Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG102017\
Data File : BG029386.D
Acq On : 21 Oct 2017 10:28
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 21 22:35:35 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 17:32:15 2017
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	104	0.00
2		1,4-Dioxane	0.486	0.529	-8.8	104	0.00
3		Pyridine	1.415	1.509	-6.6	101	-0.01
4		n-Nitrosodimethylamine	0.661	0.685	-3.6	100	-0.01
5	S	2-Fluorophenol	1.197	1.239	-3.5	99	0.00
6		Aniline	2.081	2.113	-1.5	96	0.00
7	S	Phenol-d6	1.680	1.748	-4.0	98	-0.01
8		2-Chlorophenol	1.372	1.406	-2.5	99	0.00
9		Benzaldehyde	0.845	0.828	2.0	93	0.00
10	C	Phenol	1.812	1.900	-4.9	99	-0.01
11		bis(2-Chloroethyl)ether	1.320	1.399	-6.0	99	0.00
12		1,3-Dichlorobenzene	1.541	1.534	0.5	96	-0.01
13	C	1,4-Dichlorobenzene	1.573	1.582	-0.6	96	-0.01
14		1,2-Dichlorobenzene	1.517	1.513	0.3	96	0.00
15		Benzyl Alcohol	1.184	1.253	-5.8	101	0.00
16		2,2'-oxybis(1-Chloropropane	2.242	2.243	-0.0	96	0.00
17		2-Methylphenol	1.204	1.237	-2.7	97	-0.01
18		Hexachloroethane	0.536	0.538	-0.4	95	0.00
19	P	n-Nitroso-di-n-propylamine	1.143	1.192	-4.3	100	0.00
20		3+4-Methylphenols	1.696	1.776	-4.7	99	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.482	0.490	-1.7	99	0.00
23	S	Nitrobenzene-d5	0.315	0.319	-1.3	97	0.00
24		Nitrobenzene	0.354	0.361	-2.0	98	0.00
25		Isophorone	0.657	0.690	-5.0	101	0.00
26	C	2-Nitrophenol	0.169	0.185	-9.5	102	-0.01
27		2,4-Dimethylphenol	0.306	0.309	-1.0	98	-0.01
28		bis(2-Chloroethoxy)methane	0.403	0.419	-4.0	101	-0.01
29	C	2,4-Dichlorophenol	0.326	0.339	-4.0	101	0.00
30		1,2,4-Trichlorobenzene	0.353	0.349	1.1	97	0.00
31		Naphthalene	0.997	0.971	2.6	95	-0.01
32		Benzoic acid	0.256	0.302	-18.0	108	0.00
33		4-Chloroaniline	0.419	0.440	-5.0	102	-0.01
34	C	Hexachlorobutadiene	0.219	0.227	-3.7	101	-0.01
35		Caprolactam	0.108	0.135	-25.0#	122	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.389	-8.4	105	-0.01
37		2-Methylnaphthalene	0.726	0.724	0.3	97	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
39		1,2,4,5-Tetrachlorobenzene	0.730	0.694	4.9	103	-0.01
40	P	Hexachlorocyclopentadiene	0.247	0.297	-20.2	127	0.00
41	S	2,4,6-Tribromophenol	0.258	0.320	-24.0	131	0.00
42	C	2,4,6-Trichlorophenol	0.458	0.468	-2.2	109	0.00
43		2,4,5-Trichlorophenol	0.471	0.495	-5.1	112	-0.01
44	S	2-Fluorobiphenyl	1.364	1.281	6.1	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.599	7.0	100	0.00
46		2-Chloronaphthalene	1.254	1.180	5.9	102	-0.01
47		2-Nitroaniline	0.344	0.376	-9.3	112	0.00
48		Acenaphthylene	1.962	1.899	3.2	105	0.00
49		Dimethylphthalate	1.560	1.653	-6.0	113	0.00
50		2,6-Dinitrotoluene	0.327	0.373	-14.1	117	0.00
51	C	Acenaphthene	1.277	1.186	7.1	99	0.00
52		3-Nitroaniline	0.353	0.397	-12.5	117	0.00
53	P	2,4-Dinitrophenol	0.153	0.268	-75.2#	186#	0.00
54		Dibenzofuran	1.823	1.813	0.5	108	0.00
55	P	4-Nitrophenol	0.291	0.333	-14.4	119	0.00
56		2,4-Dinitrotoluene	0.443	0.541	-22.1	125	0.00
57		Fluorene	1.506	1.527	-1.4	110	0.00
58		2,3,4,6-Tetrachlorophenol	0.393	0.445	-13.2	119	0.00
59		Diethylphthalate	1.555	1.701	-9.4	117	0.00
60		4-Chlorophenyl-phenylether	0.803	0.852	-6.1	114	0.00
61		4-Nitroaniline	0.362	0.419	-15.7	123	0.00
62		Azobenzene	1.311	1.325	-1.1	109	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	131	-0.01
64		4,6-Dinitro-2-methylphenol	0.105	0.129	-22.9	147	-0.01
65	c	n-Nitrosodiphenylamine	0.599	0.568	5.2	117	0.00
66		4-Bromophenyl-phenylether	0.229	0.229	0.0	123	0.00
67		Hexachlorobenzene	0.260	0.263	-1.2	125	0.00
68		Atrazine	0.214	0.226	-5.6	126	0.00
69	C	Pentachlorophenol	0.149	0.166	-11.4	129	-0.01
70		Phenanthrene	1.033	1.004	2.8	120	0.00
71		Anthracene	1.037	1.022	1.4	120	-0.01
72		Carbazole	0.997	0.998	-0.1	122	0.00
73		Di-n-butylphthalate	1.146	1.173	-2.4	124	-0.01
74	C	Fluoranthene	1.208	1.249	-3.4	126	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	137	0.00
76		Benzidine	0.604	0.512	15.2	106	-0.01
77		Pyrene	1.213	1.177	3.0	125	0.00
78	S	Terphenyl-d14	0.879	0.870	1.0	128	-0.01
79		Butylbenzylphthalate	0.517	0.508	1.7	126	0.00
80		Benzo(a)anthracene	1.174	1.198	-2.0	132	-0.01
81		3,3'-Dichlorobenzidine	0.485	0.489	-0.8	130	-0.01
82		Chrysene	1.125	1.142	-1.5	132	0.00
83		Bis(2-ethylhexyl)phthalate	0.742	0.716	3.5	124	-0.01
84	c	Di-n-octyl phthalate	1.293	1.262	2.4	125	-0.02
85		Indeno(1,2,3-cd)pyrene	1.383	1.508	-9.0	141	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	138	-0.02
87		Benzo(b)fluoranthene	1.145	1.202	-5.0	136	-0.02

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88	Benzo(k)fluoranthene	1.141	1.179	-3.3	135	-0.02
89 C	Benzo(a)pyrene	1.101	1.143	-3.8	135	-0.02
90	Dibenzo(a,h)anthracene	1.114	1.182	-6.1	136	-0.02
91	Benzo(g,h,i)perylene	1.035	1.176	-13.6	144	-0.04

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