Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081296.D
Acq On : 30 Aug 2015 4:05
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 31 03:03:01 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 31 00:56:51 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	121	0.00
2		1,4-Dioxane	0.627	0.656	-4.6	123	0.01
3		Pyridine	1.769	1.660	6.2	101	0.02
4		n-Nitrosodimethylamine	0.985	1.029	-4.5	124	0.02
5	S	2-Fluorophenol	1.330	1.411	-6.1	133	0.02
6		Aniline	2.523	2.408	4.6	118	0.00
7	S	Phenol-d6	1.735	1.670	3.7	110	0.02
8		2-Chlorophenol	1.455	1.526	-4.9	117	0.01
9		Benzaldehyde	1.118	1.088	2.7	110	0.00
10	C	Phenol	2.049	2.088	-1.9	111	0.01
11		bis(2-Chloroethyl)ether	1.572	1.633	-3.9	123	0.00
12		1,3-Dichlorobenzene	1.564	1.563	0.1	120	0.00
13	C	1,4-Dichlorobenzene	1.549	1.589	-2.6	125	0.00
14		1,2-Dichlorobenzene	1.449	1.320	8.9	102	0.01
15		Benzyl Alcohol	1.404	1.303	7.2	120	0.00
16		2,2'-oxybis(1-Chloropropane	3.089	2.904	6.0	109	0.00
17		2-Methylphenol	1.249	1.144	8.4	106	0.01
18		Hexachloroethane	0.626	0.640	-2.2	119	0.00
19	P	n-Nitroso-di-n-propylamine	1.202	1.351	-12.4	127	0.01
20		3+4-Methylphenols	1.585	1.491	5.9	114	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
22		Acetophenone	0.525	0.519	1.1	114	0.00
23	S	Nitrobenzene-d5	0.438	0.446	-1.8	127	0.00
24		Nitrobenzene	0.458	0.482	-5.2	123	0.00
25		Isophorone	0.816	0.849	-4.0	127	0.01
26	C	2-Nitrophenol	0.190	0.188	1.1	126	0.00
27		2,4-Dimethylphenol	0.337	0.320	5.0	106	0.01
28		bis(2-Chloroethoxy)methane	0.482	0.526	-9.1	134	0.00
29	C	2,4-Dichlorophenol	0.284	0.275	3.2	113	0.01
30		1,2,4-Trichlorobenzene	0.315	0.326	-3.5	121	0.00
31		Naphthalene	1.115	0.982	11.9	107	0.00
32		Benzoic acid	0.189	0.192	-1.6	145	0.01
33		4-Chloroaniline	0.470	0.471	-0.2	132	0.01
34	C	Hexachlorobutadiene	0.178	0.187	-5.1	127	0.00
35		Caprolactam	0.099	0.098	1.0	116	0.01
36	C	4-Chloro-3-methylphenol	0.338	0.357	-5.6	128	0.01
37		2-Methylnaphthalene	0.704	0.733	-4.1	124	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
39		1,2,4,5-Tetrachlorobenzene	0.645	0.560	13.2	111	0.00
40	P	Hexachlorocyclopentadiene	0.334	0.234	29.9#	84	0.00
41	S	2,4,6-Tribromophenol	0.173	0.174	-0.6	134	0.00
42	C	2,4,6-Trichlorophenol	0.456	0.442	3.1	124	0.01
43		2,4,5-Trichlorophenol	0.456	0.455	0.2	114	0.01
44	S	2-Fluorobiphenyl	1.526	1.282	16.0	119	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.762	1.653	6.2	108	0.01
46		2-Chloronaphthalene	1.415	1.223	13.6	113	0.00
47		2-Nitroaniline	0.579	0.604	-4.3	131	0.01
48		Acenaphthylene	2.532	2.306	8.9	124	0.00
49		Dimethylphthalate	1.647	1.577	4.3	114	0.00
50		2,6-Dinitrotoluene	0.354	0.288	18.6	96	0.00
51	C	Acenaphthene	1.516	1.301	14.2	112	0.00
52		3-Nitroaniline	0.443	0.453	-2.3	134	0.01
53	P	2,4-Dinitrophenol	0.167	0.074	55.7#	54	0.00
54		Dibenzofuran	2.031	1.917	5.6	128	0.00
55	P	4-Nitrophenol	0.311	0.258	17.0	111	0.01
56		2,4-Dinitrotoluene	0.469	0.368	21.5	97	0.00
57		Fluorene	1.562	1.420	9.1	123	0.01
58		2,3,4,6-Tetrachlorophenol	0.353	0.316	10.5	105	0.00
59		Diethylphthalate	1.743	1.788	-2.6	127	0.00
60		4-Chlorophenyl-phenylether	0.661	0.625	5.4	119	0.01
61		4-Nitroaniline	0.452	0.371	17.9	107	0.00
62		Azobenzene	2.009	1.698	15.5	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
64		4,6-Dinitro-2-methylphenol	0.119	0.068	42.9#	59	0.01
65	c	n-Nitrosodiphenylamine	0.714	0.711	0.4	110	0.00
66		4-Bromophenyl-phenylether	0.196	0.220	-12.2	132	0.00
67		Hexachlorobenzene	0.199	0.220	-10.6	118	0.00
68		Atrazine	0.199	0.210	-5.5	121	0.00
69	C	Pentachlorophenol	0.119	0.061	48.7#	58	0.00
70		Phenanthrene	1.185	1.128	4.8	112	0.01
71		Anthracene	1.153	1.141	1.0	110	0.00
72		Carbazole	1.110	0.931	16.1	85	0.00
73		Di-n-butylphthalate	1.344	1.537	-14.4	126	0.01
74	C	Fluoranthene	1.279	1.004	21.5#	87	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
76		Benzidine	0.756	0.669	11.5	68	0.00
77		Pyrene	1.626	1.787	-9.9	111	0.00
78	S	Terphenyl-d14	0.933	1.099	-17.8	107	0.00
79		Butylbenzylphthalate	0.699	0.845	-20.9	104	0.00
80		Benzo(a)anthracene	1.341	1.272	5.1	85	0.00
81		3,3'-Dichlorobenzidine	0.434	0.389	10.4	85	0.00
82		Chrysene	1.209	0.983	18.7	72	0.00
83		Bis(2-ethylhexyl)phthalate	0.895	1.101	-23.0	112	0.00
84	c	Di-n-octyl phthalate	1.361	1.798	-32.1#	116	0.01
85		Indeno(1,2,3-cd)pyrene	0.849	1.210	-42.5#	129	0.01
86	I	Perylene-d12	1.000	1.000	0.0	112	0.01
87		Benzo(b)fluoranthene	1.415	1.322	6.6	95	-0.02

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88	Benzo(k)fluoranthene	1.318	1.030	21.9	95	0.00
89 C	Benzo(a)pyrene	1.233	1.194	3.2	107	0.00
90	Dibenzo(a,h)anthracene	0.960	1.120	-16.7	127	0.00
91	Benzo(g,h,i)perylene	0.974	1.121	-15.1	127	0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 3