Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
Data File : BF080019.D
Acq On : 17 Jun 2015 7:53
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 17 14:54:17 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 15 15:44:31 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	146	0.00
2		1,4-Dioxane	0.496	0.563	-13.5	162#	0.01
3		Pyridine	1.272	1.404	-10.4	172#	0.00
4		n-Nitrosodimethylamine	0.627	0.769	-22.6	195#	0.00
5	S	2-Fluorophenol	1.091	1.233	-13.0	161#	0.01
6		Aniline	2.151	2.361	-9.8	152#	0.00
7	S	Phenol-d6	1.486	1.613	-8.5	147	0.00
8		2-Chlorophenol	1.326	1.364	-2.9	143	0.00
9		Benzaldehyde	0.854	0.842	1.4	153#	0.01
10	C	Phenol	1.634	1.656	-1.3	137	0.01
11		bis(2-Chloroethyl)ether	1.315	1.250	4.9	128	0.00
12		1,3-Dichlorobenzene	1.546	1.619	-4.7	147	0.00
13	C	1,4-Dichlorobenzene	1.606	1.807	-12.5	160#	0.00
14		1,2-Dichlorobenzene	1.427	1.557	-9.1	155#	0.00
15		Benzyl Alcohol	1.191	1.324	-11.2	154#	0.00
16		2,2'-oxybis(1-Chloropropane	1.948	2.104	-8.0	150#	0.00
17		2-Methylphenol	1.158	1.130	2.4	135	0.00
18		Hexachloroethane	0.497	0.544	-9.5	153#	0.00
19	P	n-Nitroso-di-n-propylamine	1.062	1.088	-2.4	141	0.01
20		3+4-Methylphenols	1.480	1.576	-6.5	145	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
22		Acetophenone	0.472	0.505	-7.0	141	0.00
23	S	Nitrobenzene-d5	0.303	0.403	-33.0#	178#	0.00
24		Nitrobenzene	0.322	0.383	-18.9	158#	0.00
25		Isophorone	0.705	0.689	2.3	129	0.00
26	C	2-Nitrophenol	0.100	0.179	-79.0#	233#	0.00
27		2,4-Dimethylphenol	0.316	0.329	-4.1	145	0.00
28		bis(2-Chloroethoxy)methane	0.419	0.457	-9.1	145	0.00
29	C	2,4-Dichlorophenol	0.265	0.303	-14.3	156#	0.00
30		1,2,4-Trichlorobenzene	0.337	0.365	-8.3	149	0.00
31		Naphthalene	1.093	1.106	-1.2	136	0.00
32		Benzoic acid	0.107	0.206	-92.5#	245#	0.00
33		4-Chloroaniline	0.424	0.419	1.2	130	0.00
34	C	Hexachlorobutadiene	0.187	0.200	-7.0	150	0.00
35		Caprolactam	0.084	0.092	-9.5	131	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.305	-1.0	130	0.00
37		2-Methylnaphthalene	0.739	0.757	-2.4	139	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.00
39		1,2,4,5-Tetrachlorobenzene	0.670	0.679	-1.3	138	0.00
40	P	Hexachlorocyclopentadiene	0.211	0.122	42.2#	70	0.00
41	S	2,4,6-Tribromophenol	0.182	0.226	-24.2	147	0.00
42	C	2,4,6-Trichlorophenol	0.413	0.479	-16.0	149	0.00
43		2,4,5-Trichlorophenol	0.456	0.467	-2.4	131	0.00
44	S	2-Fluorobiphenyl	1.428	1.325	7.2	130	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.619	1.608	0.7	135	0.01
46		2-Chloronaphthalene	1.378	1.338	2.9	132	0.00
47		2-Nitroaniline	0.256	0.364	-42.2#	176#	0.00
48		Acenaphthylene	2.326	2.388	-2.7	137	0.00
49		Dimethylphthalate	1.602	1.453	9.3	122	0.00
50		2,6-Dinitrotoluene	0.250	0.316	-26.4#	146	0.00
51	C	Acenaphthene	1.322	1.367	-3.4	140	0.00
52		3-Nitroaniline	0.319	0.365	-14.4	140	0.00
53	P	2,4-Dinitrophenol	0.037	0.071	-91.9#	244#	0.01
54		Dibenzofuran	1.932	1.863	3.6	128	0.00
55	P	4-Nitrophenol	0.215	0.313	-45.6#	172#	0.00
56		2,4-Dinitrotoluene	0.314	0.456	-45.2#	163#	0.00
57		Fluorene	1.684	1.662	1.3	127	0.00
58		2,3,4,6-Tetrachlorophenol	0.345	0.388	-12.5	138	0.00
59		Diethylphthalate	1.602	1.476	7.9	119	0.00
60		4-Chlorophenyl-phenylether	0.752	0.741	1.5	132	0.00
61		4-Nitroaniline	0.326	0.401	-23.0	141	0.00
62		Azobenzene	1.542	1.591	-3.2	135	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	127	0.01
64		4,6-Dinitro-2-methylphenol	0.038	0.072	-89.5#	229#	0.00
65	c	n-Nitrosodiphenylamine	0.646	0.677	-4.8	129	0.00
66		4-Bromophenyl-phenylether	0.224	0.233	-4.0	126	0.00
67		Hexachlorobenzene	0.244	0.248	-1.6	123	0.00
68		Atrazine	0.199	0.224	-12.6	130	0.01
69	C	Pentachlorophenol	0.095	0.128	-34.7#	165#	0.01
70		Phenanthrene	1.192	1.220	-2.3	125	0.01
71		Anthracene	1.153	1.192	-3.4	127	0.01
72		Carbazole	1.119	1.157	-3.4	123	0.01
73		Di-n-butylphthalate	1.219	1.243	-2.0	122	0.01
74	C	Fluoranthene	1.318	1.313	0.4	119	0.03
75	I	Chrysene-d12	1.000	1.000	0.0	121	0.09
76		Benzidine	0.554	0.651	-17.5	145	0.05
77		Pyrene	1.256	1.208	3.8	120	0.03
78	S	Terphenyl-d14	0.867	0.863	0.5	123	0.05
79		Butylbenzylphthalate	0.424	0.519	-22.4	138	0.07
80		Benzo(a)anthracene	1.270	1.233	2.9	122	0.09
81		3,3'-Dichlorobenzidine	0.406	0.444	-9.4	128	0.09
82		Chrysene	1.229	1.150	6.4	121	0.09
83		Bis(2-ethylhexyl)phthalate	0.679	0.786	-15.8	131	0.09
84	c	Di-n-octyl phthalate	1.107	1.377	-24.4#	137	0.13
85		Indeno(1,2,3-cd)pyrene	1.330	1.393	-4.7	129	0.16
86	I	Perylene-d12	1.000	1.000	0.0	128	0.14
87		Benzo(b)fluoranthene	1.225	1.230	-0.4	133	0.14

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.250	1.176	5.9	115	0.15
89 C	Benzo(a)pyrene	1.165	1.137	2.4	125	0.15
90	Dibenzo(a,h)anthracene	1.146	1.145	0.1	133	0.16
91	Benzo(g,h,i)perylene	1.185	1.139	3.9	128	0.17

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