Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF090618\
Data File : BF108824.D
Acq On : 6 Sep 2018 15:38
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 06 16:11:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 05 18:03:31 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	102	0.00
2		1,4-Dioxane	0.580	0.579	0.2	104	0.00
3		Pyridine	1.589	1.655	-4.2	107	0.00
4		n-Nitrosodimethylamine	0.610	0.621	-1.8	103	0.01
5	S	2-Fluorophenol	1.207	1.156	4.2	102	0.01
6		Aniline	2.089	2.054	1.7	102	0.00
7	S	Phenol-d6	1.553	1.535	1.2	105	0.01
8		2-Chlorophenol	1.336	1.336	0.0	104	0.00
9		Benzaldehyde	1.102	1.098	0.4	107	0.00
10	C	Phenol	1.779	1.707	4.0	100	0.01
11		bis(2-Chloroethyl)ether	1.425	1.405	1.4	104	0.00
12		1,3-Dichlorobenzene	1.525	1.496	1.9	103	0.00
13	C	1,4-Dichlorobenzene	1.525	1.514	0.7	105	0.00
14		1,2-Dichlorobenzene	1.406	1.362	3.1	102	0.00
15		Benzyl Alcohol	1.192	1.208	-1.3	105	0.00
16		2,2'-oxybis(1-Chloropropane	2.130	2.146	-0.8	105	0.00
17		2-Methylphenol	1.129	1.136	-0.6	104	0.00
18		Hexachloroethane	0.550	0.563	-2.4	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.080	1.070	0.9	104	0.00
20		3+4-Methylphenols	1.323	1.274	3.7	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.454	0.430	5.3	104	0.00
23	S	Nitrobenzene-d5	0.320	0.351	-9.7	111	0.00
24		Nitrobenzene	0.343	0.369	-7.6	108	0.00
25		Isophorone	0.637	0.637	0.0	107	0.00
26	C	2-Nitrophenol	0.135	0.172	-27.4#	127	0.00
27		2,4-Dimethylphenol	0.274	0.278	-1.5	106	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.422	0.9	105	0.00
29	C	2,4-Dichlorophenol	0.283	0.291	-2.8	107	0.00
30		1,2,4-Trichlorobenzene	0.307	0.303	1.3	105	0.00
31		Naphthalene	0.900	0.876	2.7	104	0.00
32		Benzoic acid	0.168	0.224	-33.3#	132	0.02
33		4-Chloroaniline	0.413	0.412	0.2	106	0.00
34	C	Hexachlorobutadiene	0.183	0.182	0.5	105	0.00
35		Caprolactam	0.096	0.101	-5.2	115	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.302	-0.3	106	0.01
37		2-Methylnaphthalene	0.594	0.579	2.5	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.598	0.580	3.0	108	0.00
40	P	Hexachlorocyclopentadiene	0.301	0.325	-8.0	113	0.00
41	S	2,4,6-Tribromophenol	0.191	0.204	-6.8	116	0.00
42	C	2,4,6-Trichlorophenol	0.407	0.423	-3.9	110	0.01
43		2,4,5-Trichlorophenol	0.422	0.443	-5.0	111	0.00
44	S	2-Fluorobiphenyl	1.172	1.143	2.5	106	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.540	1.454	5.6	107	0.00
46		2-Chloronaphthalene	1.245	1.190	4.4	106	0.00
47		2-Nitroaniline	0.340	0.396	-16.5	120	0.00
48		Acenaphthylene	1.856	1.754	5.5	106	0.00
49		Dimethylphthalate	1.399	1.369	2.1	111	0.01
50		2,6-Dinitrotoluene	0.267	0.313	-17.2	117	0.01
51	C	Acenaphthene	1.149	1.130	1.7	111	0.00
52		3-Nitroaniline	0.317	0.373	-17.7	118	0.00
53	P	2,4-Dinitrophenol	0.078	0.112	-43.6#	156#	0.01
54		Dibenzofuran	1.674	1.575	5.9	108	0.00
55	P	4-Nitrophenol	0.237	0.284	-19.8	119	0.02
56		2,4-Dinitrotoluene	0.341	0.399	-17.0	120	0.01
57		Fluorene	1.074	1.068	0.6	110	0.00
58		2,3,4,6-Tetrachlorophenol	0.340	0.364	-7.1	113	0.00
59		Diethylphthalate	1.444	1.378	4.6	106	0.00
60		4-Chlorophenyl-phenylether	0.574	0.578	-0.7	111	0.00
61		4-Nitroaniline	0.287	0.345	-20.2	122	0.01
62		Azobenzene	1.490	1.444	3.1	109	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
64		4,6-Dinitro-2-methylphenol	0.070	0.092	-31.4#	139	0.01
65	c	n-Nitrosodiphenylamine	0.665	0.630	5.3	109	0.00
66		4-Bromophenyl-phenylether	0.226	0.218	3.5	111	0.00
67		Hexachlorobenzene	0.241	0.235	2.5	112	0.01
68		Atrazine	0.213	0.211	0.9	114	0.00
69	C	Pentachlorophenol	0.147	0.163	-10.9	115	0.00
70		Phenanthrene	0.955	0.885	7.3	108	0.01
71		Anthracene	0.914	0.909	0.5	110	0.00
72		Carbazole	0.963	0.908	5.7	111	0.00
73		Di-n-butylphthalate	1.069	1.058	1.0	110	0.00
74	C	Fluoranthene	0.946	0.943	0.3	110	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	116	0.01
76		Benzidine	0.948	0.918	3.2	109	0.01
77		Pyrene	1.539	1.488	3.3	110	0.00
78	S	Terphenyl-d14	0.858	0.797	7.1	111	0.00
79		Butylbenzylphthalate	0.693	0.718	-3.6	114	0.00
80		Benzo(a)anthracene	1.282	1.192	7.0	108	0.01
81		3,3'-Dichlorobenzidine	0.499	0.524	-5.0	119	0.00
82		Chrysene	1.241	1.205	2.9	111	0.01
83		Bis(2-ethylhexyl)phthalate	0.871	0.827	5.1	104	0.00
84	c	Di-n-octyl phthalate	1.452	1.482	-2.1	115	0.00
85		Indeno(1,2,3-cd)pyrene	1.098	0.901	17.9	97	0.02
86	I	Perylene-d12	1.000	1.000	0.0	118	0.02
87		Benzo(b)fluoranthene	1.088	1.050	3.5	119	0.01

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88	Benzo(k)fluoranthene	1.001	0.995	0.6	117	0.01
89 C	Benzo(a)pyrene	0.981	0.952	3.0	117	0.01
90	Dibenzo(a,h)anthracene	0.865	0.716	17.2	97	0.02
91	Benzo(g,h,i)perylene	0.866	0.693	20.0	93	0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1