Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073115\
Data File : BF080723.D
Acq On : 31 Jul 2015 8:48
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 01 00:09:29 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 31 01:45:22 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	105	0.00
2		1,4-Dioxane	0.575	0.562	2.3	107	0.01
3		Pyridine	1.595	1.554	2.6	99	0.01
4		n-Nitrosodimethylamine	0.912	0.939	-3.0	106	0.01
5	S	2-Fluorophenol	1.166	1.127	3.3	107	0.00
6		Aniline	2.285	2.237	2.1	103	0.00
7	S	Phenol-d6	1.587	1.606	-1.2	102	0.00
8		2-Chlorophenol	1.307	1.220	6.7	102	0.00
9		Benzaldehyde	1.000	0.956	4.4	103	0.00
10	C	Phenol	1.837	1.983	-7.9	103	0.00
11		bis(2-Chloroethyl)ether	1.312	1.161	11.5	102	0.00
12		1,3-Dichlorobenzene	1.455	1.421	2.3	104	0.00
13	C	1,4-Dichlorobenzene	1.484	1.514	-2.0	105	0.00
14		1,2-Dichlorobenzene	1.357	1.242	8.5	104	0.00
15		Benzyl Alcohol	1.318	1.293	1.9	97	0.00
16		2,2'-oxybis(1-Chloropropane	2.803	2.657	5.2	104	0.00
17		2-Methylphenol	1.073	1.047	2.4	106	0.00
18		Hexachloroethane	0.554	0.537	3.1	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.069	1.007	5.8	102	0.00
20		3+4-Methylphenols	1.323	1.296	2.0	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.483	0.464	3.9	109	0.00
23	S	Nitrobenzene-d5	0.413	0.422	-2.2	104	0.00
24		Nitrobenzene	0.414	0.377	8.9	107	0.00
25		Isophorone	0.734	0.717	2.3	105	0.01
26	C	2-Nitrophenol	0.168	0.184	-9.5	105	0.00
27		2,4-Dimethylphenol	0.316	0.294	7.0	103	0.00
28		bis(2-Chloroethoxy)methane	0.435	0.432	0.7	102	0.00
29	C	2,4-Dichlorophenol	0.280	0.289	-3.2	106	0.00
30		1,2,4-Trichlorobenzene	0.306	0.292	4.6	101	0.00
31		Naphthalene	0.998	0.945	5.3	107	0.00
32		Benzoic acid	0.209	0.238	-13.9	114	0.01
33		4-Chloroaniline	0.421	0.423	-0.5	107	0.00
34	C	Hexachlorobutadiene	0.197	0.194	1.5	105	0.00
35		Caprolactam	0.084	0.087	-3.6	106	0.01
36	C	4-Chloro-3-methylphenol	0.302	0.321	-6.3	118	0.01
37		2-Methylnaphthalene	0.627	0.632	-0.8	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.628	0.549	12.6	108	0.00
40	P	Hexachlorocyclopentadiene	0.386	0.363	6.0	105	0.00
41	S	2,4,6-Tribromophenol	0.208	0.218	-4.8	109	0.00
42	C	2,4,6-Trichlorophenol	0.437	0.395	9.6	99	0.00
43		2,4,5-Trichlorophenol	0.433	0.409	5.5	105	0.00
44	S	2-Fluorobiphenyl	1.341	1.195	10.9	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.534	1.409	8.1	108	0.00
46		2-Chloronaphthalene	1.253	1.172	6.5	108	0.00
47		2-Nitroaniline	0.487	0.505	-3.7	112	0.00
48		Acenaphthylene	1.992	1.902	4.5	104	0.00
49		Dimethylphthalate	1.409	1.211	14.1	98	0.00
50		2,6-Dinitrotoluene	0.299	0.260	13.0	97	0.00
51	C	Acenaphthene	1.215	1.158	4.7	104	0.00
52		3-Nitroaniline	0.345	0.351	-1.7	111	0.00
53	P	2,4-Dinitrophenol	0.159	0.148	6.9	109	0.01
54		Dibenzofuran	1.625	1.582	2.6	107	0.00
55	P	4-Nitrophenol	0.254	0.251	1.2	103	0.00
56		2,4-Dinitrotoluene	0.392	0.352	10.2	98	0.00
57		Fluorene	1.261	1.220	3.3	111	0.00
58		2,3,4,6-Tetrachlorophenol	0.327	0.339	-3.7	108	0.00
59		Diethylphthalate	1.359	1.242	8.6	106	0.00
60		4-Chlorophenyl-phenylether	0.596	0.527	11.6	109	0.00
61		4-Nitroaniline	0.312	0.277	11.2	100	0.00
62		Azobenzene	1.472	1.483	-0.7	119	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	0.120	0.123	-2.5	112	0.01
65	c	n-Nitrosodiphenylamine	0.656	0.646	1.5	108	0.00
66		4-Bromophenyl-phenylether	0.206	0.217	-5.3	106	0.00
67		Hexachlorobenzene	0.239	0.224	6.3	107	0.00
68		Atrazine	0.158	0.171	-8.2	105	0.01
69	C	Pentachlorophenol	0.144	0.144	0.0	101	0.00
70		Phenanthrene	1.008	0.919	8.8	102	0.00
71		Anthracene	0.990	0.996	-0.6	104	0.00
72		Carbazole	0.859	0.764	11.1	101	0.00
73		Di-n-butylphthalate	1.066	1.086	-1.9	104	0.00
74	C	Fluoranthene	0.974	0.975	-0.1	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76		Benzidine	0.728	0.741	-1.8	93	0.00
77		Pyrene	1.551	1.642	-5.9	98	0.00
78	S	Terphenyl-d14	1.057	1.121	-6.1	99	0.00
79		Butylbenzylphthalate	0.592	0.619	-4.6	98	0.00
80		Benzo(a)anthracene	1.203	1.128	6.2	93	0.00
81		3,3'-Dichlorobenzidine	0.411	0.391	4.9	95	0.00
82		Chrysene	1.168	1.189	-1.8	97	0.00
83		Bis(2-ethylhexyl)phthalate	0.772	0.801	-3.8	96	0.00
84	c	Di-n-octyl phthalate	1.295	1.352	-4.4	98	0.00
85		Indeno(1,2,3-cd)pyrene	0.958	1.140	-19.0	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	96	0.00
87		Benzo(b)fluoranthene	1.195	1.127	5.7	96	0.00

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88	Benzo(k)fluoranthene	0.983	0.953	3.1	93	0.00
89 C	Benzo(a)pyrene	0.981	0.946	3.6	95	0.00
90	Dibenzo(a,h)anthracene	0.878	1.001	-14.0	110	0.00
91	Benzo(g,h,i)perylene	0.864	1.026	-18.8	115	0.01

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