Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022017\
Data File : BF093055.D
Acq On : 21 Feb 2017 5:23
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 21 06:12:45 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 17 17:31:55 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	105	-0.01
2		1,4-Dioxane	0.630	0.655	-4.0	112	-0.03
3		Pyridine	1.602	1.832	-14.4	119	-0.03
4		n-Nitrosodimethylamine	0.752	0.751	0.1	105	-0.02
5	S	2-Fluorophenol	1.166	1.234	-5.8	107	-0.01
6		Aniline	2.046	2.202	-7.6	117	0.00
7	S	Phenol-d6	1.500	1.521	-1.4	107	0.00
8		2-Chlorophenol	1.286	1.377	-7.1	112	0.00
9		Benzaldehyde	1.075	0.940	12.6	90	0.00
10	C	Phenol	1.755	1.646	6.2	104	-0.01
11		bis(2-Chloroethyl)ether	1.401	1.461	-4.3	115	0.00
12		1,3-Dichlorobenzene	1.506	1.543	-2.5	111	-0.01
13	C	1,4-Dichlorobenzene	1.525	1.623	-6.4	116	-0.01
14		1,2-Dichlorobenzene	1.458	1.375	5.7	98	0.00
15		Benzyl Alcohol	1.110	1.090	1.8	107	-0.01
16		2,2'-oxybis(1-Chloropropane	2.434	2.547	-4.6	112	-0.01
17		2-Methylphenol	1.103	1.197	-8.5	109	0.00
18		Hexachloroethane	0.520	0.523	-0.6	106	-0.01
19	P	n-Nitroso-di-n-propylamine	0.955	1.040	-8.9	126	0.00
20		3+4-Methylphenols	1.319	1.415	-7.3	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.457	0.461	-0.9	109	0.00
23	S	Nitrobenzene-d5	0.359	0.385	-7.2	111	0.00
24		Nitrobenzene	0.369	0.357	3.3	108	-0.01
25		Isophorone	0.669	0.705	-5.4	113	-0.01
26	C	2-Nitrophenol	0.175	0.198	-13.1	109	0.00
27		2,4-Dimethylphenol	0.308	0.338	-9.7	110	0.00
28		bis(2-Chloroethoxy)methane	0.443	0.501	-13.1	119	0.00
29	C	2,4-Dichlorophenol	0.287	0.282	1.7	108	-0.01
30		1,2,4-Trichlorobenzene	0.309	0.310	-0.3	107	-0.01
31		Naphthalene	1.014	0.993	2.1	106	-0.01
32		Benzoic acid	0.156	0.192	-23.1	116	0.01
33		4-Chloroaniline	0.398	0.409	-2.8	105	0.00
34	C	Hexachlorobutadiene	0.170	0.166	2.4	102	0.00
35		Caprolactam	0.090	0.105	-16.7	114	0.01
36	C	4-Chloro-3-methylphenol	0.299	0.323	-8.0	111	0.00
37		2-Methylnaphthalene	0.651	0.672	-3.2	108	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.561	0.535	4.6	106	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.183	27.7#	78	0.00
41	S	2,4,6-Tribromophenol	0.145	0.127	12.4	88	0.00
42	C	2,4,6-Trichlorophenol	0.369	0.368	0.3	103	0.00
43		2,4,5-Trichlorophenol	0.404	0.383	5.2	107	0.00
44	S	2-Fluorobiphenyl	1.104	1.069	3.2	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.448	1.515	-4.6	110	0.00
46		2-Chloronaphthalene	1.133	1.190	-5.0	108	0.00
47		2-Nitroaniline	0.381	0.398	-4.5	122	0.00
48		Acenaphthylene	1.957	1.789	8.6	108	0.00
49		Dimethylphthalate	1.347	1.274	5.4	103	0.00
50		2,6-Dinitrotoluene	0.291	0.321	-10.3	117	0.00
51	C	Acenaphthene	1.170	1.072	8.4	106	0.00
52		3-Nitroaniline	0.333	0.377	-13.2	116	0.00
53	P	2,4-Dinitrophenol	0.130	0.126	3.1	101	0.00
54		Dibenzofuran	1.565	1.416	9.5	106	0.00
55	P	4-Nitrophenol	0.240	0.206	14.2	95	0.00
56		2,4-Dinitrotoluene	0.387	0.334	13.7	84	0.00
57		Fluorene	1.204	1.190	1.2	111	0.00
58		2,3,4,6-Tetrachlorophenol	0.270	0.259	4.1	94	0.00
59		Diethylphthalate	1.270	1.181	7.0	99	0.00
60		4-Chlorophenyl-phenylether	0.578	0.504	12.8	98	0.00
61		4-Nitroaniline	0.341	0.323	5.3	104	0.00
62		Azobenzene	1.312	1.288	1.8	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.103	0.124	-20.4	104	0.00
65	c	n-Nitrosodiphenylamine	0.639	0.722	-13.0	113	0.00
66		4-Bromophenyl-phenylether	0.209	0.216	-3.3	107	0.00
67		Hexachlorobenzene	0.206	0.216	-4.9	108	0.00
68		Atrazine	0.205	0.211	-2.9	110	0.00
69	C	Pentachlorophenol	0.089	0.104	-16.9	111	0.00
70		Phenanthrene	1.146	1.133	1.1	99	0.00
71		Anthracene	1.151	1.091	5.2	98	-0.01
72		Carbazole	1.100	1.002	8.9	87	0.00
73		Di-n-butylphthalate	1.190	1.325	-11.3	113	0.00
74	C	Fluoranthene	1.182	1.032	12.7	86	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	81	-0.01
76		Benzidine	0.852	0.686	19.5	66	-0.01
77		Pyrene	1.497	1.520	-1.5	78	-0.01
78	S	Terphenyl-d14	0.851	0.899	-5.6	89	-0.01
79		Butylbenzylphthalate	0.608	0.683	-12.3	91	-0.01
80		Benzo(a)anthracene	1.223	1.095	10.5	72	-0.01
81		3,3'-Dichlorobenzidine	0.436	0.431	1.1	78	-0.01
82		Chrysene	1.145	1.066	6.9	70	-0.01
83		Bis(2-ethylhexyl)phthalate	0.831	0.944	-13.6	88	-0.01
84	c	Di-n-octyl phthalate	1.354	1.443	-6.6	82	-0.01
85		Indeno(1,2,3-cd)pyrene	0.887	1.264	-42.5#	121	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	-0.01
87		Benzo(b)fluoranthene	1.316	1.094	16.9	74	-0.01

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88	Benzo(k)fluoranthene	1.137	1.168	-2.7	105	-0.01
89 C	Benzo(a)pyrene	1.139	1.109	2.6	92	-0.01
90	Dibenzo(a,h)anthracene	0.920	1.118	-21.5	121	0.00
91	Benzo(g,h,i)perylene	0.930	1.132	-21.7	122	-0.02

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