Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
Data File : BF085998.D
Acq On : 2 Apr 2016 8:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 04 01:43:18 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 01 23:17:06 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	102	0.00
2		1,4-Dioxane	0.555	0.551	0.7	106	0.00
3		Pyridine	1.242	1.349	-8.6	100	0.01
4		n-Nitrosodimethylamine	0.546	0.551	-0.9	102	0.00
5	S	2-Fluorophenol	1.170	1.176	-0.5	102	0.00
6		Aniline	1.950	1.955	-0.3	101	0.00
7	S	Phenol-d6	1.486	1.529	-2.9	103	0.00
8		2-Chlorophenol	1.395	1.371	1.7	103	0.00
9		Benzaldehyde	0.800	0.760	5.0	97	0.00
10	C	Phenol	1.695	1.801	-6.3	101	0.00
11		bis(2-Chloroethyl)ether	1.342	1.219	9.2	104	0.00
12		1,3-Dichlorobenzene	1.479	1.436	2.9	102	0.00
13	C	1,4-Dichlorobenzene	1.524	1.452	4.7	101	-0.01
14		1,2-Dichlorobenzene	1.402	1.390	0.9	100	0.00
15		Benzyl Alcohol	0.962	0.929	3.4	102	0.00
16		2,2'-oxybis(1-Chloropropane	1.640	1.619	1.3	102	0.00
17		2-Methylphenol	1.100	1.048	4.7	103	0.00
18		Hexachloroethane	0.560	0.543	3.0	103	-0.01
19	P	n-Nitroso-di-n-propylamine	0.921	0.871	5.4	105	0.00
20		3+4-Methylphenols	1.338	1.304	2.5	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.470	0.464	1.3	105	0.00
23	S	Nitrobenzene-d5	0.339	0.329	2.9	106	0.00
24		Nitrobenzene	0.371	0.362	2.4	101	0.00
25		Isophorone	0.669	0.654	2.2	102	0.00
26	C	2-Nitrophenol	0.178	0.176	1.1	104	0.00
27		2,4-Dimethylphenol	0.319	0.296	7.2	104	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.366	6.9	105	0.00
29	C	2,4-Dichlorophenol	0.270	0.261	3.3	102	0.00
30		1,2,4-Trichlorobenzene	0.303	0.293	3.3	103	0.00
31		Naphthalene	1.006	0.931	7.5	99	-0.01
32		Benzoic acid	0.234	0.225	3.8	96	0.00
33		4-Chloroaniline	0.406	0.386	4.9	103	0.00
34	C	Hexachlorobutadiene	0.163	0.159	2.5	103	0.00
35		Caprolactam	0.086	0.097	-12.8	108	0.01
36	C	4-Chloro-3-methylphenol	0.303	0.303	0.0	102	0.00
37		2-Methylnaphthalene	0.657	0.617	6.1	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	0.601	0.559	7.0	99	0.00
40	P	Hexachlorocyclopentadiene	0.163	0.170	-4.3	108	0.00
41	S	2,4,6-Tribromophenol	0.188	0.193	-2.7	107	0.00
42	C	2,4,6-Trichlorophenol	0.386	0.377	2.3	104	0.00
43		2,4,5-Trichlorophenol	0.392	0.379	3.3	104	0.00
44	S	2-Fluorobiphenyl	1.203	1.255	-4.3	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.724	1.622	5.9	102	0.00
46		2-Chloronaphthalene	1.250	1.234	1.3	103	0.00
47		2-Nitroaniline	0.366	0.363	0.8	101	0.00
48		Acenaphthylene	2.002	2.041	-1.9	104	0.00
49		Dimethylphthalate	1.482	1.397	5.7	108	0.00
50		2,6-Dinitrotoluene	0.314	0.336	-7.0	105	0.00
51	C	Acenaphthene	1.191	1.194	-0.3	104	0.00
52		3-Nitroaniline	0.378	0.399	-5.6	104	0.00
53	P	2,4-Dinitrophenol	0.127	0.148	-16.5	101	0.00
54		Dibenzofuran	1.573	1.583	-0.6	105	0.00
55	P	4-Nitrophenol	0.223	0.215	3.6	103	0.00
56		2,4-Dinitrotoluene	0.396	0.377	4.8	106	0.00
57		Fluorene	1.344	1.366	-1.6	104	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.288	1.4	106	0.00
59		Diethylphthalate	1.496	1.457	2.6	111	0.00
60		4-Chlorophenyl-phenylether	0.626	0.635	-1.4	108	0.00
61		4-Nitroaniline	0.372	0.395	-6.2	113	0.01
62		Azobenzene	1.433	1.464	-2.2	109	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	-0.01
64		4,6-Dinitro-2-methylphenol	0.121	0.121	0.0	109	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.633	-1.3	107	0.00
66		4-Bromophenyl-phenylether	0.204	0.208	-2.0	107	0.00
67		Hexachlorobenzene	0.211	0.220	-4.3	109	0.00
68		Atrazine	0.202	0.194	4.0	110	0.00
69	C	Pentachlorophenol	0.099	0.100	-1.0	104	0.00
70		Phenanthrene	1.091	1.080	1.0	105	0.00
71		Anthracene	1.143	1.014	11.3	106	0.00
72		Carbazole	0.982	0.923	6.0	108	0.00
73		Di-n-butylphthalate	1.217	1.265	-3.9	110	0.00
74	C	Fluoranthene	1.135	1.060	6.6	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
76		Benzidine	0.706	0.646	8.5	102	0.00
77		Pyrene	1.409	1.267	10.1	106	0.00
78	S	Terphenyl-d14	0.851	0.809	4.9	115	0.00
79		Butylbenzylphthalate	0.635	0.630	0.8	112	-0.01
80		Benzo(a)anthracene	1.179	1.106	6.2	110	0.00
81		3,3'-Dichlorobenzidine	0.861	0.850	1.3	105	-0.01
82		Chrysene	1.136	1.160	-2.1	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.842	0.837	0.6	113	0.00
84	c	Di-n-octyl phthalate	1.345	1.473	-9.5	115	0.00
85		Indeno(1,2,3-cd)pyrene	0.921	0.904	1.8	112	0.00
86	I	Perylene-d12	1.000	1.000	0.0	114	0.00
87		Benzo(b)fluoranthene	1.258	1.476	-17.3	119	-0.01

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88	Benzo(k)fluoranthene	1.114	0.847	24.0	104	0.00
89 C	Benzo(a)pyrene	1.115	1.060	4.9	113	0.00
90	Dibenzo(a,h)anthracene	0.897	0.862	3.9	113	-0.01
91	Benzo(g,h,i)perylene	0.868	0.808	6.9	111	-0.01

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