Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012517\
Data File : BF092481.D
Acq On : 25 Jan 2017 17:22
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 25 18:25:55 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	108	0.00
2		1,4-Dioxane	0.684	0.645	5.7	103	-0.01
3		Pyridine	1.946	1.813	6.8	102	0.00
4		n-Nitrosodimethylamine	0.955	0.879	8.0	99	0.00
5	S	2-Fluorophenol	1.285	1.291	-0.5	111	0.01
6		Aniline	2.442	2.423	0.8	106	0.00
7	S	Phenol-d6	1.637	1.538	6.0	103	0.00
8		2-Chlorophenol	1.350	1.346	0.3	107	0.00
9		Benzaldehyde	1.162	1.061	8.7	102	0.00
10	C	Phenol	1.976	1.813	8.2	95	0.00
11		bis(2-Chloroethyl)ether	1.479	1.501	-1.5	104	0.00
12		1,3-Dichlorobenzene	1.597	1.605	-0.5	108	0.00
13	C	1,4-Dichlorobenzene	1.607	1.683	-4.7	108	0.00
14		1,2-Dichlorobenzene	1.525	1.413	7.3	107	0.00
15		Benzyl Alcohol	1.234	1.185	4.0	102	0.00
16		2,2'-oxybis(1-Chloropropane	2.925	2.748	6.1	102	0.00
17		2-Methylphenol	1.154	1.108	4.0	102	0.00
18		Hexachloroethane	0.554	0.573	-3.4	109	0.00
19	P	n-Nitroso-di-n-propylamine	1.115	1.038	6.9	103	0.00
20		3+4-Methylphenols	1.400	1.405	-0.4	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	-0.01
22		Acetophenone	0.476	0.470	1.3	106	0.00
23	S	Nitrobenzene-d5	0.366	0.362	1.1	106	0.00
24		Nitrobenzene	0.384	0.402	-4.7	104	0.00
25		Isophorone	0.721	0.694	3.7	104	0.00
26	C	2-Nitrophenol	0.161	0.176	-9.3	120	0.00
27		2,4-Dimethylphenol	0.334	0.327	2.1	107	0.00
28		bis(2-Chloroethoxy)methane	0.474	0.437	7.8	106	0.00
29	C	2,4-Dichlorophenol	0.290	0.288	0.7	107	0.00
30		1,2,4-Trichlorobenzene	0.313	0.300	4.2	109	0.00
31		Naphthalene	1.024	0.954	6.8	105	0.00
32		Benzoic acid	0.231	0.220	4.8	93	0.00
33		4-Chloroaniline	0.433	0.412	4.8	108	0.00
34	C	Hexachlorobutadiene	0.171	0.174	-1.8	110	0.00
35		Caprolactam	0.097	0.090	7.2	98	0.00
36	C	4-Chloro-3-methylphenol	0.313	0.332	-6.1	109	0.00
37		2-Methylnaphthalene	0.648	0.601	7.3	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.504	0.527	-4.6	113	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.268	-5.9	111	0.00
41	S	2,4,6-Tribromophenol	0.136	0.137	-0.7	112	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.373	3.6	109	-0.01
43		2,4,5-Trichlorophenol	0.354	0.363	-2.5	108	0.00
44	S	2-Fluorobiphenyl	1.082	1.011	6.6	101	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.446	1.272	12.0	107	0.00
46		2-Chloronaphthalene	1.181	1.028	13.0	99	0.00
47		2-Nitroaniline	0.398	0.354	11.1	99	0.00
48		Acenaphthylene	1.732	1.770	-2.2	113	0.00
49		Dimethylphthalate	1.272	1.286	-1.1	109	0.00
50		2,6-Dinitrotoluene	0.260	0.297	-14.2	113	0.00
51	C	Acenaphthene	1.076	1.103	-2.5	123	0.00
52		3-Nitroaniline	0.326	0.318	2.5	91	-0.01
53	P	2,4-Dinitrophenol	0.093	0.127	-36.6#	136	-0.01
54		Dibenzofuran	1.461	1.451	0.7	117	0.00
55	P	4-Nitrophenol	0.259	0.259	0.0	97	0.00
56		2,4-Dinitrotoluene	0.345	0.394	-14.2	130	0.00
57		Fluorene	1.084	1.111	-2.5	122	0.00
58		2,3,4,6-Tetrachlorophenol	0.265	0.284	-7.2	120	0.00
59		Diethylphthalate	1.220	1.214	0.5	114	0.00
60		4-Chlorophenyl-phenylether	0.526	0.470	10.6	99	-0.01
61		4-Nitroaniline	0.326	0.336	-3.1	111	-0.01
62		Azobenzene	1.388	1.236	11.0	94	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	-0.01
64		4,6-Dinitro-2-methylphenol	0.099	0.134	-35.4#	131	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.655	-4.8	121	0.00
66		4-Bromophenyl-phenylether	0.202	0.186	7.9	101	-0.01
67		Hexachlorobenzene	0.204	0.212	-3.9	119	0.00
68		Atrazine	0.214	0.218	-1.9	106	-0.01
69	C	Pentachlorophenol	0.131	0.132	-0.8	104	-0.01
70		Phenanthrene	1.101	0.950	13.7	96	-0.01
71		Anthracene	1.076	1.119	-4.0	116	0.00
72		Carbazole	1.028	0.910	11.5	94	-0.01
73		Di-n-butylphthalate	1.175	1.105	6.0	105	-0.01
74	C	Fluoranthene	1.077	0.993	7.8	105	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	109	-0.01
76		Benzidine	0.742	0.725	2.3	101	-0.01
77		Pyrene	1.449	1.225	15.5	104	-0.01
78	S	Terphenyl-d14	0.751	0.661	12.0	109	-0.01
79		Butylbenzylphthalate	0.587	0.576	1.9	102	-0.02
80		Benzo(a)anthracene	1.075	1.011	6.0	108	-0.01
81		3,3'-Dichlorobenzidine	0.408	0.396	2.9	104	-0.01
82		Chrysene	1.148	0.959	16.5	100	-0.01
83		Bis(2-ethylhexyl)phthalate	0.739	0.727	1.6	109	-0.02
84	c	Di-n-octyl phthalate	1.342	1.348	-0.4	103	-0.02
85		Indeno(1,2,3-cd)pyrene	0.849	0.783	7.8	104	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	102	-0.01
87		Benzo(b)fluoranthene	1.284	1.140	11.2	87	-0.01

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88	Benzo(k)fluoranthene	1.068	1.164	-9.0	124	-0.01
89 C	Benzo(a)pyrene	1.086	1.089	-0.3	103	-0.01
90	Dibenzo(a,h)anthracene	0.878	0.870	0.9	104	-0.01
91	Benzo(g,h,i)perylene	0.888	0.850	4.3	105	-0.02

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