Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085734.D
Acq On : 25 Mar 2016 1:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 25 02:30:54 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 24 14:15:39 2016
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	98	0.00
2		1,4-Dioxane	0.573	0.553	3.5	95	0.00
3		Pyridine	1.375	1.306	5.0	88	0.01
4		n-Nitrosodimethylamine	0.550	0.567	-3.1	97	0.01
5	S	2-Fluorophenol	1.201	1.263	-5.2	108	0.01
6		Aniline	2.055	2.035	1.0	101	0.00
7	S	Phenol-d6	1.527	1.496	2.0	102	0.00
8		2-Chlorophenol	1.395	1.443	-3.4	101	0.00
9		Benzaldehyde	0.920	0.917	0.3	102	0.00
10	C	Phenol	1.730	1.611	6.9	95	0.01
11		bis(2-Chloroethyl)ether	1.331	1.276	4.1	98	0.00
12		1,3-Dichlorobenzene	1.503	1.529	-1.7	99	0.00
13	C	1,4-Dichlorobenzene	1.524	1.529	-0.3	105	0.00
14		1,2-Dichlorobenzene	1.420	1.468	-3.4	101	0.00
15		Benzyl Alcohol	1.019	1.052	-3.2	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.766	1.813	-2.7	103	0.00
17		2-Methylphenol	1.117	1.128	-1.0	97	0.00
18		Hexachloroethane	0.558	0.573	-2.7	103	0.00
19	P	n-Nitroso-di-n-propylamine	0.914	0.952	-4.2	98	0.00
20		3+4-Methylphenols	1.343	1.379	-2.7	104	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22		Acetophenone	0.466	0.465	0.2	104	0.00
23	S	Nitrobenzene-d5	0.355	0.358	-0.8	100	0.00
24		Nitrobenzene	0.366	0.362	1.1	105	0.01
25		Isophorone	0.684	0.698	-2.0	105	0.00
26	C	2-Nitrophenol	0.183	0.182	0.5	101	0.00
27		2,4-Dimethylphenol	0.320	0.352	-10.0	103	0.00
28		bis(2-Chloroethoxy)methane	0.390	0.400	-2.6	99	0.00
29	C	2,4-Dichlorophenol	0.269	0.268	0.4	102	0.01
30		1,2,4-Trichlorobenzene	0.299	0.304	-1.7	103	0.00
31		Naphthalene	1.008	1.000	0.8	104	0.00
32		Benzoic acid	0.213	0.245	-15.0	102	0.00
33		4-Chloroaniline	0.411	0.412	-0.2	101	0.00
34	C	Hexachlorobutadiene	0.162	0.170	-4.9	106	0.00
35		Caprolactam	0.095	0.105	-10.5	116	0.01
36	C	4-Chloro-3-methylphenol	0.314	0.323	-2.9	109	0.01
37		2-Methylnaphthalene	0.635	0.637	-0.3	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	0.593	0.601	-1.3	102	0.00
40	P	Hexachlorocyclopentadiene	0.196	0.193	1.5	99	-0.01
41	S	2,4,6-Tribromophenol	0.190	0.193	-1.6	102	0.00
42	C	2,4,6-Trichlorophenol	0.401	0.401	0.0	104	0.00
43		2,4,5-Trichlorophenol	0.420	0.408	2.9	100	0.00
44	S	2-Fluorobiphenyl	1.197	1.206	-0.8	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.688	1.752	-3.8	106	0.00
46		2-Chloronaphthalene	1.241	1.321	-6.4	104	0.00
47		2-Nitroaniline	0.379	0.356	6.1	103	0.00
48		Acenaphthylene	2.022	2.050	-1.4	101	0.00
49		Dimethylphthalate	1.517	1.439	5.1	103	0.00
50		2,6-Dinitrotoluene	0.327	0.359	-9.8	108	0.00
51	C	Acenaphthene	1.235	1.160	6.1	102	0.00
52		3-Nitroaniline	0.374	0.379	-1.3	111	0.00
53	P	2,4-Dinitrophenol	0.140	0.138	1.4	91	0.00
54		Dibenzofuran	1.599	1.592	0.4	102	0.00
55	P	4-Nitrophenol	0.244	0.244	0.0	98	0.00
56		2,4-Dinitrotoluene	0.415	0.430	-3.6	97	0.00
57		Fluorene	1.329	1.362	-2.5	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.298	-2.1	107	0.00
59		Diethylphthalate	1.499	1.442	3.8	102	0.00
60		4-Chlorophenyl-phenylether	0.650	0.589	9.4	102	0.00
61		4-Nitroaniline	0.358	0.326	8.9	88	0.00
62		Azobenzene	1.441	1.456	-1.0	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.117	0.120	-2.6	96	0.00
65	c	n-Nitrosodiphenylamine	0.633	0.653	-3.2	103	0.00
66		4-Bromophenyl-phenylether	0.205	0.209	-2.0	103	0.00
67		Hexachlorobenzene	0.214	0.230	-7.5	104	0.00
68		Atrazine	0.214	0.205	4.2	99	0.00
69	C	Pentachlorophenol	0.105	0.110	-4.8	97	0.00
70		Phenanthrene	1.027	1.120	-9.1	103	0.00
71		Anthracene	1.078	1.068	0.9	106	0.00
72		Carbazole	0.905	0.912	-0.8	97	0.00
73		Di-n-butylphthalate	1.242	1.250	-0.6	105	0.00
74	C	Fluoranthene	1.092	1.071	1.9	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
76		Benzidine	0.704	0.727	-3.3	105	0.00
77		Pyrene	1.644	1.506	8.4	105	0.00
78	S	Terphenyl-d14	0.972	0.865	11.0	103	0.00
79		Butylbenzylphthalate	0.688	0.682	0.9	105	0.00
80		Benzo(a)anthracene	1.140	1.103	3.2	104	0.00
81		3,3'-Dichlorobenzidine	0.791	0.712	10.0	103	0.00
82		Chrysene	1.149	1.097	4.5	105	0.00
83		Bis(2-ethylhexyl)phthalate	0.824	0.824	0.0	104	0.00
84	c	Di-n-octyl phthalate	1.226	1.247	-1.7	104	0.00
85		Indeno(1,2,3-cd)pyrene	0.887	0.847	4.5	100	0.00
86	I	Perylene-d12	1.000	1.000	0.0	100	0.00
87		Benzo(b)fluoranthene	1.260	1.373	-9.0	107	0.00

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88	Benzo(k)fluoranthene	1.136	1.101	3.1	104	0.00
89 C	Benzo(a)pyrene	1.119	1.149	-2.7	104	0.00
90	Dibenzo(a,h)anthracene	1.039	1.012	2.6	100	0.00
91	Benzo(g,h,i)perylene	1.052	1.024	2.7	99	0.00

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