Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\
Data File : BG022431.D
Acq On : 18 Jun 2016 19:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 20 13:14:04 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 17:40:23 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	126	-0.05
2		1,4-Dioxane	0.496	0.517	-4.2	143	-0.06
3		Pyridine	1.340	1.413	-5.4	132	-0.06
4		n-Nitrosodimethylamine	0.300	0.377	-25.7#	153#	-0.06
5	S	2-Fluorophenol	1.034	1.143	-10.5	135	-0.04
6		Aniline	2.066	2.252	-9.0	130	-0.04
7	S	Phenol-d6	1.626	1.774	-9.1	131	-0.02
8		2-Chlorophenol	1.430	1.481	-3.6	128	-0.04
9		Benzaldehyde	0.895	0.927	-3.6	123	-0.04
10	C	Phenol	1.677	1.825	-8.8	133	-0.02
11		bis(2-Chloroethyl)ether	1.337	1.442	-7.9	132	-0.05
12		1,3-Dichlorobenzene	1.533	1.518	1.0	126	-0.05
13	C	1,4-Dichlorobenzene	1.527	1.560	-2.2	129	-0.05
14		1,2-Dichlorobenzene	1.539	1.543	-0.3	128	-0.05
15		Benzyl Alcohol	1.051	1.144	-8.8	136	-0.04
16		2,2'-oxybis(1-Chloropropane	1.387	1.682	-21.3	154#	-0.05
17		2-Methylphenol	1.251	1.333	-6.6	135	-0.02
18		Hexachloroethane	0.558	0.595	-6.6	137	-0.05
19	P	n-Nitroso-di-n-propylamine	1.101	1.228	-11.5	133	-0.05
20		3+4-Methylphenols	1.795	1.815	-1.1	127	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	122	-0.05
22		Acetophenone	0.431	0.464	-7.7	128	-0.05
23	S	Nitrobenzene-d5	0.287	0.327	-13.9	135	-0.04
24		Nitrobenzene	0.288	0.318	-10.4	132	-0.05
25		Isophorone	0.671	0.691	-3.0	127	-0.04
26	C	2-Nitrophenol	0.156	0.173	-10.9	128	-0.04
27		2,4-Dimethylphenol	0.283	0.298	-5.3	126	-0.04
28		bis(2-Chloroethoxy)methane	0.385	0.417	-8.3	131	-0.05
29	C	2,4-Dichlorophenol	0.262	0.277	-5.7	122	-0.03
30		1,2,4-Trichlorobenzene	0.293	0.291	0.7	120	-0.05
31		Naphthalene	0.998	1.011	-1.3	128	-0.05
32		Benzoic acid	0.087	0.131	-50.6#	175#	0.02
33		4-Chloroaniline	0.415	0.423	-1.9	122	-0.04
34	C	Hexachlorobutadiene	0.157	0.152	3.2	123	-0.05
35		Caprolactam	0.144	0.121	16.0	103	-0.04
36	C	4-Chloro-3-methylphenol	0.311	0.328	-5.5	125	-0.02
37		2-Methylnaphthalene	0.737	0.721	2.2	119	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	113	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.515	0.524	-1.7	114	-0.04
40	P	Hexachlorocyclopentadiene	0.233	0.246	-5.6	116	-0.05
41	S	2,4,6-Tribromophenol	0.226	0.189	16.4	92	-0.04
42	C	2,4,6-Trichlorophenol	0.345	0.356	-3.2	111	-0.04
43		2,4,5-Trichlorophenol	0.382	0.358	6.3	105	-0.02
44	S	2-Fluorobiphenyl	1.312	1.372	-4.6	117	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.505	1.608	-6.8	117	-0.04
46		2-Chloronaphthalene	1.154	1.213	-5.1	115	-0.04
47		2-Nitroaniline	0.275	0.309	-12.4	122	-0.04
48		Acenaphthylene	2.024	2.043	-0.9	114	-0.04
49		Dimethylphthalate	1.658	1.545	6.8	107	-0.04
50		2,6-Dinitrotoluene	0.360	0.350	2.8	106	-0.04
51	C	Acenaphthene	1.213	1.210	0.2	113	-0.04
52		3-Nitroaniline	0.400	0.363	9.3	109	-0.03
53	P	2,4-Dinitrophenol	0.108	0.190	-75.9#	177#	-0.03
54		Dibenzofuran	1.893	1.875	1.0	111	-0.04
55	P	4-Nitrophenol	0.276	0.299	-8.3	113	0.00
56		2,4-Dinitrotoluene	0.484	0.490	-1.2	107	-0.04
57		Fluorene	1.631	1.577	3.3	109	-0.04
58		2,3,4,6-Tetrachlorophenol	0.342	0.300	12.3	101	-0.04
59		Diethylphthalate	1.771	1.650	6.8	108	-0.04
60		4-Chlorophenyl-phenylether	0.738	0.712	3.5	108	-0.04
61		4-Nitroaniline	0.424	0.344	18.9	91	-0.03
62		Azobenzene	1.341	1.438	-7.2	122	-0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.04
64		4,6-Dinitro-2-methylphenol	0.092	0.100	-8.7	103	-0.03
65	c	n-Nitrosodiphenylamine	0.576	0.615	-6.8	105	-0.04
66		4-Bromophenyl-phenylether	0.187	0.190	-1.6	101	-0.04
67		Hexachlorobenzene	0.218	0.207	5.0	96	-0.04
68		Atrazine	0.213	0.200	6.1	95	-0.04
69	C	Pentachlorophenol	0.084	0.091	-8.3	110	-0.03
70		Phenanthrene	1.086	1.097	-1.0	104	-0.04
71		Anthracene	1.077	1.097	-1.9	102	-0.04
72		Carbazole	1.020	1.015	0.5	101	-0.04
73		Di-n-butylphthalate	1.334	1.357	-1.7	105	-0.04
74	C	Fluoranthene	1.249	1.199	4.0	100	-0.04
75	I	Chrysene-d12	1.000	1.000	0.0	92	-0.05
76		Benzidine	0.523	0.500	4.4	79	-0.03
77		Pyrene	1.243	1.336	-7.5	99	-0.04
78	S	Terphenyl-d14	0.842	0.890	-5.7	96	-0.04
79		Butylbenzylphthalate	0.577	0.693	-20.1	110	-0.04
80		Benzo(a)anthracene	1.121	1.153	-2.9	96	-0.05
81		3,3'-Dichlorobenzidine	0.315	0.311	1.3	87	-0.05
82		Chrysene	1.091	1.107	-1.5	96	-0.05
83		Bis(2-ethylhexyl)phthalate	0.848	0.992	-17.0	108	-0.05
84	c	Di-n-octyl phthalate	1.408	1.649	-17.1	107	-0.06
85		Indeno(1,2,3-cd)pyrene	1.224	1.200	2.0	89	-0.12
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.08
87		Benzo(b)fluoranthene	1.166	1.210	-3.8	92	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.148	1.198	-4.4	93	-0.08
89 C	Benzo(a)pyrene	1.115	1.165	-4.5	92	-0.08
90	Dibenzo(a,h)anthracene	1.050	1.057	-0.7	88	-0.13
91	Benzo(g,h,i)perylene	1.093	1.088	0.5	89	-0.12

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