Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060615\
Data File : BG017504.D
Acq On : 6 Jun 2015 18:54
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 05:13:20 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 01:31:15 2015
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	100	0.00
2		1,4-Dioxane	0.511	0.500	2.2	103	-0.02
3		Pyridine	1.365	1.405	-2.9	99	-0.02
4		n-Nitrosodimethylamine	0.895	1.104	-23.4	118	-0.02
5	S	2-Fluorophenol	1.135	1.152	-1.5	101	0.00
6		Aniline	2.076	2.225	-7.2	107	0.00
7	S	Phenol-d6	1.547	1.688	-9.1	106	0.00
8		2-Chlorophenol	1.342	1.413	-5.3	106	0.00
9		Benzaldehyde	1.050	1.058	-0.8	96	0.00
10	C	Phenol	1.589	1.775	-11.7	107	0.00
11		bis(2-Chloroethyl)ether	1.210	1.373	-13.5	111	0.00
12		1,3-Dichlorobenzene	1.501	1.603	-6.8	108	0.00
13	C	1,4-Dichlorobenzene	1.573	1.649	-4.8	105	0.00
14		1,2-Dichlorobenzene	1.472	1.554	-5.6	106	0.00
15		Benzyl Alcohol	1.440	1.618	-12.4	107	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	2.284	-12.0	111	0.00
17		2-Methylphenol	1.205	1.361	-12.9	111	0.00
18		Hexachloroethane	0.627	0.714	-13.9	111	0.00
19	P	n-Nitroso-di-n-propylamine	1.308	1.450	-10.9	110	-0.01
20		3+4-Methylphenols	1.668	1.866	-11.9	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.490	0.479	2.2	109	0.00
23	S	Nitrobenzene-d5	0.401	0.408	-1.7	112	0.00
24		Nitrobenzene	0.416	0.422	-1.4	113	0.00
25		Isophorone	0.838	0.792	5.5	107	0.00
26	C	2-Nitrophenol	0.193	0.199	-3.1	112	0.00
27		2,4-Dimethylphenol	0.314	0.301	4.1	107	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.375	2.1	107	0.00
29	C	2,4-Dichlorophenol	0.306	0.302	1.3	109	0.00
30		1,2,4-Trichlorobenzene	0.359	0.351	2.2	111	0.00
31		Naphthalene	1.045	1.036	0.9	111	0.00
32		Benzoic acid	0.198	0.183	7.6	100	0.03
33		4-Chloroaniline	0.464	0.438	5.6	105	0.00
34	C	Hexachlorobutadiene	0.234	0.250	-6.8	120	0.00
35		Caprolactam	0.168	0.135	19.6	85	0.00
36	C	4-Chloro-3-methylphenol	0.398	0.372	6.5	102	0.01
37		2-Methylnaphthalene	0.803	0.812	-1.1	112	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
39		1,2,4,5-Tetrachlorobenzene	0.609	0.609	0.0	116	0.00
40	P	Hexachlorocyclopentadiene	0.304	0.297	2.3	105	0.00
41	S	2,4,6-Tribromophenol	0.278	0.260	6.5	103	0.00
42	C	2,4,6-Trichlorophenol	0.436	0.444	-1.8	109	0.00
43		2,4,5-Trichlorophenol	0.477	0.449	5.9	107	0.01
44	S	2-Fluorobiphenyl	1.326	1.351	-1.9	115	0.00

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Quant Time: Jun 08 05:13:20 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 01:31:15 2015
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)
45		1,1'-Biphenyl	1.462	1.488	-1.8	113	0.00
46		2-Chloronaphthalene	1.222	1.246	-2.0	110	0.00
47		2-Nitroaniline	0.451	0.443	1.8	107	0.00
48		Acenaphthylene	2.153	2.122	1.4	109	0.00
49		Dimethylphthalate	1.741	1.647	5.4	104	0.00
50		2,6-Dinitrotoluene	0.391	0.372	4.9	106	0.00
51	C	Acenaphthene	1.269	1.284	-1.2	115	0.00
52		3-Nitroaniline	0.424	0.365	13.9	95	0.00
53	P	2,4-Dinitrophenol	0.205	0.177	13.7	93	0.00
54		Dibenzofuran	1.873	1.860	0.7	109	0.00
55	P	4-Nitrophenol	0.306	0.271	11.4	89	0.02
56		2,4-Dinitrotoluene	0.557	0.531	4.7	101	0.00
57		Fluorene	1.698	1.738	-2.4	113	0.00
58		2,3,4,6-Tetrachlorophenol	0.435	0.416	4.4	101	0.00
59		Diethylphthalate	1.890	1.747	7.6	103	0.00
60		4-Chlorophenyl-phenylether	0.860	0.867	-0.8	113	0.00
61		4-Nitroaniline	0.466	0.372	20.2	84	0.00
62		Azobenzene	1.796	1.832	-2.0	112	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.148	0.142	4.1	95	0.00
65	c	n-Nitrosodiphenylamine	0.597	0.639	-7.0	105	0.00
66		4-Bromophenyl-phenylether	0.221	0.237	-7.2	110	0.00
67		Hexachlorobenzene	0.238	0.252	-5.9	107	0.00
68		Atrazine	0.263	0.260	1.1	96	0.00
69	C	Pentachlorophenol	0.128	0.124	3.1	94	0.00
70		Phenanthrene	1.115	1.151	-3.2	105	0.00
71		Anthracene	1.114	1.141	-2.4	104	0.00
72		Carbazole	1.100	1.059	3.7	97	0.00
73		Di-n-butylphthalate	1.409	1.357	3.7	96	0.00
74	C	Fluoranthene	1.458	1.399	4.0	97	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	96	0.00
76		Benzidine	0.687	0.705	-2.6	92	0.00
77		Pyrene	1.232	1.270	-3.1	97	0.00
78	S	Terphenyl-d14	0.968	0.996	-2.9	98	0.00
79		Butylbenzylphthalate	0.533	0.551	-3.4	96	0.00
80		Benzo(a)anthracene	1.230	1.239	-0.7	95	0.00
81		3,3'-Dichlorobenzidine	0.460	0.477	-3.7	97	0.00
82		Chrysene	1.114	1.124	-0.9	95	0.00
83		Bis(2-ethylhexyl)phthalate	0.769	0.788	-2.5	96	0.00
84	c	Di-n-octyl phthalate	1.289	1.355	-5.1	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.399	1.456	-4.1	98	0.02
86	I	Perylene-d12	1.000	1.000	0.0	98	0.01
87		Benzo(b)fluoranthene	1.267	1.265	0.2	94	0.01

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QLast Update : Thu Jun 04 01:31:15 2015
Response via : Initial Calibration

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.213	1.266	-4.4	104	0.01
89 C	Benzo(a)pyrene	1.162	1.180	-1.5	98	0.01
90	Dibenzo(a,h)anthracene	1.208	1.233	-2.1	98	0.01
91	Benzo(g,h,i)perylene	1.150	1.156	-0.5	98	0.02

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