000#	100.0#	0#	-0.08
55		Dibenzofuran	1.723	0.000	100.0#	0#	0.07
56	P	4-Nitrophenol	0.251	0.000#	100.0#	0#	0.02
57		2,4-Dinitrotoluene	0.467	0.067	85.7#	14#	-0.37
58		Fluorene	1.471	0.036	97.6#	2#	0.43
59		2,3,4,6-Tetrachlorophenol	0.483	0.000	100.0#	0#	-0.17
60		Diethylphthalate	1.525	0.003	99.8#	0#	-0.02
61		4-Chlorophenyl-phenylether	0.882	0.000	100.0#	0#	-15.73#
62		4-Nitroaniline	0.360	0.000	100.0#	0#	0.17
63		Azobenzene	1.388	0.001	99.9#	0#	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	96	-0.01
65		4,6-Dinitro-2-methylphenol	0.113	0.000	100.0#	0#	0.03
66	c	n-Nitrosodiphenylamine	0.528	0.015	97.2#	3#	0.22
67		4-Bromophenyl-phenylether	0.251	0.033	86.9#	12#	-0.45
68		Hexachlorobenzene	0.279	0.000	100.0#	0#	-16.75#
69		Atrazine	0.222	0.000	100.0#	0#	0.15
70	C	Pentachlorophenol	0.161	0.000	100.0#	0#	0.27
71		Phenanthrene	0.976	0.002	99.8#	0#	-0.02
72		Anthracene	0.975	0.002	99.8#	0#	-0.11
73		Carbazole	0.904	0.002	99.8#	0#	-0.01
74		Di-n-butylphthalate	1.078	0.006	99.4#	1#	-0.02
75	C	Fluoranthene	1.291	0.003	99.8#	0#	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	98	-0.02
77		Benzidine	0.501	0.036	92.8#	7#	0.36
78		Pyrene	1.194	0.002	99.8#	0#	-0.37
79	S	Terphenyl-d14	0.939	1.122	-19.5	110	-0.01
80		Butylbenzylphthalate	0.453	0.003	99.3#	1#	-0.02
81		Benzo(a)anthracene	1.246	0.004	99.7#	0#	-0.02
82		3,3'-Dichlorobenzidine	0.489	0.000	100.0#	0#	-0.03
83		Chrysene	1.209	0.002	99.8#	0#	-0.02
84		Bis(2-ethylhexyl)phthalate	0.645	0.007	98.9#	1#	-0.02
85	c	Di-n-octyl phthalate	1.054	0.000	100.0#	0#	-0.02
86		Indeno(1,2,3-cd)pyrene	1.506	0.000	100.0#	0#	-0.10
87	I	Perylene-d12	1.000	1.000	0.0	100	-0.04

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	Compound	AvgRF	CCRF	%Dev Area%	Dev(min)
88	Benzo(b)fluoranthene	1.132	0.001	99.9#	0# -0.07
89	Benzo(k)fluoranthene	1.120	0.000	100.0#	0# -0.07
90 C	Benzo(a)pyrene	1.068	0.000	100.0#	0# -0.04
91	Dibenzo(a,h)anthracene	1.095	0.000	100.0#	0# -0.02
92	Benzo(g,h,i)perylene	1.061	0.000	100.0#	0# -0.05

(#) = Out of Range SPCC's out = 3 CCC's out = 13