Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080615\
Data File : BF080904.D
Acq On : 7 Aug 2015 00:08
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 07 07:19:11 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 07 03:17:12 2015
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	100	0.00
2		1,4-Dioxane	0.608	0.597	1.8	97	0.00
3		Pyridine	1.709	1.828	-7.0	97	-0.01
4		n-Nitrosodimethylamine	0.884	0.908	-2.7	97	0.00
5	S	2-Fluorophenol	1.255	1.280	-2.0	98	0.00
6		Aniline	2.562	2.435	5.0	95	0.00
7	S	Phenol-d6	1.727	1.773	-2.7	100	0.00
8		2-Chlorophenol	1.425	1.512	-6.1	101	0.00
9		Benzaldehyde	1.190	1.190	0.0	100	0.00
10	C	Phenol	2.072	2.261	-9.1	103	0.00
11		bis(2-Chloroethyl)ether	1.467	1.407	4.1	102	0.00
12		1,3-Dichlorobenzene	1.545	1.547	-0.1	96	0.00
13	C	1,4-Dichlorobenzene	1.573	1.530	2.7	91	0.00
14		1,2-Dichlorobenzene	1.442	1.433	0.6	100	0.00
15		Benzyl Alcohol	1.370	1.460	-6.6	95	0.00
16		2,2'-oxybis(1-Chloropropane	3.011	2.882	4.3	99	0.00
17		2-Methylphenol	1.277	1.304	-2.1	97	0.00
18		Hexachloroethane	0.598	0.594	0.7	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.230	1.234	-0.3	98	0.00
20		3+4-Methylphenols	1.548	1.434	7.4	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.500	0.454	9.2	92	0.00
23	S	Nitrobenzene-d5	0.425	0.439	-3.3	99	0.00
24		Nitrobenzene	0.433	0.388	10.4	94	0.00
25		Isophorone	0.814	0.799	1.8	96	0.00
26	C	2-Nitrophenol	0.182	0.189	-3.8	104	0.00
27		2,4-Dimethylphenol	0.348	0.353	-1.4	92	0.00
28		bis(2-Chloroethoxy)methane	0.482	0.423	12.2	96	0.00
29	C	2,4-Dichlorophenol	0.278	0.268	3.6	98	0.00
30		1,2,4-Trichlorobenzene	0.308	0.315	-2.3	97	0.00
31		Naphthalene	1.085	1.102	-1.6	96	0.00
32		Benzoic acid	0.222	0.240	-8.1	100	0.00
33		4-Chloroaniline	0.449	0.464	-3.3	101	0.00
34	C	Hexachlorobutadiene	0.175	0.168	4.0	95	0.00
35		Caprolactam	0.104	0.109	-4.8	96	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.341	-1.8	100	0.00
37		2-Methylnaphthalene	0.669	0.687	-2.7	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	0.609	0.617	-1.3	95	0.00
40	P	Hexachlorocyclopentadiene	0.350	0.364	-4.0	93	0.00
41	S	2,4,6-Tribromophenol	0.215	0.240	-11.6	101	0.00
42	C	2,4,6-Trichlorophenol	0.449	0.410	8.7	94	0.00
43		2,4,5-Trichlorophenol	0.456	0.451	1.1	95	0.00
44	S	2-Fluorobiphenyl	1.403	1.427	-1.7	106	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.692	1.790	-5.8	97	0.00
46		2-Chloronaphthalene	1.323	1.391	-5.1	97	0.00
47		2-Nitroaniline	0.538	0.595	-10.6	104	0.00
48		Acenaphthylene	2.320	2.446	-5.4	101	0.00
49		Dimethylphthalate	1.588	1.471	7.4	93	0.00
50		2,6-Dinitrotoluene	0.341	0.322	5.6	96	0.00
51	C	Acenaphthene	1.410	1.485	-5.3	103	0.00
52		3-Nitroaniline	0.421	0.441	-4.8	106	0.00
53	P	2,4-Dinitrophenol	0.187	0.213	-13.9	106	0.00
54		Dibenzofuran	1.910	1.998	-4.6	101	0.00
55	P	4-Nitrophenol	0.318	0.345	-8.5	93	0.00
56		2,4-Dinitrotoluene	0.458	0.418	8.7	91	0.00
57		Fluorene	1.494	1.607	-7.6	99	0.00
58		2,3,4,6-Tetrachlorophenol	0.364	0.359	1.4	100	0.00
59		Diethylphthalate	1.730	1.638	5.3	95	0.00
60		4-Chlorophenyl-phenylether	0.700	0.638	8.9	95	0.00
61		4-Nitroaniline	0.432	0.449	-3.9	95	-0.01
62		Azobenzene	1.896	1.888	0.4	102	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.130	0.144	-10.8	103	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.749	-5.0	100	0.00
66		4-Bromophenyl-phenylether	0.217	0.212	2.3	100	0.00
67		Hexachlorobenzene	0.235	0.239	-1.7	96	0.00
68		Atrazine	0.206	0.211	-2.4	103	0.00
69	C	Pentachlorophenol	0.132	0.134	-1.5	99	0.00
70		Phenanthrene	1.170	1.140	2.6	95	0.00
71		Anthracene	1.152	1.144	0.7	106	0.00
72		Carbazole	1.163	1.131	2.8	92	0.00
73		Di-n-butylphthalate	1.418	1.367	3.6	101	0.00
74	C	Fluoranthene	1.280	1.249	2.4	104	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
76		Benzidine	0.817	0.908	-11.1	101	0.00
77		Pyrene	1.487	1.576	-6.0	106	0.00
78	S	Terphenyl-d14	0.949	0.978	-3.1	105	0.00
79		Butylbenzylphthalate	0.711	0.771	-8.4	95	0.00
80		Benzo(a)anthracene	1.307	1.336	-2.2	100	0.00
81		3,3'-Dichlorobenzidine	0.489	0.560	-14.5	101	0.00
82		Chrysene	1.249	1.357	-8.6	106	0.00
83		Bis(2-ethylhexyl)phthalate	0.974	1.039	-6.7	99	0.00
84	c	Di-n-octyl phthalate	1.717	1.861	-8.4	102	0.00
85		Indeno(1,2,3-cd)pyrene	1.216	1.347	-10.8	104	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	101	0.00
87		Benzo(b)fluoranthene	1.405	1.512	-7.6	108	0.00

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88	Benzo(k)fluoranthene	1.183	1.021	13.7	91	0.00
89 C	Benzo(a)pyrene	1.210	1.202	0.7	97	-0.01
90	Dibenzo(a,h)anthracene	1.043	1.080	-3.5	103	0.00
91	Benzo(g,h,i)perylene	1.011	1.040	-2.9	102	0.00

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