Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032017\
Data File : BG026335.D
Acq On : 20 Mar 2017 20:25
Operator : SJ/MA
Sample : SSTDCCC040EC
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040EC

Quant Time: Mar 21 04:56:54 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	101	0.00
2		1,4-Dioxane	0.506	0.500	1.2	102	0.00
3		Pyridine	1.385	1.408	-1.7	100	0.00
4		n-Nitrosodimethylamine	0.379	0.372	1.8	98	0.00
5	S	2-Fluorophenol	1.127	1.135	-0.7	102	0.00
6		Aniline	2.021	1.999	1.1	98	0.00
7	S	Phenol-d6	1.513	1.543	-2.0	101	0.00
8		2-Chlorophenol	1.269	1.280	-0.9	101	0.00
9		Benzaldehyde	1.021	1.015	0.6	99	0.00
10	C	Phenol	1.614	1.619	-0.3	99	0.00
11		bis(2-Chloroethyl)ether	1.324	1.311	1.0	101	0.00
12		1,3-Dichlorobenzene	1.510	1.515	-0.3	101	0.00
13	C	1,4-Dichlorobenzene	1.528	1.560	-2.1	102	0.00
14		1,2-Dichlorobenzene	1.462	1.488	-1.8	102	0.00
15		Benzyl Alcohol	0.992	0.987	0.5	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.470	1.395	5.1	95	0.00
17		2-Methylphenol	1.147	1.143	0.3	101	0.00
18		Hexachloroethane	0.500	0.499	0.2	100	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	0.946	2.8	97	0.00
20		3+4-Methylphenols	1.578	1.604	-1.6	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.460	0.472	-2.6	100	0.00
23	S	Nitrobenzene-d5	0.333	0.335	-0.6	99	0.00
24		Nitrobenzene	0.328	0.329	-0.3	98	0.00
25		Isophorone	0.627	0.626	0.2	99	0.00
26	C	2-Nitrophenol	0.171	0.177	-3.5	99	0.00
27		2,4-Dimethylphenol	0.281	0.284	-1.1	99	0.00
28		bis(2-Chloroethoxy)methane	0.407	0.405	0.5	98	0.00
29	C	2,4-Dichlorophenol	0.284	0.298	-4.9	102	0.00
30		1,2,4-Trichlorobenzene	0.319	0.327	-2.5	101	0.00
31		Naphthalene	1.011	1.022	-1.1	101	0.00
32		Benzoic acid	0.216	0.208	3.7	94	0.00
33		4-Chloroaniline	0.411	0.415	-1.0	99	0.00
34	C	Hexachlorobutadiene	0.188	0.191	-1.6	101	0.00
35		Caprolactam	0.112	0.112	0.0	98	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.304	-0.3	98	0.00
37		2-Methylnaphthalene	0.741	0.755	-1.9	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	0.602	0.624	-3.7	102	0.00
40	P	Hexachlorocyclopentadiene	0.317	0.332	-4.7	101	0.00
41	S	2,4,6-Tribromophenol	0.229	0.238	-3.9	102	0.00
42	C	2,4,6-Trichlorophenol	0.394	0.404	-2.5	99	0.00
43		2,4,5-Trichlorophenol	0.436	0.453	-3.9	101	0.00
44	S	2-Fluorobiphenyl	1.426	1.455	-2.0	101	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.656	1.692	-2.2	100	0.00
46		2-Chloronaphthalene	1.210	1.232	-1.8	101	0.00
47		2-Nitroaniline	0.344	0.355	-3.2	97	0.00
48		Acenaphthylene	2.109	2.154	-2.1	100	0.00
49		Dimethylphthalate	1.510	1.541	-2.1	100	0.00
50		2,6-Dinitrotoluene	0.352	0.372	-5.7	98	0.00
51	C	Acenaphthene	1.299	1.318	-1.5	100	0.00
52		3-Nitroaniline	0.388	0.409	-5.4	99	0.00
53	P	2,4-Dinitrophenol	0.204	0.210	-2.9	99	0.00
54		Dibenzofuran	1.828	1.862	-1.9	101	0.00
55	P	4-Nitrophenol	0.318	0.335	-5.3	99	0.00
56		2,4-Dinitrotoluene	0.496	0.517	-4.2	98	0.00
57		Fluorene	1.627	1.653	-1.6	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.380	0.388	-2.1	100	0.00
59		Diethylphthalate	1.543	1.556	-0.8	99	0.00
60		4-Chlorophenyl-phenylether	0.771	0.773	-0.3	99	0.00
61		4-Nitroaniline	0.410	0.431	-5.1	99	0.00
62		Azobenzene	1.255	1.254	0.1	97	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.126	0.133	-5.6	99	0.00
65	c	n-Nitrosodiphenylamine	0.587	0.600	-2.2	100	0.00
66		4-Bromophenyl-phenylether	0.206	0.216	-4.9	101	0.00
67		Hexachlorobenzene	0.220	0.227	-3.2	102	0.00
68		Atrazine	0.222	0.228	-2.7	98	0.00
69	C	Pentachlorophenol	0.143	0.146	-2.1	98	0.00
70		Phenanthrene	1.074	1.082	-0.7	99	0.00
71		Anthracene	1.096	1.123	-2.5	101	0.00
72		Carbazole	1.128	1.149	-1.9	100	0.00
73		Di-n-butylphthalate	1.159	1.178	-1.6	99	0.00
74	C	Fluoranthene	1.263	1.288	-2.0	101	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.690	0.695	-0.7	95	0.00
77		Pyrene	1.158	1.184	-2.2	101	0.00
78	S	Terphenyl-d14	0.824	0.836	-1.5	102	0.00
79		Butylbenzylphthalate	0.463	0.475	-2.6	100	0.00
80		Benzo(a)anthracene	1.125	1.122	0.3	100	0.00
81		3,3'-Dichlorobenzidine	0.461	0.464	-0.7	100	0.00
82		Chrysene	1.075	1.081	-0.6	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.700	0.702	-0.3	100	0.00
84	c	Di-n-octyl phthalate	1.194	1.210	-1.3	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.328	1.369	-3.1	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.179	1.208	-2.5	104	0.00

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88	Benzo(k)fluoranthene	1.143	1.156	-1.1	100	0.00
89 C	Benzo(a)pyrene	1.131	1.138	-0.6	102	0.00
90	Dibenzo(a,h)anthracene	1.143	1.162	-1.7	102	0.00
91	Benzo(g,h,i)perylene	1.152	1.175	-2.0	103	0.00

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