Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110716\
Data File : BG024750.D
Acq On : 8 Nov 2016 6:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 08 06:58:34 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 26 11:14:13 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	103	0.00
2		1,4-Dioxane	0.499	0.485	2.8	104	0.00
3		Pyridine	1.493	1.553	-4.0	111	-0.01
4		n-Nitrosodimethylamine	0.773	0.778	-0.6	105	0.00
5	S	2-Fluorophenol	1.220	1.211	0.7	108	0.00
6		Aniline	2.121	2.156	-1.7	107	0.00
7	S	Phenol-d6	1.674	1.735	-3.6	109	0.00
8		2-Chlorophenol	1.241	1.249	-0.6	110	0.00
9		Benzaldehyde	1.034	0.995	3.8	102	0.00
10	C	Phenol	1.725	1.794	-4.0	111	0.00
11		bis(2-Chloroethyl)ether	1.296	1.302	-0.5	110	0.00
12		1,3-Dichlorobenzene	1.536	1.568	-2.1	112	-0.01
13	C	1,4-Dichlorobenzene	1.517	1.557	-2.6	111	-0.01
14		1,2-Dichlorobenzene	1.459	1.492	-2.3	111	0.00
15		Benzyl Alcohol	1.557	1.548	0.6	106	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.349	2.3	104	-0.01
17		2-Methylphenol	1.143	1.144	-0.1	108	0.00
18		Hexachloroethane	0.583	0.597	-2.4	115	-0.01
19	P	n-Nitroso-di-n-propylamine	1.234	1.260	-2.1	106	-0.01
20		3+4-Methylphenols	1.535	1.680	-9.4	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	-0.01
22		Acetophenone	0.514	0.538	-4.7	109	0.00
23	S	Nitrobenzene-d5	0.439	0.461	-5.0	110	-0.01
24		Nitrobenzene	0.489	0.505	-3.3	107	0.00
25		Isophorone	0.817	0.856	-4.8	107	0.00
26	C	2-Nitrophenol	0.181	0.198	-9.4	113	0.00
27		2,4-Dimethylphenol	0.318	0.350	-10.1	111	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.435	-2.8	109	0.00
29	C	2,4-Dichlorophenol	0.333	0.368	-10.5	112	0.00
30		1,2,4-Trichlorobenzene	0.395	0.420	-6.3	112	0.00
31		Naphthalene	0.998	1.069	-7.1	111	0.00
32		Benzoic acid	0.264	0.265	-0.4	104	0.03
33		4-Chloroaniline	0.433	0.463	-6.9	106	0.00
34	C	Hexachlorobutadiene	0.309	0.335	-8.4	117	-0.01
35		Caprolactam	0.122	0.132	-8.2	110	0.00
36	C	4-Chloro-3-methylphenol	0.402	0.443	-10.2	110	0.00
37		2-Methylnaphthalene	0.728	0.788	-8.2	113	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.694	-2.8	114	-0.01
40	P	Hexachlorocyclopentadiene	0.380	0.385	-1.3	109	0.00
41	S	2,4,6-Tribromophenol	0.299	0.336	-12.4	114	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.430	-6.2	115	0.00
43		2,4,5-Trichlorophenol	0.438	0.462	-5.5	112	0.00
44	S	2-Fluorobiphenyl	1.203	1.244	-3.4	112	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.400	-0.9	109	-0.01
46		2-Chloronaphthalene	1.057	1.081	-2.3	112	0.00
47		2-Nitroaniline	0.433	0.451	-4.2	104	0.00
48		Acenaphthylene	1.621	1.674	-3.3	109	0.00
49		Dimethylphthalate	1.420	1.473	-3.7	108	0.00
50		2,6-Dinitrotoluene	0.303	0.329	-8.6	109	0.00
51	C	Acenaphthene	1.208	1.106	8.4	98	-0.01
52		3-Nitroaniline	0.314	0.317	-1.0	103	0.00
53	P	2,4-Dinitrophenol	0.220	0.272	-23.6	122	0.00
54		Dibenzofuran	1.670	1.730	-3.6	110	0.00
55	P	4-Nitrophenol	0.273	0.273	0.0	103	0.00
56		2,4-Dinitrotoluene	0.440	0.485	-10.2	108	0.00
57		Fluorene	1.385	1.449	-4.6	109	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.488	-7.5	109	0.00
59		Diethylphthalate	1.489	1.514	-1.7	105	0.00
60		4-Chlorophenyl-phenylether	0.777	0.801	-3.1	110	0.00
61		4-Nitroaniline	0.343	0.371	-8.2	111	0.00
62		Azobenzene	1.485	1.492	-0.5	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.157	-13.8	113	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.573	-4.8	108	0.00
66		4-Bromophenyl-phenylether	0.243	0.260	-7.0	111	0.00
67		Hexachlorobenzene	0.268	0.300	-11.9	114	0.00
68		Atrazine	0.235	0.233	0.9	100	0.00
69	C	Pentachlorophenol	0.177	0.189	-6.8	104	0.00
70		Phenanthrene	1.019	1.070	-5.0	109	0.00
71		Anthracene	1.028	1.085	-5.5	108	0.00
72		Carbazole	0.938	0.994	-6.0	110	0.00
73		Di-n-butylphthalate	1.081	1.139	-5.4	108	0.00
74	C	Fluoranthene	1.289	1.399	-8.5	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.471	0.425	9.8	95	0.00
77		Pyrene	0.978	1.018	-4.1	107	0.00
78	S	Terphenyl-d14	0.669	0.675	-0.9	107	-0.01
79		Butylbenzylphthalate	0.408	0.415	-1.7	107	-0.01
80		Benzo(a)anthracene	1.099	1.121	-2.0	110	0.00
81		3,3'-Dichlorobenzidine	0.443	0.459	-3.6	110	0.00
82		Chrysene	1.046	1.067	-2.0	108	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.598	-2.6	109	-0.02
84	c	Di-n-octyl phthalate	0.968	0.996	-2.9	108	-0.03
85		Indeno(1,2,3-cd)pyrene	1.444	1.450	-0.4	111	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	99	-0.02
87		Benzo(b)fluoranthene	1.081	1.121	-3.7	111	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.044	1.075	-3.0	109	-0.01
89 C	Benzo(a)pyrene	1.056	1.096	-3.8	110	-0.02
90	Dibenzo(a,h)anthracene	1.159	1.168	-0.8	111	-0.03
91	Benzo(g,h,i)perylene	1.158	1.168	-0.9	112	-0.04

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