Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG022716\
Data File : BG021134.D
Acq On : 28 Feb 2016 3:32
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 29 00:40:12 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 26 16:39:44 2016
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	113	0.00
2		1,4-Dioxane	0.529	0.580	-9.6	121	-0.01
3		Pyridine	1.547	1.743	-12.7	122	-0.01
4		n-Nitrosodimethylamine	0.784	0.854	-8.9	115	-0.01
5	S	2-Fluorophenol	1.142	1.261	-10.4	120	0.00
6		Aniline	2.103	2.727	-29.7#	137	-0.01
7	S	Phenol-d6	1.700	2.114	-24.4	133	-0.01
8		2-Chlorophenol	1.425	1.606	-12.7	122	-0.01
9		Benzaldehyde	0.921	1.179	-28.0#	141	0.00
10	C	Phenol	1.741	2.283	-31.1#	140	-0.01
11		bis(2-Chloroethyl)ether	1.335	1.680	-25.8#	138	-0.01
12		1,3-Dichlorobenzene	1.516	1.605	-5.9	114	0.00
13	C	1,4-Dichlorobenzene	1.620	1.641	-1.3	114	0.00
14		1,2-Dichlorobenzene	1.534	1.596	-4.0	113	-0.01
15		Benzyl Alcohol	1.485	1.894	-27.5#	135	-0.01
16		2,2'-oxybis(1-Chloropropane	1.530	2.739	-79.0#	198#	0.00
17		2-Methylphenol	1.229	1.592	-29.5#	137	-0.01
18		Hexachloroethane	0.613	0.676	-10.3	118	-0.01
19	P	n-Nitroso-di-n-propylamine	1.311	1.887	-43.9#	149	-0.02
20		3+4-Methylphenols	1.704	2.229	-30.8#	138	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	130	-0.01
22		Acetophenone	0.552	0.580	-5.1	132	-0.01
23	S	Nitrobenzene-d5	0.420	0.468	-11.4	141	-0.01
24		Nitrobenzene	0.434	0.510	-17.5	146	-0.01
25		Isophorone	0.787	0.956	-21.5	153#	-0.01
26	C	2-Nitrophenol	0.183	0.180	1.6	124	0.00
27		2,4-Dimethylphenol	0.323	0.345	-6.8	135	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.460	-20.1	153#	-0.01
29	C	2,4-Dichlorophenol	0.317	0.301	5.0	124	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.307	14.5	108	0.00
31		Naphthalene	1.027	1.037	-1.0	129	0.00
32		Benzoic acid	0.297	0.321	-8.1	132	-0.03
33		4-Chloroaniline	0.429	0.461	-7.5	136	0.00
34	C	Hexachlorobutadiene	0.240	0.202	15.8	102	-0.01
35		Caprolactam	0.107	0.133	-24.3	155#	-0.03
36	C	4-Chloro-3-methylphenol	0.388	0.443	-14.2	142	0.00
37		2-Methylnaphthalene	0.721	0.738	-2.4	130	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.00
39		1,2,4,5-Tetrachlorobenzene	0.739	0.622	15.8	109	-0.01
40	P	Hexachlorocyclopentadiene	0.486	0.387	20.4	102	-0.01
41	S	2,4,6-Tribromophenol	0.262	0.214	18.3	103	0.00
42	C	2,4,6-Trichlorophenol	0.463	0.440	5.0	121	0.00
43		2,4,5-Trichlorophenol	0.476	0.466	2.1	122	0.00
44	S	2-Fluorobiphenyl	1.563	1.427	8.7	121	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.653	1.8	131	0.00
46		2-Chloronaphthalene	1.278	1.243	2.7	129	-0.01
47		2-Nitroaniline	0.414	0.556	-34.3#	171#	0.00
48		Acenaphthylene	2.022	2.079	-2.8	137	-0.01
49		Dimethylphthalate	1.724	1.727	-0.2	133	0.00
50		2,6-Dinitrotoluene	0.355	0.368	-3.7	135	0.00
51	C	Acenaphthene	1.379	1.380	-0.1	130	0.00
52		3-Nitroaniline	0.345	0.398	-15.4	150	0.00
53	P	2,4-Dinitrophenol	0.247	0.234	5.3	119	0.00
54		Dibenzofuran	1.747	1.743	0.2	131	-0.01
55	P	4-Nitrophenol	0.306	0.350	-14.4	152#	0.00
56		2,4-Dinitrotoluene	0.509	0.531	-4.3	133	0.00
57		Fluorene	1.700	1.682	1.1	131	0.00
58		2,3,4,6-Tetrachlorophenol	0.414	0.389	6.0	121	0.00
59		Diethylphthalate	1.817	1.893	-4.2	139	-0.01
60		4-Chlorophenyl-phenylether	0.935	0.857	8.3	121	-0.01
61		4-Nitroaniline	0.393	0.435	-10.7	147	0.00
62		Azobenzene	1.574	2.010	-27.7#	169#	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	134	0.00
64		4,6-Dinitro-2-methylphenol	0.147	0.142	3.4	123	0.00
65	c	n-Nitrosodiphenylamine	0.590	0.607	-2.9	134	0.00
66		4-Bromophenyl-phenylether	0.235	0.211	10.2	115	0.00
67		Hexachlorobenzene	0.248	0.211	14.9	110	0.00
68		Atrazine	0.219	0.213	2.7	123	-0.01
69	C	Pentachlorophenol	0.168	0.145	13.7	110	0.00
70		Phenanthrene	1.076	1.088	-1.1	132	0.00
71		Anthracene	1.084	1.106	-2.0	132	-0.01
72		Carbazole	0.936	0.981	-4.8	136	0.00
73		Di-n-butylphthalate	1.233	1.375	-11.5	142	-0.01
74	C	Fluoranthene	1.350	1.284	4.9	125	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	117	-0.01
76		Benzidine	0.526	0.587	-11.6	123	0.00
77		Pyrene	1.107	1.201	-8.5	125	0.00
78	S	Terphenyl-d14	0.856	0.849	0.8	113	-0.01
79		Butylbenzylphthalate	0.462	0.577	-24.9	139	-0.01
80		Benzo(a)anthracene	1.151	1.168	-1.5	116	-0.01
81		3,3'-Dichlorobenzidine	0.450	0.438	2.7	107	-0.01
82		Chrysene	1.108	1.110	-0.2	116	-0.01
83		Bis(2-ethylhexyl)phthalate	0.714	0.836	-17.1	130	-0.02
84	c	Di-n-octyl phthalate	1.196	1.397	-16.8	130	-0.03
85		Indeno(1,2,3-cd)pyrene	1.336	1.230	7.9	105	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	112	-0.03
87		Benzo(b)fluoranthene	1.250	1.261	-0.9	110	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.232	1.251	-1.5	113	-0.03
89 C	Benzo(a)pyrene	1.200	1.203	-0.3	110	-0.03
90	Dibenzo(a,h)anthracene	1.210	1.165	3.7	106	-0.06
91	Benzo(g,h,i)perylene	1.165	1.115	4.3	106	-0.04

(#) = Out of Range SPCC's out = 0 CCC's out = 1