Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG120515\
Data File : BG019942.D
Acq On : 5 Dec 2015 15:04
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 07 01:08:04 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	173#	-0.02
2		1,4-Dioxane	0.431	0.477	-10.7	193#	0.00
3		Pyridine	1.309	1.448	-10.6	190#	-0.02
4		n-Nitrosodimethylamine	0.689	0.676	1.9	161#	-0.01
5	S	2-Fluorophenol	1.107	1.173	-6.0	184#	-0.01
6		Aniline	2.164	2.287	-5.7	180#	-0.01
7	S	Phenol-d6	1.671	1.808	-8.2	187#	-0.01
8		2-Chlorophenol	1.349	1.409	-4.4	179#	-0.02
9		Benzaldehyde	1.114	1.077	3.3	169#	-0.02
10	C	Phenol	1.758	1.903	-8.2	184#	-0.01
11		bis(2-Chloroethyl)ether	1.302	1.367	-5.0	183#	-0.02
12		1,3-Dichlorobenzene	1.529	1.580	-3.3	180#	-0.02
13	C	1,4-Dichlorobenzene	1.580	1.599	-1.2	178#	-0.02
14		1,2-Dichlorobenzene	1.507	1.533	-1.7	177#	-0.02
15		Benzyl Alcohol	1.466	1.440	1.8	172#	-0.02
16		2,2'-oxybis(1-Chloropropane	2.124	2.235	-5.2	184#	-0.01
17		2-Methylphenol	1.241	1.301	-4.8	178#	-0.02
18		Hexachloroethane	0.591	0.602	-1.9	173#	-0.02
19	P	n-Nitroso-di-n-propylamine	1.318	1.280	2.9	169#	-0.02
20		3+4-Methylphenols	1.748	1.817	-3.9	174#	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	174#	-0.02
22		Acetophenone	0.548	0.553	-0.9	173#	-0.02
23	S	Nitrobenzene-d5	0.392	0.413	-5.4	180#	-0.02
24		Nitrobenzene	0.425	0.457	-7.5	181#	-0.02
25		Isophorone	0.840	0.827	1.5	169#	-0.02
26	C	2-Nitrophenol	0.164	0.189	-15.2	194#	-0.01
27		2,4-Dimethylphenol	0.342	0.346	-1.2	175#	-0.02
28		bis(2-Chloroethoxy)methane	0.411	0.424	-3.2	179#	-0.02
29	C	2,4-Dichlorophenol	0.349	0.366	-4.9	177#	-0.01
30		1,2,4-Trichlorobenzene	0.395	0.388	1.8	167#	-0.02
31		Naphthalene	1.071	1.080	-0.8	173#	-0.02
32		Benzoic acid	0.211	0.289	-37.0#	237#	0.00
33		4-Chloroaniline	0.472	0.475	-0.6	174#	-0.02
34	C	Hexachlorobutadiene	0.291	0.277	4.8	164#	-0.02
35		Caprolactam	0.130	0.127	2.3	172#	-0.02
36	C	4-Chloro-3-methylphenol	0.435	0.438	-0.7	175#	-0.02
37		2-Methylnaphthalene	0.785	0.772	1.7	170#	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	163#	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.759	0.758	0.1	164#	-0.02
40	P	Hexachlorocyclopentadiene	0.433	0.410	5.3	151#	-0.02
41	S	2,4,6-Tribromophenol	0.306	0.300	2.0	161#	-0.02
42	C	2,4,6-Trichlorophenol	0.502	0.508	-1.2	169#	-0.02
43		2,4,5-Trichlorophenol	0.531	0.536	-0.9	169#	-0.02
44	S	2-Fluorobiphenyl	1.465	1.433	2.2	162#	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.641	1.642	-0.1	163#	-0.02
46		2-Chloronaphthalene	1.265	1.267	-0.2	164#	-0.02
47		2-Nitroaniline	0.400	0.448	-12.0	180#	-0.02
48		Acenaphthylene	2.155	2.129	1.2	163#	-0.02
49		Dimethylphthalate	1.799	1.727	4.0	160#	-0.02
50		2,6-Dinitrotoluene	0.338	0.367	-8.6	171#	-0.02
51	C	Acenaphthene	1.308	1.335	-2.1	167#	-0.02
52		3-Nitroaniline	0.351	0.390	-11.1	178#	-0.02
53	P	2,4-Dinitrophenol	0.143	0.178	-24.5	208#	-0.02
54		Dibenzofuran	1.945	1.911	1.7	163#	-0.02
55	P	4-Nitrophenol	0.306	0.326	-6.5	170#	-0.01
56		2,4-Dinitrotoluene	0.472	0.537	-13.8	177#	-0.02
57		Fluorene	1.741	1.669	4.1	160#	-0.02
58		2,3,4,6-Tetrachlorophenol	0.491	0.487	0.8	159#	-0.02
59		Diethylphthalate	1.947	1.815	6.8	158#	-0.02
60		4-Chlorophenyl-phenylether	1.000	0.945	5.5	156#	-0.02
61		4-Nitroaniline	0.395	0.419	-6.1	169#	-0.02
62		Azobenzene	1.620	1.563	3.5	160#	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	155#	-0.02
64		4,6-Dinitro-2-methylphenol	0.103	0.125	-21.4	186#	-0.02
65	c	n-Nitrosodiphenylamine	0.579	0.584	-0.9	157#	-0.02
66		4-Bromophenyl-phenylether	0.249	0.248	0.4	154#	-0.02
67		Hexachlorobenzene	0.259	0.254	1.9	153#	-0.02
68		Atrazine	0.251	0.246	2.0	150	-0.02
69	C	Pentachlorophenol	0.151	0.170	-12.6	173#	-0.02
70		Phenanthrene	1.132	1.127	0.4	155#	-0.02
71		Anthracene	1.153	1.149	0.3	154#	-0.02
72		Carbazole	1.015	1.010	0.5	154#	-0.02
73		Di-n-butylphthalate	1.245	1.242	0.2	154#	-0.02
74	C	Fluoranthene	1.447	1.437	0.7	153#	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	156#	-0.03
76		Benzidine	0.657	0.616	6.2	141	-0.02
77		Pyrene	1.177	1.196	-1.6	155#	-0.02
78	S	Terphenyl-d14	0.820	0.788	3.9	146	-0.02
79		Butylbenzylphthalate	0.461	0.482	-4.6	160#	-0.03
80		Benzo(a)anthracene	1.224	1.207	1.4	154#	-0.03
81		3,3'-Dichlorobenzidine	0.466	0.448	3.9	147	-0.04
82		Chrysene	1.162	1.156	0.5	154#	-0.03
83		Bis(2-ethylhexyl)phthalate	0.677	0.692	-2.2	159#	-0.04
84	c	Di-n-octyl phthalate	1.100	1.181	-7.4	163#	-0.06
85		Indeno(1,2,3-cd)pyrene	1.280	1.397	-9.1	169#	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	162#	-0.05
87		Benzo(b)fluoranthene	1.268	1.262	0.5	163#	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.255	1.200	4.4	158#	-0.05
89 C	Benzo(a)pyrene	1.203	1.188	1.2	162#	-0.05
90	Dibenzo(a,h)anthracene	1.189	1.183	0.5	165#	-0.08
91	Benzo(g,h,i)perylene	1.167	1.195	-2.4	168#	-0.09

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