Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032217\
Data File : BG026393.D
Acq On : 23 Mar 2017 2:31
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 23 03:09:27 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	172#	0.00
2		1,4-Dioxane	0.506	0.480	5.1	166#	0.00
3		Pyridine	1.385	1.276	7.9	154#	0.00
4		n-Nitrosodimethylamine	0.379	0.337	11.1	151#	0.00
5	S	2-Fluorophenol	1.127	1.095	2.8	166#	0.00
6		Aniline	2.021	1.972	2.4	164#	0.00
7	S	Phenol-d6	1.513	1.449	4.2	161#	0.00
8		2-Chlorophenol	1.269	1.271	-0.2	170#	0.00
9		Benzaldehyde	1.021	0.702	31.2#	116	0.00
10	C	Phenol	1.614	1.551	3.9	161#	0.00
11		bis(2-Chloroethyl)ether	1.324	1.272	3.9	166#	0.00
12		1,3-Dichlorobenzene	1.510	1.545	-2.3	174#	0.00
13	C	1,4-Dichlorobenzene	1.528	1.562	-2.2	173#	0.00
14		1,2-Dichlorobenzene	1.462	1.479	-1.2	172#	0.00
15		Benzyl Alcohol	0.992	0.931	6.1	157#	0.00
16		2,2'-oxybis(1-Chloropropane	1.470	1.239	15.7	143	0.00
17		2-Methylphenol	1.147	1.144	0.3	171#	0.00
18		Hexachloroethane	0.500	0.500	0.0	169#	0.00
19	P	n-Nitroso-di-n-propylamine	0.973	0.898	7.7	156#	0.00
20		3+4-Methylphenols	1.578	1.594	-1.0	167#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	174#	0.00
22		Acetophenone	0.460	0.449	2.4	163#	0.00
23	S	Nitrobenzene-d5	0.333	0.288	13.5	146	0.00
24		Nitrobenzene	0.328	0.307	6.4	158#	0.00
25		Isophorone	0.627	0.603	3.8	163#	0.00
26	C	2-Nitrophenol	0.171	0.182	-6.4	175#	0.00
27		2,4-Dimethylphenol	0.281	0.284	-1.1	170#	0.00
28		bis(2-Chloroethoxy)methane	0.407	0.394	3.2	164#	0.00
29	C	2,4-Dichlorophenol	0.284	0.301	-6.0	177#	0.00
30		1,2,4-Trichlorobenzene	0.319	0.337	-5.6	179#	0.00
31		Naphthalene	1.011	0.985	2.6	167#	0.00
32		Benzoic acid	0.216	0.231	-6.9	180#	0.02
33		4-Chloroaniline	0.411	0.423	-2.9	173#	0.00
34	C	Hexachlorobutadiene	0.188	0.198	-5.3	181#	0.00
35		Caprolactam	0.112	0.114	-1.8	172#	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.303	0.0	169#	0.00
37		2-Methylnaphthalene	0.741	0.741	0.0	170#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	203#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.602	0.541	10.1	182#	0.00
40	P	Hexachlorocyclopentadiene	0.317	0.298	6.0	188#	0.00
41	S	2,4,6-Tribromophenol	0.229	0.214	6.6	190#	0.00
42	C	2,4,6-Trichlorophenol	0.394	0.360	8.6	182#	0.00
43		2,4,5-Trichlorophenol	0.436	0.403	7.6	186#	0.00
44	S	2-Fluorobiphenyl	1.426	1.080	24.3	154#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.656	1.396	15.7	170#	0.00
46		2-Chloronaphthalene	1.210	1.043	13.8	175#	0.00
47		2-Nitroaniline	0.344	0.278	19.2	157#	0.00
48		Acenaphthylene	2.109	1.776	15.8	171#	0.00
49		Dimethylphthalate	1.510	1.305	13.6	174#	0.00
50		2,6-Dinitrotoluene	0.352	0.326	7.4	178#	0.00
51	C	Acenaphthene	1.299	1.108	14.7	173#	0.00
52		3-Nitroaniline	0.388	0.348	10.3	173#	0.00
53	P	2,4-Dinitrophenol	0.204	0.193	5.4	188#	0.00
54		Dibenzofuran	1.828	1.550	15.2	172#	0.00
55	P	4-Nitrophenol	0.318	0.290	8.8	176#	0.00
56		2,4-Dinitrotoluene	0.496	0.457	7.9	179#	0.00
57		Fluorene	1.627	1.382	15.1	172#	0.00
58		2,3,4,6-Tetrachlorophenol	0.380	0.344	9.5	183#	0.00
59		Diethylphthalate	1.543	1.311	15.0	171#	0.00
60		4-Chlorophenyl-phenylether	0.771	0.685	11.2	182#	0.00
61		4-Nitroaniline	0.410	0.374	8.8	177#	0.00
62		Azobenzene	1.255	0.988	21.3	158#	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	179#	0.00
64		4,6-Dinitro-2-methylphenol	0.126	0.142	-12.7	191#	0.00
65	c	n-Nitrosodiphenylamine	0.587	0.583	0.7	174#	0.00
66		4-Bromophenyl-phenylether	0.206	0.220	-6.8	184#	0.00
67		Hexachlorobenzene	0.220	0.235	-6.8	189#	0.00
68		Atrazine	0.222	0.216	2.7	168#	0.00
69	C	Pentachlorophenol	0.143	0.150	-4.9	181#	0.00
70		Phenanthrene	1.074	1.008	6.1	166#	0.00
71		Anthracene	1.096	1.043	4.8	168#	0.00
72		Carbazole	1.128	1.068	5.3	167#	0.00
73		Di-n-butylphthalate	1.159	1.085	6.4	164#	0.00
74	C	Fluoranthene	1.263	1.185	6.2	167#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	178#	0.00
76		Benzidine	0.690	0.399	42.2#	95	0.00
77		Pyrene	1.158	1.106	4.5	164#	0.00
78	S	Terphenyl-d14	0.824	0.699	15.2	149	0.00
79		Butylbenzylphthalate	0.463	0.471	-1.7	173#	0.00
80		Benzo(a)anthracene	1.125	1.091	3.0	170#	0.00
81		3,3'-Dichlorobenzidine	0.461	0.440	4.6	165#	0.00
82		Chrysene	1.075	1.034	3.8	169#	0.00
83		Bis(2-ethylhexyl)phthalate	0.700	0.674	3.7	168#	0.00
84	c	Di-n-octyl phthalate	1.194	1.148	3.9	166#	0.00
85		Indeno(1,2,3-cd)pyrene	1.328	1.419	-6.9	184#	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	187#	-0.02
87		Benzo(b)fluoranthene	1.179	1.154	2.1	183#	0.00

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88	Benzo(k)fluoranthene	1.143	1.075	5.9	170#	0.00
89 C	Benzo(a)pyrene	1.131	1.094	3.3	179#	0.00
90	Dibenzo(a,h)anthracene	1.143	1.152	-0.8	186#	0.00
91	Benzo(g,h,i)perylene	1.152	1.148	0.3	183#	0.00

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