Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF080315\
Data File : BF080827.D
Acq On : 4 Aug 2015 6:40
Operator : TP
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 04 15:32:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 14:50:50 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	103	-0.01
2		1,4-Dioxane	0.617	0.561	9.1	100	-0.01
3		Pyridine	1.626	1.626	0.0	108	-0.01
4		n-Nitrosodimethylamine	0.903	0.853	5.5	103	-0.01
5	S	2-Fluorophenol	1.205	1.241	-3.0	101	0.00
6		Aniline	2.316	2.231	3.7	105	0.00
7	S	Phenol-d6	1.620	1.542	4.8	104	0.00
8		2-Chlorophenol	1.327	1.263	4.8	103	0.00
9		Benzaldehyde	0.999	0.948	5.1	103	0.00
10	C	Phenol	1.885	1.597	15.3	105	0.00
11		bis(2-Chloroethyl)ether	1.355	1.383	-2.1	104	0.00
12		1,3-Dichlorobenzene	1.490	1.364	8.5	104	0.00
13	C	1,4-Dichlorobenzene	1.486	1.471	1.0	106	0.00
14		1,2-Dichlorobenzene	1.376	1.239	10.0	103	-0.01
15		Benzyl Alcohol	1.104	1.060	4.0	100	0.00
16		2,2'-oxybis(1-Chloropropane	2.678	2.521	5.9	105	0.00
17		2-Methylphenol	1.171	1.167	0.3	105	0.00
18		Hexachloroethane	0.548	0.521	4.9	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.046	0.965	7.7	109	0.00
20		3+4-Methylphenols	1.360	1.244	8.5	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.488	0.475	2.7	106	0.00
23	S	Nitrobenzene-d5	0.407	0.359	11.8	107	0.00
24		Nitrobenzene	0.414	0.394	4.8	102	0.00
25		Isophorone	0.728	0.705	3.2	108	0.00
26	C	2-Nitrophenol	0.177	0.166	6.2	106	0.00
27		2,4-Dimethylphenol	0.337	0.333	1.2	108	0.00
28		bis(2-Chloroethoxy)methane	0.439	0.378	13.9	104	0.00
29	C	2,4-Dichlorophenol	0.283	0.261	7.8	108	0.00
30		1,2,4-Trichlorobenzene	0.308	0.299	2.9	105	0.00
31		Naphthalene	0.984	0.952	3.3	103	0.00
32		Benzoic acid	0.204	0.205	-0.5	100	0.01
33		4-Chloroaniline	0.414	0.414	0.0	109	0.00
34	C	Hexachlorobutadiene	0.189	0.173	8.5	106	0.00
35		Caprolactam	0.079	0.084	-6.3	111	0.00
36	C	4-Chloro-3-methylphenol	0.304	0.298	2.0	103	0.00
37		2-Methylnaphthalene	0.617	0.584	5.3	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	0.637	0.617	3.1	104	0.00
40	P	Hexachlorocyclopentadiene	0.387	0.381	1.6	100	0.00
41	S	2,4,6-Tribromophenol	0.192	0.211	-9.9	107	0.00
42	C	2,4,6-Trichlorophenol	0.451	0.427	5.3	104	0.00
43		2,4,5-Trichlorophenol	0.442	0.435	1.6	109	0.00
44	S	2-Fluorobiphenyl	1.329	1.222	8.1	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.514	1.585	-4.7	108	0.00
46		2-Chloronaphthalene	1.253	1.266	-1.0	104	0.00
47		2-Nitroaniline	0.464	0.501	-8.0	110	0.00
48		Acenaphthylene	1.962	1.971	-0.5	107	0.00
49		Dimethylphthalate	1.333	1.225	8.1	110	0.00
50		2,6-Dinitrotoluene	0.289	0.269	6.9	108	0.00
51	C	Acenaphthene	1.183	1.159	2.0	104	0.00
52		3-Nitroaniline	0.339	0.348	-2.7	104	0.00
53	P	2,4-Dinitrophenol	0.157	0.170	-8.3	107	0.00
54		Dibenzofuran	1.560	1.573	-0.8	108	0.00
55	P	4-Nitrophenol	0.238	0.219	8.0	103	0.00
56		2,4-Dinitrotoluene	0.364	0.333	8.5	108	0.00
57		Fluorene	1.212	1.224	-1.0	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.317	0.327	-3.2	110	0.00
59		Diethylphthalate	1.281	1.200	6.3	109	0.00
60		4-Chlorophenyl-phenylether	0.599	0.536	10.5	107	0.00
61		4-Nitroaniline	0.308	0.278	9.7	109	0.00
62		Azobenzene	1.483	1.431	3.5	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
64		4,6-Dinitro-2-methylphenol	0.130	0.135	-3.8	108	0.00
65	c	n-Nitrosodiphenylamine	0.697	0.695	0.3	108	0.00
66		4-Bromophenyl-phenylether	0.219	0.204	6.8	109	0.00
67		Hexachlorobenzene	0.248	0.254	-2.4	115	0.00
68		Atrazine	0.160	0.180	-12.5	108	0.00
69	C	Pentachlorophenol	0.133	0.133	0.0	111	0.00
70		Phenanthrene	1.041	1.023	1.7	109	0.00
71		Anthracene	1.032	0.931	9.8	107	0.00
72		Carbazole	0.883	0.859	2.7	106	0.00
73		Di-n-butylphthalate	1.069	0.961	10.1	108	0.00
74	C	Fluoranthene	0.988	0.864	12.6	106	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76		Benzidine	0.590	0.651	-10.3	101	0.00
77		Pyrene	1.484	1.498	-0.9	110	0.00
78	S	Terphenyl-d14	0.999	0.945	5.4	107	0.00
79		Butylbenzylphthalate	0.583	0.618	-6.0	109	0.00
80		Benzo(a)anthracene	1.133	1.139	-0.5	108	0.00
81		3,3'-Dichlorobenzidine	0.403	0.433	-7.4	106	0.00
82		Chrysene	1.108	1.166	-5.2	109	0.00
83		Bis(2-ethylhexyl)phthalate	0.719	0.763	-6.1	111	0.00
84	c	Di-n-octyl phthalate	1.136	1.191	-4.8	109	0.00
85		Indeno(1,2,3-cd)pyrene	1.014	1.198	-18.1	125	0.00
86	I	Perylene-d12	1.000	1.000	0.0	109	0.00
87		Benzo(b)fluoranthene	1.172	1.087	7.3	110	0.00

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88	Benzo(k)fluoranthene	1.023	1.027	-0.4	104	0.00
89 C	Benzo(a)pyrene	1.000	1.021	-2.1	113	0.00
90	Dibenzo(a,h)anthracene	0.970	1.095	-12.9	121	0.00
91	Benzo(g,h,i)perylene	1.014	1.168	-15.2	126	0.00

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