Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032017\
Data File : BG026337.D
Acq On : 20 Mar 2017 21:43
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 21 04:58:31 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG032017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 20 17:43:27 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	107	0.00
2		1,4-Dioxane	0.506	0.513	-1.4	110	0.00
3		Pyridine	1.385	1.391	-0.4	104	0.00
4		n-Nitrosodimethylamine	0.379	0.370	2.4	103	0.00
5	S	2-Fluorophenol	1.127	1.146	-1.7	108	0.00
6		Aniline	2.021	2.034	-0.6	105	0.00
7	S	Phenol-d6	1.513	1.535	-1.5	106	0.00
8		2-Chlorophenol	1.269	1.292	-1.8	107	0.00
9		Benzaldehyde	1.021	1.018	0.3	104	0.00
10	C	Phenol	1.614	1.615	-0.1	105	0.00
11		bis(2-Chloroethyl)ether	1.324	1.331	-0.5	108	0.00
12		1,3-Dichlorobenzene	1.510	1.549	-2.6	109	0.00
13	C	1,4-Dichlorobenzene	1.528	1.559	-2.0	108	0.00
14		1,2-Dichlorobenzene	1.462	1.485	-1.6	107	0.00
15		Benzyl Alcohol	0.992	1.004	-1.2	105	0.00
16		2,2'-oxybis(1-Chloropropane	1.470	1.407	4.3	101	0.00
17		2-Methylphenol	1.147	1.165	-1.6	108	0.00
18		Hexachloroethane	0.500	0.505	-1.0	106	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	0.951	2.3	103	0.00
20		3+4-Methylphenols	1.578	1.629	-3.2	106	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.460	0.467	-1.5	105	0.00
23	S	Nitrobenzene-d5	0.333	0.334	-0.3	105	0.00
24		Nitrobenzene	0.328	0.327	0.3	104	0.00
25		Isophorone	0.627	0.625	0.3	105	0.00
26	C	2-Nitrophenol	0.171	0.178	-4.1	106	0.00
27		2,4-Dimethylphenol	0.281	0.289	-2.8	107	0.00
28		bis(2-Chloroethoxy)methane	0.407	0.406	0.2	105	0.00
29	C	2,4-Dichlorophenol	0.284	0.300	-5.6	109	0.00
30		1,2,4-Trichlorobenzene	0.319	0.330	-3.4	109	0.00
31		Naphthalene	1.011	1.022	-1.1	107	0.00
32		Benzoic acid	0.216	0.215	0.5	103	0.00
33		4-Chloroaniline	0.411	0.413	-0.5	105	0.00
34	C	Hexachlorobutadiene	0.188	0.192	-2.1	109	0.00
35		Caprolactam	0.112	0.111	0.9	104	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.306	-1.0	106	0.00
37		2-Methylnaphthalene	0.741	0.752	-1.5	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.602	0.616	-2.3	108	0.00
40	P	Hexachlorocyclopentadiene	0.317	0.330	-4.1	109	0.00
41	S	2,4,6-Tribromophenol	0.229	0.235	-2.6	109	0.00
42	C	2,4,6-Trichlorophenol	0.394	0.405	-2.8	107	0.00
43		2,4,5-Trichlorophenol	0.436	0.446	-2.3	107	0.00
44	S	2-Fluorobiphenyl	1.426	1.427	-0.1	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.656	1.658	-0.1	106	0.00
46		2-Chloronaphthalene	1.210	1.218	-0.7	107	0.00
47		2-Nitroaniline	0.344	0.348	-1.2	103	0.00
48		Acenaphthylene	2.109	2.127	-0.9	107	0.00
49		Dimethylphthalate	1.510	1.513	-0.2	106	0.00
50		2,6-Dinitrotoluene	0.352	0.370	-5.1	105	0.00
51	C	Acenaphthene	1.299	1.297	0.2	106	0.00
52		3-Nitroaniline	0.388	0.402	-3.6	105	0.00
53	P	2,4-Dinitrophenol	0.204	0.200	2.0	102	0.00
54		Dibenzofuran	1.828	1.836	-0.4	107	0.00
55	P	4-Nitrophenol	0.318	0.322	-1.3	103	0.00
56		2,4-Dinitrotoluene	0.496	0.509	-2.6	104	0.00
57		Fluorene	1.627	1.629	-0.1	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.380	0.385	-1.3	107	0.00
59		Diethylphthalate	1.543	1.533	0.6	105	0.00
60		4-Chlorophenyl-phenylether	0.771	0.776	-0.6	108	0.00
61		4-Nitroaniline	0.410	0.417	-1.7	104	0.00
62		Azobenzene	1.255	1.240	1.2	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
64		4,6-Dinitro-2-methylphenol	0.126	0.133	-5.6	106	0.00
65	c	n-Nitrosodiphenylamine	0.587	0.595	-1.4	106	0.00
66		4-Bromophenyl-phenylether	0.206	0.213	-3.4	106	0.00
67		Hexachlorobenzene	0.220	0.224	-1.8	107	0.00
68		Atrazine	0.222	0.224	-0.9	103	0.00
69	C	Pentachlorophenol	0.143	0.146	-2.1	105	0.00
70		Phenanthrene	1.074	1.063	1.0	104	0.00
71		Anthracene	1.096	1.105	-0.8	106	0.00
72		Carbazole	1.128	1.117	1.0	104	0.00
73		Di-n-butylphthalate	1.159	1.148	0.9	103	0.00
74	C	Fluoranthene	1.263	1.243	1.6	104	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
76		Benzidine	0.690	0.719	-4.2	100	0.00
77		Pyrene	1.158	1.194	-3.1	103	0.00
78	S	Terphenyl-d14	0.824	0.838	-1.7	104	0.00
79		Butylbenzylphthalate	0.463	0.476	-2.8	102	0.00
80		Benzo(a)anthracene	1.125	1.124	0.1	102	0.00
81		3,3'-Dichlorobenzidine	0.461	0.460	0.2	101	0.00
82		Chrysene	1.075	1.095	-1.9	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.700	0.699	0.1	101	0.00
84	c	Di-n-octyl phthalate	1.194	1.212	-1.5	102	0.00
85		Indeno(1,2,3-cd)pyrene	1.328	1.374	-3.5	104	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	105	0.00
87		Benzo(b)fluoranthene	1.179	1.199	-1.7	106	0.00

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88	Benzo(k)fluoranthene	1.143	1.152	-0.8	102	0.00
89 C	Benzo(a)pyrene	1.131	1.135	-0.4	104	0.00
90	Dibenzo(a,h)anthracene	1.143	1.153	-0.9	104	0.00
91	Benzo(g,h,i)perylene	1.152	1.163	-1.0	104	0.00

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