Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF012617\
Data File : BF092532.D
Acq On : 26 Jan 2017 18:36
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 27 01:14:45 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 24 13:48:07 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	111	0.01
2		1,4-Dioxane	0.684	0.628	8.2	103	0.00
3		Pyridine	1.946	1.775	8.8	103	0.01
4		n-Nitrosodimethylamine	0.955	0.838	12.3	97	0.01
5	S	2-Fluorophenol	1.285	1.228	4.4	109	0.01
6		Aniline	2.442	2.380	2.5	107	0.01
7	S	Phenol-d6	1.637	1.553	5.1	107	0.01
8		2-Chlorophenol	1.350	1.392	-3.1	114	0.01
9		Benzaldehyde	1.162	1.087	6.5	107	0.01
10	C	Phenol	1.976	1.902	3.7	103	0.01
11		bis(2-Chloroethyl)ether	1.479	1.305	11.8	93	0.00
12		1,3-Dichlorobenzene	1.597	1.541	3.5	107	0.01
13	C	1,4-Dichlorobenzene	1.607	1.521	5.4	101	0.00
14		1,2-Dichlorobenzene	1.525	1.511	0.9	117	0.01
15		Benzyl Alcohol	1.234	1.182	4.2	105	0.01
16		2,2'-oxybis(1-Chloropropane	2.925	2.722	6.9	104	0.00
17		2-Methylphenol	1.154	1.143	1.0	108	0.01
18		Hexachloroethane	0.554	0.535	3.4	104	0.00
19	P	n-Nitroso-di-n-propylamine	1.115	1.016	8.9	104	0.00
20		3+4-Methylphenols	1.400	1.352	3.4	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.476	0.442	7.1	103	0.00
23	S	Nitrobenzene-d5	0.366	0.382	-4.4	116	0.01
24		Nitrobenzene	0.384	0.336	12.5	90	0.00
25		Isophorone	0.721	0.677	6.1	105	0.00
26	C	2-Nitrophenol	0.161	0.179	-11.2	126	0.01
27		2,4-Dimethylphenol	0.334	0.294	12.0	99	0.00
28		bis(2-Chloroethoxy)methane	0.474	0.466	1.7	117	0.01
29	C	2,4-Dichlorophenol	0.290	0.277	4.5	107	0.01
30		1,2,4-Trichlorobenzene	0.313	0.314	-0.3	118	0.01
31		Naphthalene	1.024	0.951	7.1	109	0.01
32		Benzoic acid	0.231	0.139	39.8#	61	0.00
33		4-Chloroaniline	0.433	0.410	5.3	112	0.01
34	C	Hexachlorobutadiene	0.171	0.160	6.4	104	0.00
35		Caprolactam	0.097	0.090	7.2	101	0.01
36	C	4-Chloro-3-methylphenol	0.313	0.302	3.5	103	0.00
37		2-Methylnaphthalene	0.648	0.634	2.2	117	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.01
39		1,2,4,5-Tetrachlorobenzene	0.504	0.525	-4.2	112	0.01
40	P	Hexachlorocyclopentadiene	0.253	0.245	3.2	101	0.00
41	S	2,4,6-Tribromophenol	0.136	0.145	-6.6	118	0.01
42	C	2,4,6-Trichlorophenol	0.387	0.380	1.8	111	0.00
43		2,4,5-Trichlorophenol	0.354	0.384	-8.5	114	0.01
44	S	2-Fluorobiphenyl	1.082	1.024	5.4	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.446	1.448	-0.1	121	0.01
46		2-Chloronaphthalene	1.181	1.112	5.8	107	0.01
47		2-Nitroaniline	0.398	0.363	8.8	102	0.01
48		Acenaphthylene	1.732	1.737	-0.3	111	0.01
49		Dimethylphthalate	1.272	1.168	8.2	98	0.00
50		2,6-Dinitrotoluene	0.260	0.284	-9.2	108	0.01
51	C	Acenaphthene	1.076	1.113	-3.4	124	0.01
52		3-Nitroaniline	0.326	0.322	1.2	92	0.00
53	P	2,4-Dinitrophenol	0.093	0.140	-50.5#	149	0.00
54		Dibenzofuran	1.461	1.461	0.0	117	0.01
55	P	4-Nitrophenol	0.259	0.296	-14.3	110	0.01
56		2,4-Dinitrotoluene	0.345	0.410	-18.8	135	0.01
57		Fluorene	1.084	1.165	-7.5	128	0.01
58		2,3,4,6-Tetrachlorophenol	0.265	0.286	-7.9	120	0.01
59		Diethylphthalate	1.220	1.152	5.6	108	0.00
60		4-Chlorophenyl-phenylether	0.526	0.539	-2.5	113	0.00
61		4-Nitroaniline	0.326	0.326	0.0	107	0.00
62		Azobenzene	1.388	1.331	4.1	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.099	0.137	-38.4#	139	0.01
65	c	n-Nitrosodiphenylamine	0.625	0.597	4.5	114	0.00
66		4-Bromophenyl-phenylether	0.202	0.203	-0.5	114	0.00
67		Hexachlorobenzene	0.204	0.208	-2.0	120	0.01
68		Atrazine	0.214	0.232	-8.4	117	0.00
69	C	Pentachlorophenol	0.131	0.134	-2.3	108	0.01
70		Phenanthrene	1.101	0.967	12.2	101	0.00
71		Anthracene	1.076	1.113	-3.4	119	0.01
72		Carbazole	1.028	0.939	8.7	101	0.00
73		Di-n-butylphthalate	1.175	1.167	0.7	115	0.00
74	C	Fluoranthene	1.077	0.999	7.2	110	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.742	0.836	-12.7	107	0.00
77		Pyrene	1.449	1.453	-0.3	113	0.00
78	S	Terphenyl-d14	0.751	0.758	-0.9	115	0.00
79		Butylbenzylphthalate	0.587	0.638	-8.7	104	-0.01
80		Benzo(a)anthracene	1.075	1.119	-4.1	110	0.00
81		3,3'-Dichlorobenzidine	0.408	0.437	-7.1	105	0.00
82		Chrysene	1.148	1.197	-4.3	115	0.01
83		Bis(2-ethylhexyl)phthalate	0.739	0.832	-12.6	114	-0.01
84	c	Di-n-octyl phthalate	1.342	1.491	-11.1	104	-0.01
85		Indeno(1,2,3-cd)pyrene	0.849	0.896	-5.5	109	0.02
86	I	Perylene-d12	1.000	1.000	0.0	100	0.01
87		Benzo(b)fluoranthene	1.284	1.163	9.4	88	0.01

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88	Benzo(k)fluoranthene	1.068	1.164	-9.0	123	0.01
89 C	Benzo(a)pyrene	1.086	1.094	-0.7	102	0.01
90	Dibenzo(a,h)anthracene	0.878	0.901	-2.6	107	0.02
91	Benzo(g,h,i)perylene	0.888	0.907	-2.1	111	0.02

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