Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031816\
Data File : BF085546.D
Acq On : 19 Mar 2016 1:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 19 03:38:16 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 23:31:39 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	102	0.00
2		1,4-Dioxane	0.579	0.582	-0.5	100	-0.01
3		Pyridine	1.421	1.473	-3.7	105	-0.01
4		n-Nitrosodimethylamine	0.593	0.588	0.8	102	-0.01
5	S	2-Fluorophenol	1.191	1.243	-4.4	101	0.00
6		Aniline	2.051	2.062	-0.5	102	0.00
7	S	Phenol-d6	1.527	1.579	-3.4	103	0.00
8		2-Chlorophenol	1.377	1.450	-5.3	105	0.00
9		Benzaldehyde	0.768	0.718	6.5	102	0.00
10	C	Phenol	1.739	1.935	-11.3	103	0.00
11		bis(2-Chloroethyl)ether	1.311	1.398	-6.6	104	0.00
12		1,3-Dichlorobenzene	1.501	1.489	0.8	102	-0.01
13	C	1,4-Dichlorobenzene	1.531	1.455	5.0	102	0.00
14		1,2-Dichlorobenzene	1.359	1.436	-5.7	101	0.00
15		Benzyl Alcohol	1.046	1.004	4.0	100	0.00
16		2,2'-oxybis(1-Chloropropane	1.725	1.783	-3.4	101	0.00
17		2-Methylphenol	1.142	1.128	1.2	101	0.00
18		Hexachloroethane	0.553	0.563	-1.8	103	0.00
19	P	n-Nitroso-di-n-propylamine	0.912	0.928	-1.8	100	0.00
20		3+4-Methylphenols	1.323	1.281	3.2	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.475	0.431	9.3	102	0.00
23	S	Nitrobenzene-d5	0.345	0.349	-1.2	100	-0.01
24		Nitrobenzene	0.377	0.360	4.5	104	0.00
25		Isophorone	0.690	0.670	2.9	103	0.00
26	C	2-Nitrophenol	0.186	0.210	-12.9	104	0.00
27		2,4-Dimethylphenol	0.330	0.306	7.3	102	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.406	-3.6	101	0.00
29	C	2,4-Dichlorophenol	0.272	0.285	-4.8	105	0.00
30		1,2,4-Trichlorobenzene	0.305	0.288	5.6	101	0.00
31		Naphthalene	1.010	0.903	10.6	99	0.00
32		Benzoic acid	0.276	0.273	1.1	101	0.00
33		4-Chloroaniline	0.414	0.433	-4.6	102	0.00
34	C	Hexachlorobutadiene	0.162	0.162	0.0	100	0.00
35		Caprolactam	0.094	0.095	-1.1	104	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.336	-5.7	106	0.00
37		2-Methylnaphthalene	0.615	0.651	-5.9	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	0.607	0.542	10.7	104	0.00
40	P	Hexachlorocyclopentadiene	0.273	0.284	-4.0	105	0.00
41	S	2,4,6-Tribromophenol	0.196	0.211	-7.7	113	0.00
42	C	2,4,6-Trichlorophenol	0.415	0.414	0.2	108	0.00
43		2,4,5-Trichlorophenol	0.423	0.412	2.6	101	-0.01
44	S	2-Fluorobiphenyl	1.250	1.260	-0.8	103	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.725	1.539	10.8	98	0.00
46		2-Chloronaphthalene	1.260	1.246	1.1	102	0.00
47		2-Nitroaniline	0.381	0.425	-11.5	109	0.00
48		Acenaphthylene	2.040	2.128	-4.3	106	0.00
49		Dimethylphthalate	1.509	1.461	3.2	105	0.00
50		2,6-Dinitrotoluene	0.319	0.303	5.0	107	0.00
51	C	Acenaphthene	1.280	1.326	-3.6	107	0.00
52		3-Nitroaniline	0.399	0.398	0.3	112	0.00
53	P	2,4-Dinitrophenol	0.195	0.191	2.1	107	0.00
54		Dibenzofuran	1.559	1.637	-5.0	107	0.00
55	P	4-Nitrophenol	0.308	0.331	-7.5	107	0.00
56		2,4-Dinitrotoluene	0.417	0.454	-8.9	106	0.00
57		Fluorene	1.299	1.357	-4.5	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.304	0.308	-1.3	110	0.00
59		Diethylphthalate	1.501	1.508	-0.5	108	0.00
60		4-Chlorophenyl-phenylether	0.634	0.610	3.8	110	0.00
61		4-Nitroaniline	0.393	0.431	-9.7	105	0.00
62		Azobenzene	1.440	1.455	-1.0	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.01
64		4,6-Dinitro-2-methylphenol	0.128	0.132	-3.1	107	0.00
65	c	n-Nitrosodiphenylamine	0.623	0.624	-0.2	101	0.00
66		4-Bromophenyl-phenylether	0.200	0.215	-7.5	108	0.00
67		Hexachlorobenzene	0.212	0.232	-9.4	106	0.00
68		Atrazine	0.190	0.214	-12.6	109	0.00
69	C	Pentachlorophenol	0.129	0.135	-4.7	104	-0.01
70		Phenanthrene	1.055	1.116	-5.8	106	0.00
71		Anthracene	1.106	1.006	9.0	105	0.00
72		Carbazole	0.954	1.043	-9.3	109	0.00
73		Di-n-butylphthalate	1.195	1.307	-9.4	110	0.00
74	C	Fluoranthene	1.104	1.043	5.5	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
76		Benzidine	0.673	0.676	-0.4	111	0.00
77		Pyrene	1.436	1.339	6.8	101	-0.01
78	S	Terphenyl-d14	0.831	0.745	10.3	110	0.00
79		Butylbenzylphthalate	0.685	0.653	4.7	104	-0.01
80		Benzo(a)anthracene	1.161	1.137	2.1	112	0.00
81		3,3'-Dichlorobenzidine	0.426	0.415	2.6	111	0.00
82		Chrysene	1.117	1.107	0.9	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.850	0.801	5.8	106	0.00
84	c	Di-n-octyl phthalate	1.386	1.505	-8.6	118	0.00
85		Indeno(1,2,3-cd)pyrene	0.933	0.895	4.1	115	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	118	0.00
87		Benzo(b)fluoranthene	1.305	1.173	10.1	118	0.00

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88	Benzo(k)fluoranthene	1.075	1.223	-13.8	112	0.00
89 C	Benzo(a)pyrene	1.106	1.122	-1.4	118	0.00
90	Dibenzo(a,h)anthracene	0.923	0.861	6.7	115	0.00
91	Benzo(g,h,i)perylene	0.893	0.808	9.5	113	0.00

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