Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF032317\
Data File : BF093877.D
Acq On : 23 Mar 2017 14:12
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 24 01:03:46 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 23 02:02:22 2017
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	100	-0.01
2		1,4-Dioxane	0.569	0.559	1.8	97	-0.02
3		Pyridine	1.550	1.492	3.7	93	-0.02
4		n-Nitrosodimethylamine	0.638	0.622	2.5	94	-0.01
5	S	2-Fluorophenol	1.184	1.181	0.3	100	-0.01
6		Aniline	2.038	2.075	-1.8	99	-0.01
7	S	Phenol-d6	1.465	1.474	-0.6	100	0.00
8		2-Chlorophenol	1.302	1.325	-1.8	99	-0.01
9		Benzaldehyde	0.989	0.955	3.4	95	-0.01
10	C	Phenol	1.667	1.733	-4.0	98	-0.01
11		bis(2-Chloroethyl)ether	1.303	1.309	-0.5	99	-0.01
12		1,3-Dichlorobenzene	1.460	1.479	-1.3	99	-0.01
13	C	1,4-Dichlorobenzene	1.479	1.490	-0.7	99	-0.01
14		1,2-Dichlorobenzene	1.362	1.386	-1.8	101	-0.01
15		Benzyl Alcohol	1.072	1.114	-3.9	100	-0.01
16		2,2'-oxybis(1-Chloropropane	2.007	1.991	0.8	96	0.00
17		2-Methylphenol	1.115	1.141	-2.3	100	-0.01
18		Hexachloroethane	0.520	0.528	-1.5	98	-0.01
19	P	n-Nitroso-di-n-propylamine	0.941	0.959	-1.9	100	-0.01
20		3+4-Methylphenols	1.350	1.343	0.5	100	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	100	-0.01
22		Acetophenone	0.446	0.435	2.5	100	-0.01
23	S	Nitrobenzene-d5	0.350	0.350	0.0	99	-0.01
24		Nitrobenzene	0.346	0.342	1.2	97	-0.01
25		Isophorone	0.616	0.620	-0.6	100	-0.01
26	C	2-Nitrophenol	0.165	0.178	-7.9	101	-0.01
27		2,4-Dimethylphenol	0.284	0.284	0.0	99	-0.01
28		bis(2-Chloroethoxy)methane	0.406	0.406	0.0	99	-0.01
29	C	2,4-Dichlorophenol	0.266	0.274	-3.0	101	-0.01
30		1,2,4-Trichlorobenzene	0.274	0.273	0.4	99	-0.01
31		Naphthalene	0.962	0.952	1.0	100	-0.01
32		Benzoic acid	0.189	0.196	-3.7	91	0.00
33		4-Chloroaniline	0.390	0.395	-1.3	100	-0.01
34	C	Hexachlorobutadiene	0.140	0.140	0.0	99	-0.01
35		Caprolactam	0.085	0.088	-3.5	101	0.00
36	C	4-Chloro-3-methylphenol	0.277	0.284	-2.5	101	0.00
37		2-Methylnaphthalene	0.641	0.646	-0.8	101	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.547	0.541	1.1	102	0.00
40	P	Hexachlorocyclopentadiene	0.256	0.270	-5.5	100	-0.01
41	S	2,4,6-Tribromophenol	0.158	0.166	-5.1	105	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.396	-2.3	104	0.00
43		2,4,5-Trichlorophenol	0.419	0.430	-2.6	101	-0.01
44	S	2-Fluorobiphenyl	1.311	1.309	0.2	103	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.580	2.0	100	-0.01
46		2-Chloronaphthalene	1.263	1.234	2.3	101	0.00
47		2-Nitroaniline	0.375	0.387	-3.2	101	-0.01
48		Acenaphthylene	2.099	2.078	1.0	101	-0.01
49		Dimethylphthalate	1.422	1.414	0.6	102	0.00
50		2,6-Dinitrotoluene	0.326	0.334	-2.5	101	0.00
51	C	Acenaphthene	1.225	1.185	3.3	101	0.00
52		3-Nitroaniline	0.383	0.395	-3.1	104	0.00
53	P	2,4-Dinitrophenol	0.155	0.167	-7.7	105	0.00
54		Dibenzofuran	1.667	1.632	2.1	99	0.00
55	P	4-Nitrophenol	0.300	0.316	-5.3	102	0.00
56		2,4-Dinitrotoluene	0.409	0.427	-4.4	101	-0.01
57		Fluorene	1.382	1.292	6.5	103	-0.01
58		2,3,4,6-Tetrachlorophenol	0.287	0.302	-5.2	105	0.00
59		Diethylphthalate	1.405	1.422	-1.2	104	0.00
60		4-Chlorophenyl-phenylether	0.589	0.558	5.3	102	-0.01
61		4-Nitroaniline	0.367	0.392	-6.8	106	0.00
62		Azobenzene	1.329	1.312	1.3	100	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.134	-9.8	106	0.00
65	c	n-Nitrosodiphenylamine	0.692	0.682	1.4	103	-0.01
66		4-Bromophenyl-phenylether	0.197	0.200	-1.5	104	-0.01
67		Hexachlorobenzene	0.199	0.202	-1.5	105	-0.01
68		Atrazine	0.207	0.212	-2.4	105	0.00
69	C	Pentachlorophenol	0.124	0.129	-4.0	104	0.00
70		Phenanthrene	1.113	1.079	3.1	102	0.00
71		Anthracene	1.146	1.128	1.6	103	0.00
72		Carbazole	1.175	1.153	1.9	102	0.00
73		Di-n-butylphthalate	1.222	1.241	-1.6	104	0.00
74	C	Fluoranthene	1.139	1.127	1.1	104	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76		Benzidine	0.792	0.831	-4.9	107	0.00
77		Pyrene	1.451	1.431	1.4	103	0.00
78	S	Terphenyl-d14	0.892	0.865	3.0	103	0.00
79		Butylbenzylphthalate	0.618	0.656	-6.1	105	0.00
80		Benzo(a)anthracene	1.166	1.176	-0.9	104	0.00
81		3,3'-Dichlorobenzidine	0.417	0.441	-5.8	105	0.00
82		Chrysene	1.082	1.074	0.7	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.844	0.884	-4.7	105	0.00
84	c	Di-n-octyl phthalate	1.410	1.531	-8.6	107	0.00
85		Indeno(1,2,3-cd)pyrene	0.937	0.980	-4.6	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	115	0.00
87		Benzo(b)fluoranthene	1.226	1.264	-3.1	116	0.00

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88	Benzo(k)fluoranthene	1.199	1.132	5.6	105	0.00
89 C	Benzo(a)pyrene	1.129	1.131	-0.2	112	0.00
90	Dibenzo(a,h)anthracene	0.956	0.955	0.1	115	0.00
91	Benzo(g,h,i)perylene	0.964	0.930	3.5	112	0.00

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