Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060516\
Data File : BF087728.D
Acq On : 6 Jun 2016 1:22
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 06 04:41:39 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jun 04 11:07:28 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	106	-0.01
2		1,4-Dioxane	0.598	0.608	-1.7	105	0.00
3		Pyridine	1.580	1.360	13.9	88	0.00
4		n-Nitrosodimethylamine	0.668	0.652	2.4	101	0.00
5	S	2-Fluorophenol	1.237	1.165	5.8	96	0.00
6		Aniline	2.200	2.141	2.7	102	0.00
7	S	Phenol-d6	1.552	1.440	7.2	95	0.00
8		2-Chlorophenol	1.424	1.360	4.5	106	0.00
9		Benzaldehyde	1.049	0.959	8.6	93	0.00
10	C	Phenol	1.767	1.684	4.7	92	0.00
11		bis(2-Chloroethyl)ether	1.284	1.265	1.5	114	0.01
12		1,3-Dichlorobenzene	1.524	1.409	7.5	105	0.00
13	C	1,4-Dichlorobenzene	1.530	1.519	0.7	107	0.00
14		1,2-Dichlorobenzene	1.434	1.313	8.4	101	0.00
15		Benzyl Alcohol	1.146	1.077	6.0	95	0.00
16		2,2'-oxybis(1-Chloropropane	1.872	1.729	7.6	100	0.00
17		2-Methylphenol	1.143	1.156	-1.1	107	0.01
18		Hexachloroethane	0.535	0.527	1.5	102	0.00
19	P	n-Nitroso-di-n-propylamine	0.969	0.892	7.9	106	0.00
20		3+4-Methylphenols	1.401	1.349	3.7	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.453	0.464	-2.4	102	0.00
23	S	Nitrobenzene-d5	0.346	0.328	5.2	102	0.00
24		Nitrobenzene	0.346	0.378	-9.2	106	0.00
25		Isophorone	0.629	0.622	1.1	105	0.00
26	C	2-Nitrophenol	0.161	0.165	-2.5	112	0.00
27		2,4-Dimethylphenol	0.333	0.344	-3.3	105	0.00
28		bis(2-Chloroethoxy)methane	0.399	0.379	5.0	103	0.00
29	C	2,4-Dichlorophenol	0.274	0.267	2.6	105	0.00
30		1,2,4-Trichlorobenzene	0.287	0.297	-3.5	109	0.00
31		Naphthalene	0.972	0.970	0.2	106	0.00
32		Benzoic acid	0.201	0.151	24.9	71	0.01
33		4-Chloroaniline	0.422	0.400	5.2	103	0.00
34	C	Hexachlorobutadiene	0.149	0.150	-0.7	103	0.00
35		Caprolactam	0.090	0.098	-8.9	116	0.01
36	C	4-Chloro-3-methylphenol	0.283	0.271	4.2	97	0.00
37		2-Methylnaphthalene	0.642	0.619	3.6	103	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
39		1,2,4,5-Tetrachlorobenzene	0.551	0.556	-0.9	108	0.00
40	P	Hexachlorocyclopentadiene	0.324	0.284	12.3	98	0.00
41	S	2,4,6-Tribromophenol	0.201	0.196	2.5	110	0.01
42	C	2,4,6-Trichlorophenol	0.416	0.413	0.7	107	0.00
43		2,4,5-Trichlorophenol	0.387	0.411	-6.2	106	0.00
44	S	2-Fluorobiphenyl	1.285	1.334	-3.8	105	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.623	1.471	9.4	105	0.00
46		2-Chloronaphthalene	1.292	1.142	11.6	104	0.00
47		2-Nitroaniline	0.364	0.352	3.3	110	0.00
48		Acenaphthylene	2.015	1.974	2.0	107	0.01
49		Dimethylphthalate	1.454	1.460	-0.4	103	0.00
50		2,6-Dinitrotoluene	0.331	0.358	-8.2	106	0.00
51	C	Acenaphthene	1.296	1.247	3.8	104	0.00
52		3-Nitroaniline	0.378	0.376	0.5	120	0.01
53	P	2,4-Dinitrophenol	0.134	0.114	14.9	81	0.00
54		Dibenzofuran	1.765	1.762	0.2	104	0.00
55	P	4-Nitrophenol	0.323	0.312	3.4	123	0.01
56		2,4-Dinitrotoluene	0.421	0.472	-12.1	108	0.00
57		Fluorene	1.380	1.244	9.9	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.334	0.327	2.1	110	0.00
59		Diethylphthalate	1.460	1.411	3.4	107	0.00
60		4-Chlorophenyl-phenylether	0.605	0.548	9.4	106	0.00
61		4-Nitroaniline	0.364	0.354	2.7	95	0.00
62		Azobenzene	1.470	1.485	-1.0	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
64		4,6-Dinitro-2-methylphenol	0.091	0.094	-3.3	89	0.00
65	c	n-Nitrosodiphenylamine	0.655	0.650	0.8	106	0.00
66		4-Bromophenyl-phenylether	0.198	0.216	-9.1	104	0.00
67		Hexachlorobenzene	0.222	0.216	2.7	106	0.00
68		Atrazine	0.194	0.203	-4.6	103	0.00
69	C	Pentachlorophenol	0.132	0.125	5.3	88	0.00
70		Phenanthrene	1.084	1.024	5.5	106	0.00
71		Anthracene	1.087	1.126	-3.6	101	0.00
72		Carbazole	1.029	0.947	8.0	101	0.00
73		Di-n-butylphthalate	1.103	1.221	-10.7	107	0.00
74	C	Fluoranthene	1.067	1.068	-0.1	101	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	86	0.00
76		Benzidine	0.789	0.712	9.8	67	0.00
77		Pyrene	1.424	1.661	-16.6	99	0.00
78	S	Terphenyl-d14	0.820	0.928	-13.2	105	0.00
79		Butylbenzylphthalate	0.606	0.756	-24.8	99	0.00
80		Benzo(a)anthracene	1.154	1.191	-3.2	88	-0.01
81		3,3'-Dichlorobenzidine	0.326	0.355	-8.9	85	0.00
82		Chrysene	1.149	1.266	-10.2	95	0.00
83		Bis(2-ethylhexyl)phthalate	0.721	0.936	-29.8#	104	0.00
84	c	Di-n-octyl phthalate	1.341	1.504	-12.2	94	-0.01
85		Indeno(1,2,3-cd)pyrene	0.950	1.123	-18.2	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	93	0.00
87		Benzo(b)fluoranthene	1.301	1.367	-5.1	95	0.00

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88	Benzo(k)fluoranthene	1.082	0.984	9.1	89	0.00
89 C	Benzo(a)pyrene	1.094	1.111	-1.6	92	0.00
90	Dibenzo(a,h)anthracene	0.931	1.011	-8.6	98	0.00
91	Benzo(g,h,i)perylene	0.939	1.016	-8.2	98	0.00

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