Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022017\
Data File : BF093030.D
Acq On : 20 Feb 2017 16:02
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 20 16:49:21 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	0.00
2		1,4-Dioxane	0.630	0.644	-2.2	110	-0.02
3		Pyridine	1.602	1.705	-6.4	110	-0.02
4		n-Nitrosodimethylamine	0.752	0.751	0.1	104	-0.01
5	S	2-Fluorophenol	1.166	1.198	-2.7	103	-0.01
6		Aniline	2.046	2.202	-7.6	116	0.00
7	S	Phenol-d6	1.500	1.530	-2.0	107	0.00
8		2-Chlorophenol	1.286	1.385	-7.7	112	0.00
9		Benzaldehyde	1.075	0.982	8.7	93	0.00
10	C	Phenol	1.755	1.711	2.5	108	-0.01
11		bis(2-Chloroethyl)ether	1.401	1.521	-8.6	119	0.00
12		1,3-Dichlorobenzene	1.506	1.509	-0.2	107	-0.01
13	C	1,4-Dichlorobenzene	1.525	1.565	-2.6	111	-0.01
14		1,2-Dichlorobenzene	1.458	1.480	-1.5	105	0.00
15		Benzyl Alcohol	1.110	1.122	-1.1	109	0.00
16		2,2'-oxybis(1-Chloropropane	2.434	2.616	-7.5	114	-0.01
17		2-Methylphenol	1.103	1.237	-12.1	112	0.00
18		Hexachloroethane	0.520	0.537	-3.3	108	-0.01
19	P	n-Nitroso-di-n-propylamine	0.955	1.020	-6.8	123	0.00
20		3+4-Methylphenols	1.319	1.433	-8.6	111	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.457	0.422	7.7	105	0.00
23	S	Nitrobenzene-d5	0.359	0.360	-0.3	109	0.00
24		Nitrobenzene	0.369	0.348	5.7	111	-0.01
25		Isophorone	0.669	0.675	-0.9	113	-0.01
26	C	2-Nitrophenol	0.175	0.179	-2.3	103	0.00
27		2,4-Dimethylphenol	0.308	0.326	-5.8	111	0.00
28		bis(2-Chloroethoxy)methane	0.443	0.442	0.2	110	0.00
29	C	2,4-Dichlorophenol	0.287	0.285	0.7	114	0.00
30		1,2,4-Trichlorobenzene	0.309	0.305	1.3	111	0.00
31		Naphthalene	1.014	0.988	2.6	110	0.00
32		Benzoic acid	0.156	0.126	19.2	80	0.01
33		4-Chloroaniline	0.398	0.380	4.5	103	0.00
34	C	Hexachlorobutadiene	0.170	0.163	4.1	105	0.00
35		Caprolactam	0.090	0.092	-2.2	105	0.00
36	C	4-Chloro-3-methylphenol	0.299	0.318	-6.4	115	0.00
37		2-Methylnaphthalene	0.651	0.650	0.2	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.561	0.537	4.3	111	0.00
40	P	Hexachlorocyclopentadiene	0.253	0.244	3.6	108	0.00
41	S	2,4,6-Tribromophenol	0.145	0.136	6.2	99	0.00
42	C	2,4,6-Trichlorophenol	0.369	0.361	2.2	105	0.00
43		2,4,5-Trichlorophenol	0.404	0.379	6.2	111	0.00
44	S	2-Fluorobiphenyl	1.104	1.067	3.4	104	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.448	1.447	0.1	110	0.00
46		2-Chloronaphthalene	1.133	1.168	-3.1	111	0.00
47		2-Nitroaniline	0.381	0.374	1.8	120	-0.01
48		Acenaphthylene	1.957	1.798	8.1	113	0.00
49		Dimethylphthalate	1.347	1.240	7.9	104	0.00
50		2,6-Dinitrotoluene	0.291	0.329	-13.1	126	0.00
51	C	Acenaphthene	1.170	1.003	14.3	103	0.00
52		3-Nitroaniline	0.333	0.374	-12.3	120	0.00
53	P	2,4-Dinitrophenol	0.130	0.160	-23.1	133	0.00
54		Dibenzofuran	1.565	1.429	8.7	112	0.00
55	P	4-Nitrophenol	0.240	0.215	10.4	103	0.00
56		2,4-Dinitrotoluene	0.387	0.348	10.1	91	0.00
57		Fluorene	1.204	1.177	2.2	115	0.00
58		2,3,4,6-Tetrachlorophenol	0.270	0.276	-2.2	104	0.00
59		Diethylphthalate	1.270	1.195	5.9	104	0.00
60		4-Chlorophenyl-phenylether	0.578	0.497	14.0	101	0.00
61		4-Nitroaniline	0.341	0.342	-0.3	115	0.00
62		Azobenzene	1.312	1.264	3.7	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
64		4,6-Dinitro-2-methylphenol	0.103	0.136	-32.0#	128	0.00
65	c	n-Nitrosodiphenylamine	0.639	0.664	-3.9	117	0.00
66		4-Bromophenyl-phenylether	0.209	0.200	4.3	111	0.00
67		Hexachlorobenzene	0.206	0.205	0.5	115	0.00
68		Atrazine	0.205	0.195	4.9	115	0.00
69	C	Pentachlorophenol	0.089	0.112	-25.8#	134	0.00
70		Phenanthrene	1.146	1.075	6.2	106	0.00
71		Anthracene	1.151	1.159	-0.7	118	0.00
72		Carbazole	1.100	0.985	10.5	96	0.00
73		Di-n-butylphthalate	1.190	1.266	-6.4	122	0.00
74	C	Fluoranthene	1.182	1.195	-1.1	112	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
76		Benzidine	0.852	0.687	19.4	87	-0.01
77		Pyrene	1.497	1.579	-5.5	107	0.00
78	S	Terphenyl-d14	0.851	0.787	7.5	104	-0.01
79		Butylbenzylphthalate	0.608	0.679	-11.7	120	-0.01
80		Benzo(a)anthracene	1.223	1.211	1.0	105	-0.01
81		3,3'-Dichlorobenzidine	0.436	0.432	0.9	104	-0.01
82		Chrysene	1.145	1.230	-7.4	107	0.00
83		Bis(2-ethylhexyl)phthalate	0.831	0.851	-2.4	106	-0.01
84	c	Di-n-octyl phthalate	1.354	1.487	-9.8	113	0.00
85		Indeno(1,2,3-cd)pyrene	0.887	0.823	7.2	105	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	104	-0.01
87		Benzo(b)fluoranthene	1.316	1.521	-15.6	114	-0.01

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88	Benzo(k)fluoranthene	1.137	0.995	12.5	99	-0.01
89 C	Benzo(a)pyrene	1.139	1.153	-1.2	105	-0.01
90	Dibenzo(a,h)anthracene	0.920	0.864	6.1	103	-0.01
91	Benzo(g,h,i)perylene	0.930	0.881	5.3	105	-0.02

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