Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF050317\
Data File : BF094862.D
Acq On : 4 May 2017 7:13
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 04 07:59:02 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	129	0.00
2		1,4-Dioxane	0.588	0.572	2.7	132	-0.03
3		Pyridine	1.364	1.384	-1.5	134	-0.03
4		n-Nitrosodimethylamine	0.568	0.565	0.5	133	-0.02
5	S	2-Fluorophenol	1.200	1.250	-4.2	135	-0.01
6		Aniline	1.817	1.922	-5.8	130	0.00
7	S	Phenol-d6	1.439	1.486	-3.3	133	0.00
8		2-Chlorophenol	1.365	1.359	0.4	130	0.00
9		Benzaldehyde	0.879	0.857	2.5	124	0.00
10	C	Phenol	1.517	1.571	-3.6	134	0.00
11		bis(2-Chloroethyl)ether	1.252	1.208	3.5	124	0.00
12		1,3-Dichlorobenzene	1.549	1.550	-0.1	139	0.00
13	C	1,4-Dichlorobenzene	1.571	1.528	2.7	125	0.00
14		1,2-Dichlorobenzene	1.466	1.438	1.9	128	0.00
15		Benzyl Alcohol	0.926	0.993	-7.2	132	0.00
16		2,2'-oxybis(1-Chloropropane	1.772	1.749	1.3	131	0.00
17		2-Methylphenol	1.117	1.150	-3.0	138	0.00
18		Hexachloroethane	0.532	0.478	10.2	115	-0.01
19	P	n-Nitroso-di-n-propylamine	0.973	1.011	-3.9	137	0.00
20		3+4-Methylphenols	1.518	1.573	-3.6	134	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	124	0.00
22		Acetophenone	0.480	0.510	-6.3	134	0.00
23	S	Nitrobenzene-d5	0.317	0.333	-5.0	130	0.00
24		Nitrobenzene	0.332	0.339	-2.1	131	0.00
25		Isophorone	0.566	0.586	-3.5	130	0.00
26	C	2-Nitrophenol	0.179	0.182	-1.7	124	0.00
27		2,4-Dimethylphenol	0.290	0.289	0.3	129	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.391	0.3	126	0.00
29	C	2,4-Dichlorophenol	0.274	0.281	-2.6	133	0.00
30		1,2,4-Trichlorobenzene	0.293	0.293	0.0	128	0.00
31		Naphthalene	1.005	0.984	2.1	125	0.00
32		Benzoic acid	0.177	0.188	-6.2	138	0.02
33		4-Chloroaniline	0.375	0.387	-3.2	130	0.00
34	C	Hexachlorobutadiene	0.155	0.144	7.1	119	-0.01
35		Caprolactam	0.082	0.088	-7.3	130	0.02
36	C	4-Chloro-3-methylphenol	0.293	0.299	-2.0	129	0.00
37		2-Methylnaphthalene	0.660	0.646	2.1	125	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	124	0.00
39		1,2,4,5-Tetrachlorobenzene	0.615	0.625	-1.6	119	0.00
40	P	Hexachlorocyclopentadiene	0.221	0.063	71.5#	32#	0.00
41	S	2,4,6-Tribromophenol	0.164	0.165	-0.6	116	0.00
42	C	2,4,6-Trichlorophenol	0.411	0.432	-5.1	124	0.00
43		2,4,5-Trichlorophenol	0.441	0.467	-5.9	124	0.00
44	S	2-Fluorobiphenyl	1.390	1.384	0.4	121	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.715	-1.8	120	0.00
46		2-Chloronaphthalene	1.360	1.342	1.3	119	0.00
47		2-Nitroaniline	0.372	0.395	-6.2	129	0.00
48		Acenaphthylene	2.130	2.107	1.1	119	0.00
49		Dimethylphthalate	1.642	1.730	-5.4	126	-0.01
50		2,6-Dinitrotoluene	0.335	0.344	-2.7	121	0.00
51	C	Acenaphthene	1.272	1.244	2.2	118	0.00
52		3-Nitroaniline	0.360	0.380	-5.6	124	0.00
53	P	2,4-Dinitrophenol	0.156	0.080	48.7#	59	0.00
54		Dibenzofuran	1.760	1.777	-1.0	124	0.00
55	P	4-Nitrophenol	0.216	0.225	-4.2	117	0.00
56		2,4-Dinitrotoluene	0.430	0.423	1.6	115	0.00
57		Fluorene	1.531	1.489	2.7	121	0.00
58		2,3,4,6-Tetrachlorophenol	0.278	0.295	-6.1	129	0.00
59		Diethylphthalate	1.574	1.609	-2.2	127	0.00
60		4-Chlorophenyl-phenylether	0.673	0.651	3.3	116	0.00
61		4-Nitroaniline	0.330	0.331	-0.3	122	0.00
62		Azobenzene	1.265	1.258	0.6	117	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.082	38.8#	67	0.00
65	c	n-Nitrosodiphenylamine	0.726	0.777	-7.0	120	0.00
66		4-Bromophenyl-phenylether	0.211	0.211	0.0	108	0.00
67		Hexachlorobenzene	0.217	0.208	4.1	107	0.00
68		Atrazine	0.205	0.208	-1.5	119	0.00
69	C	Pentachlorophenol	0.085	0.095	-11.8	132	0.00
70		Phenanthrene	1.149	1.126	2.0	111	0.00
71		Anthracene	1.174	1.148	2.2	110	0.00
72		Carbazole	1.068	1.033	3.3	110	0.00
73		Di-n-butylphthalate	1.332	1.436	-7.8	119	0.00
74	C	Fluoranthene	1.179	1.099	6.8	104	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.465	0.489	-5.2	95	0.00
77		Pyrene	1.496	1.554	-3.9	105	0.00
78	S	Terphenyl-d14	0.908	0.929	-2.3	106	0.00
79		Butylbenzylphthalate	0.696	0.756	-8.6	109	0.00
80		Benzo(a)anthracene	1.164	1.137	2.3	100	0.00
81		3,3'-Dichlorobenzidine	0.402	0.409	-1.7	100	-0.01
82		Chrysene	1.125	1.114	1.0	101	-0.01
83		Bis(2-ethylhexyl)phthalate	0.991	0.979	1.2	99	0.00
84	c	Di-n-octyl phthalate	1.625	1.705	-4.9	105	-0.01
85		Indeno(1,2,3-cd)pyrene	0.914	0.892	2.4	102	0.00
86	I	Perylene-d12	1.000	1.000	0.0	108	0.00
87		Benzo(b)fluoranthene	1.236	1.271	-2.8	102	-0.01

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88	Benzo(k)fluoranthene	1.192	1.162	2.5	102	0.00
89 C	Benzo(a)pyrene	1.140	1.091	4.3	98	0.00
90	Dibenzo(a,h)anthracene	0.942	0.912	3.2	102	-0.01
91	Benzo(g,h,i)perylene	0.924	0.865	6.4	101	0.00

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