Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG061115\
Data File : BG017611.D
Acq On : 11 Jun 2015 16:04
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 12 01:16:43 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 01:16:03 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	167#	-0.08
2		1,4-Dioxane	0.511	0.568	-11.2	194#	-0.09
3		Pyridine	1.365	1.452	-6.4	171#	-0.09
4		n-Nitrosodimethylamine	0.895	1.398	-56.2#	249#	-0.09
5	S	2-Fluorophenol	1.135	1.220	-7.5	178#	-0.09
6		Aniline	2.076	2.193	-5.6	175#	-0.08
7	S	Phenol-d6	1.547	1.529	1.2	161#	-0.08
8		2-Chlorophenol	1.342	1.330	0.9	166#	-0.08
9		Benzaldehyde	1.050	0.935	11.0	142	-0.08
10	C	Phenol	1.589	1.683	-5.9	169#	-0.09
11		bis(2-Chloroethyl)ether	1.210	1.305	-7.9	176#	-0.08
12		1,3-Dichlorobenzene	1.501	1.578	-5.1	177#	-0.08
13	C	1,4-Dichlorobenzene	1.573	1.681	-6.9	179#	-0.08
14		1,2-Dichlorobenzene	1.472	1.545	-5.0	175#	-0.08
15		Benzyl Alcohol	1.440	1.739	-20.8	192#	-0.08
16		2,2'-oxybis(1-Chloropropane	2.040	2.077	-1.8	167#	-0.08
17		2-Methylphenol	1.205	1.252	-3.9	170#	-0.08
18		Hexachloroethane	0.627	0.687	-9.6	179#	-0.08
19	P	n-Nitroso-di-n-propylamine	1.308	1.374	-5.0	173#	-0.08
20		3+4-Methylphenols	1.668	1.770	-6.1	171#	-0.09
21	I	Naphthalene-d8	1.000	1.000	0.0	163#	-0.08
22		Acetophenone	0.490	0.527	-7.6	170#	-0.08
23	S	Nitrobenzene-d5	0.401	0.467	-16.5	183#	-0.08
24		Nitrobenzene	0.416	0.488	-17.3	186#	-0.08
25		Isophorone	0.838	0.841	-0.4	162#	-0.08
26	C	2-Nitrophenol	0.193	0.213	-10.4	171#	-0.08
27		2,4-Dimethylphenol	0.314	0.333	-6.1	169#	-0.08
28		bis(2-Chloroethoxy)methane	0.383	0.392	-2.3	159#	-0.08
29	C	2,4-Dichlorophenol	0.306	0.350	-14.4	179#	-0.08
30		1,2,4-Trichlorobenzene	0.359	0.428	-19.2	193#	-0.08
31		Naphthalene	1.045	1.075	-2.9	165#	-0.08
32		Benzoic acid	0.198	0.190	4.0	148	-0.08
33		4-Chloroaniline	0.464	0.465	-0.2	159#	-0.08
34	C	Hexachlorobutadiene	0.234	0.333	-42.3#	228#	-0.08
35		Caprolactam	0.168	0.149	11.3	133	-0.09
36	C	4-Chloro-3-methylphenol	0.398	0.417	-4.8	162#	-0.09
37		2-Methylnaphthalene	0.803	0.828	-3.1	163#	-0.08
38	I	Acenaphthene-d10	1.000	1.000	0.0	173#	-0.08
39		1,2,4,5-Tetrachlorobenzene	0.609	0.763	-25.3#	220#	-0.08
40	P	Hexachlorocyclopentadiene	0.304	0.222	27.0#	119	-0.08
41	S	2,4,6-Tribromophenol	0.278	0.288	-3.6	174#	-0.07
42	C	2,4,6-Trichlorophenol	0.436	0.479	-9.9	178#	-0.08
43		2,4,5-Trichlorophenol	0.477	0.511	-7.1	184#	-0.09
44	S	2-Fluorobiphenyl	1.326	1.465	-10.5	189#	-0.08

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Quant Time: Jun 12 01:16:43 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 01:16:03 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	1.506	-3.0	174#	-0.08
46		2-Chloronaphthalene	1.222	1.279	-4.7	172#	-0.08
47		2-Nitroaniline	0.451	0.488	-8.2	179#	-0.08
48		Acenaphthylene	2.153	2.233	-3.7	174#	-0.08
49		Dimethylphthalate	1.741	1.752	-0.6	169#	-0.07
50		2,6-Dinitrotoluene	0.391	0.410	-4.9	178#	-0.07
51	C	Acenaphthene	1.269	1.268	0.1	172#	-0.08
52		3-Nitroaniline	0.424	0.394	7.1	156#	-0.08
53	P	2,4-Dinitrophenol	0.205	0.202	1.5	160#	-0.06
54		Dibenzofuran	1.873	2.050	-9.5	183#	-0.07
55	P	4-Nitrophenol	0.306	0.274	10.5	136	-0.08
56		2,4-Dinitrotoluene	0.557	0.599	-7.5	174#	-0.07
57		Fluorene	1.698	1.892	-11.4	187#	-0.07
58		2,3,4,6-Tetrachlorophenol	0.435	0.533	-22.5	197#	-0.08
59		Diethylphthalate	1.890	1.901	-0.6	170#	-0.07
60		4-Chlorophenyl-phenylether	0.860	1.037	-20.6	205#	-0.07
61		4-Nitroaniline	0.466	0.410	12.0	140	-0.07
62		Azobenzene	1.796	2.058	-14.6	191#	-0.07
63	I	Phenanthrene-d10	1.000	1.000	0.0	186#	-0.08
64		4,6-Dinitro-2-methylphenol	0.148	0.145	2.0	180#	-0.06
65	c	n-Nitrosodiphenylamine	0.597	0.571	4.4	174#	-0.07
66		4-Bromophenyl-phenylether	0.221	0.245	-10.9	210#	-0.07
67		Hexachlorobenzene	0.238	0.249	-4.6	196#	-0.07
68		Atrazine	0.263	0.265	-0.8	182#	-0.07
69	C	Pentachlorophenol	0.128	0.112	12.5	156#	-0.08
70		Phenanthrene	1.115	1.066	4.4	179#	-0.08
71		Anthracene	1.114	1.067	4.2	180#	-0.08
72		Carbazole	1.100	1.004	8.7	171#	-0.08
73		Di-n-butylphthalate	1.409	1.295	8.1	170#	-0.08
74	C	Fluoranthene	1.458	1.481	-1.6	191#	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	192#	-0.07
76		Benzidine	0.687	0.528	23.1	137	-0.07
77		Pyrene	1.232	1.258	-2.1	192#	-0.08
78	S	Terphenyl-d14	0.968	1.010	-4.3	197#	-0.07
79		Butylbenzylphthalate	0.533	0.498	6.6	173#	-0.07
80		Benzo(a)anthracene	1.230	1.291	-5.0	198#	-0.07
81		3,3'-Dichlorobenzidine	0.460	0.444	3.5	179#	-0.08
82		Chrysene	1.114	1.181	-6.0	200#	-0.07
83		Bis(2-ethylhexyl)phthalate	0.769	0.710	7.7	173#	-0.06
84	c	Di-n-octyl phthalate	1.289	1.120	13.1	165#	-0.08
85		Indeno(1,2,3-cd)pyrene	1.399	1.335	4.6	179#	-0.16
86	I	Perylene-d12	1.000	1.000	0.0	172#	-0.11
87		Benzo(b)fluoranthene	1.267	1.361	-7.4	179#	-0.09

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QLast Update : Fri Jun 12 01:16:03 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	1.325	-9.2	192#	-0.10
89 C	Benzo(a)pyrene	1.162	1.251	-7.7	183#	-0.11
90	Dibenzo(a,h)anthracene	1.208	1.271	-5.2	178#	-0.16
91	Benzo(g,h,i)perylene	1.150	1.211	-5.3	181#	-0.18

(#) = Out of Range SPCC's out = 0 CCC's out = 1