Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060815\
Data File : BG017553.D
Acq On : 9 Jun 2015 1:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 09 04:10:13 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 02:10:45 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	107	-0.01
2		1,4-Dioxane	0.511	0.490	4.1	108	0.00
3		Pyridine	1.365	1.188	13.0	90	0.00
4		n-Nitrosodimethylamine	0.895	1.076	-20.2	123	0.00
5	S	2-Fluorophenol	1.135	1.086	4.3	102	0.00
6		Aniline	2.076	2.032	2.1	104	0.00
7	S	Phenol-d6	1.547	1.454	6.0	98	0.00
8		2-Chlorophenol	1.342	1.307	2.6	105	0.00
9		Benzaldehyde	1.050	0.977	7.0	95	0.00
10	C	Phenol	1.589	1.555	2.1	100	0.00
11		bis(2-Chloroethyl)ether	1.210	1.201	0.7	104	0.00
12		1,3-Dichlorobenzene	1.501	1.465	2.4	105	0.00
13	C	1,4-Dichlorobenzene	1.573	1.528	2.9	104	0.00
14		1,2-Dichlorobenzene	1.472	1.473	-0.1	107	0.00
15		Benzyl Alcohol	1.440	1.398	2.9	99	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	1.953	4.3	101	0.00
17		2-Methylphenol	1.205	1.180	2.1	103	-0.01
18		Hexachloroethane	0.627	0.669	-6.7	112	0.00
19	P	n-Nitroso-di-n-propylamine	1.308	1.290	1.4	104	0.00
20		3+4-Methylphenols	1.668	1.663	0.3	103	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.490	0.500	-2.0	104	0.00
23	S	Nitrobenzene-d5	0.401	0.420	-4.7	105	0.00
24		Nitrobenzene	0.416	0.441	-6.0	108	0.00
25		Isophorone	0.838	0.787	6.1	97	0.00
26	C	2-Nitrophenol	0.193	0.201	-4.1	104	-0.01
27		2,4-Dimethylphenol	0.314	0.313	0.3	102	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.368	3.9	96	0.00
29	C	2,4-Dichlorophenol	0.306	0.323	-5.6	106	0.00
30		1,2,4-Trichlorobenzene	0.359	0.371	-3.3	107	-0.01
31		Naphthalene	1.045	1.056	-1.1	104	0.00
32		Benzoic acid	0.198	0.202	-2.0	101	-0.01
33		4-Chloroaniline	0.464	0.447	3.7	98	-0.01
34	C	Hexachlorobutadiene	0.234	0.252	-7.7	110	0.00
35		Caprolactam	0.168	0.143	14.9	82	-0.02
36	C	4-Chloro-3-methylphenol	0.398	0.387	2.8	97	-0.01
37		2-Methylnaphthalene	0.803	0.808	-0.6	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.609	0.604	0.8	108	-0.01
40	P	Hexachlorocyclopentadiene	0.304	0.263	13.5	87	0.00
41	S	2,4,6-Tribromophenol	0.278	0.282	-1.4	105	-0.01
42	C	2,4,6-Trichlorophenol	0.436	0.443	-1.6	102	-0.01
43		2,4,5-Trichlorophenol	0.477	0.475	0.4	106	-0.02
44	S	2-Fluorobiphenyl	1.326	1.344	-1.4	107	-0.01

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Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 09 04:10:13 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 02:10:45 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	1.470	-0.5	105	-0.01
46		2-Chloronaphthalene	1.222	1.224	-0.2	102	-0.01
47		2-Nitroaniline	0.451	0.457	-1.3	103	-0.01
48		Acenaphthylene	2.153	2.153	0.0	104	-0.01
49		Dimethylphthalate	1.741	1.663	4.5	99	-0.01
50		2,6-Dinitrotoluene	0.391	0.373	4.6	100	-0.01
51	C	Acenaphthene	1.269	1.263	0.5	106	-0.01
52		3-Nitroaniline	0.424	0.391	7.8	96	-0.02
53	P	2,4-Dinitrophenol	0.205	0.202	1.5	99	-0.01
54		Dibenzofuran	1.873	1.857	0.9	102	-0.01
55	P	4-Nitrophenol	0.306	0.282	7.8	87	-0.02
56		2,4-Dinitrotoluene	0.557	0.552	0.9	99	-0.01
57		Fluorene	1.698	1.729	-1.8	106	-0.01
58		2,3,4,6-Tetrachlorophenol	0.435	0.447	-2.8	102	-0.01
59		Diethylphthalate	1.890	1.789	5.3	99	-0.01
60		4-Chlorophenyl-phenylether	0.860	0.878	-2.1	107	-0.01
61		4-Nitroaniline	0.466	0.412	11.6	87	-0.01
62		Azobenzene	1.796	1.863	-3.7	107	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.02
64		4,6-Dinitro-2-methylphenol	0.148	0.149	-0.7	100	0.00
65	c	n-Nitrosodiphenylamine	0.597	0.604	-1.2	101	-0.01
66		4-Bromophenyl-phenylether	0.221	0.231	-4.5	108	-0.02
67		Hexachlorobenzene	0.238	0.250	-5.0	107	-0.01
68		Atrazine	0.263	0.262	0.4	98	-0.01
69	C	Pentachlorophenol	0.128	0.120	6.3	91	-0.02
70		Phenanthrene	1.115	1.127	-1.1	103	-0.02
71		Anthracene	1.114	1.117	-0.3	103	-0.01
72		Carbazole	1.100	1.077	2.1	100	-0.02
73		Di-n-butylphthalate	1.409	1.404	0.4	100	-0.02
74	C	Fluoranthene	1.458	1.456	0.1	102	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	103	-0.02
76		Benzidine	0.687	0.615	10.5	85	-0.02
77		Pyrene	1.232	1.255	-1.9	102	-0.02
78	S	Terphenyl-d14	0.968	1.015	-4.9	106	-0.02
79		Butylbenzylphthalate	0.533	0.536	-0.6	100	-0.02
80		Benzo(a)anthracene	1.230	1.263	-2.7	103	-0.02
81		3,3'-Dichlorobenzidine	0.460	0.475	-3.3	102	-0.02
82		Chrysene	1.114	1.132	-1.6	102	-0.02
83		Bis(2-ethylhexyl)phthalate	0.769	0.794	-3.3	103	-0.02
84	c	Di-n-octyl phthalate	1.289	1.270	1.5	100	-0.02
85		Indeno(1,2,3-cd)pyrene	1.399	1.448	-3.5	104	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	101	-0.03
87		Benzo(b)fluoranthene	1.267	1.331	-5.1	103	-0.03

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 02:10:45 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)
88	Benzo(k)fluoranthene	1.213	1.240	-2.2	105	-0.03
89 C	Benzo(a)pyrene	1.162	1.205	-3.7	103	-0.03
90	Dibenzo(a,h)anthracene	1.208	1.257	-4.1	104	-0.05
91	Benzo(g,h,i)perylene	1.150	1.168	-1.6	103	-0.06

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