Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG061115\
Data File : BG017628.D
Acq On : 12 Jun 2015 2:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 12 03:31:54 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	113	-0.08
2		1,4-Dioxane	0.511	0.493	3.5	114	-0.09
3		Pyridine	1.365	1.271	6.9	101	-0.09
4		n-Nitrosodimethylamine	0.895	1.250	-39.7#	151#	-0.09
5	S	2-Fluorophenol	1.135	1.079	4.9	106	-0.09
6		Aniline	2.076	2.060	0.8	111	-0.08
7	S	Phenol-d6	1.547	1.475	4.7	105	-0.08
8		2-Chlorophenol	1.342	1.319	1.7	111	-0.08
9		Benzaldehyde	1.050	0.919	12.5	94	-0.08
10	C	Phenol	1.589	1.599	-0.6	109	-0.08
11		bis(2-Chloroethyl)ether	1.210	1.194	1.3	109	-0.08
12		1,3-Dichlorobenzene	1.501	1.486	1.0	113	-0.08
13	C	1,4-Dichlorobenzene	1.573	1.560	0.8	112	-0.08
14		1,2-Dichlorobenzene	1.472	1.465	0.5	112	-0.08
15		Benzyl Alcohol	1.440	1.637	-13.7	122	-0.08
16		2,2'-oxybis(1-Chloropropane	2.040	1.944	4.7	106	-0.08
17		2-Methylphenol	1.205	1.257	-4.3	115	-0.08
18		Hexachloroethane	0.627	0.687	-9.6	121	-0.08
19	P	n-Nitroso-di-n-propylamine	1.308	1.436	-9.8	122	-0.08
20		3+4-Methylphenols	1.668	1.710	-2.5	112	-0.08
21	I	Naphthalene-d8	1.000	1.000	0.0	113	-0.08
22		Acetophenone	0.490	0.525	-7.1	118	-0.08
23	S	Nitrobenzene-d5	0.401	0.460	-14.7	125	-0.08
24		Nitrobenzene	0.416	0.472	-13.5	126	-0.08
25		Isophorone	0.838	0.897	-7.0	121	-0.08
26	C	2-Nitrophenol	0.193	0.210	-8.8	118	-0.08
27		2,4-Dimethylphenol	0.314	0.326	-3.8	115	-0.08
28		bis(2-Chloroethoxy)methane	0.383	0.396	-3.4	112	-0.08
29	C	2,4-Dichlorophenol	0.306	0.350	-14.4	125	-0.09
30		1,2,4-Trichlorobenzene	0.359	0.419	-16.7	131	-0.08
31		Naphthalene	1.045	1.045	0.0	111	-0.08
32		Benzoic acid	0.198	0.214	-8.1	116	-0.08
33		4-Chloroaniline	0.464	0.471	-1.5	112	-0.08
34	C	Hexachlorobutadiene	0.234	0.324	-38.5#	154#	-0.08
35		Caprolactam	0.168	0.154	8.3	96	-0.09
36	C	4-Chloro-3-methylphenol	0.398	0.422	-6.0	114	-0.09
37		2-Methylnaphthalene	0.803	0.857	-6.7	117	-0.08
38	I	Acenaphthene-d10	1.000	1.000	0.0	128	-0.08
39		1,2,4,5-Tetrachlorobenzene	0.609	0.740	-21.5	158#	-0.08
40	P	Hexachlorocyclopentadiene	0.304	0.227	25.3#	90	-0.08
41	S	2,4,6-Tribromophenol	0.278	0.287	-3.2	128	-0.08
42	C	2,4,6-Trichlorophenol	0.436	0.476	-9.2	131	-0.08
43		2,4,5-Trichlorophenol	0.477	0.524	-9.9	140	-0.09
44	S	2-Fluorobiphenyl	1.326	1.445	-9.0	138	-0.08

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LabSampleId : SSTDCCC040

Quant Time: Jun 12 03:31:54 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	1.472	-0.7	126	-0.08
46		2-Chloronaphthalene	1.222	1.271	-4.0	127	-0.08
47		2-Nitroaniline	0.451	0.470	-4.2	127	-0.08
48		Acenaphthylene	2.153	2.175	-1.0	125	-0.08
49		Dimethylphthalate	1.741	1.707	2.0	121	-0.08
50		2,6-Dinitrotoluene	0.391	0.392	-0.3	126	-0.08
51	C	Acenaphthene	1.269	1.259	0.8	126	-0.08
52		3-Nitroaniline	0.424	0.379	10.6	111	-0.08
53	P	2,4-Dinitrophenol	0.205	0.200	2.4	117	-0.07
54		Dibenzofuran	1.873	2.021	-7.9	133	-0.08
55	P	4-Nitrophenol	0.306	0.252	17.6	93	-0.09
56		2,4-Dinitrotoluene	0.557	0.580	-4.1	124	-0.08
57		Fluorene	1.698	1.830	-7.8	134	-0.08
58		2,3,4,6-Tetrachlorophenol	0.435	0.545	-25.3#	149	-0.08
59		Diethylphthalate	1.890	1.841	2.6	121	-0.08
60		4-Chlorophenyl-phenylether	0.860	1.028	-19.5	150	-0.08
61		4-Nitroaniline	0.466	0.394	15.5	99	-0.08
62		Azobenzene	1.796	2.016	-12.2	139	-0.08
63	I	Phenanthrene-d10	1.000	1.000	0.0	132	-0.08
64		4,6-Dinitro-2-methylphenol	0.148	0.150	-1.4	132	-0.07
65	c	n-Nitrosodiphenylamine	0.597	0.601	-0.7	130	-0.08
66		4-Bromophenyl-phenylether	0.221	0.243	-10.0	147	-0.08
67		Hexachlorobenzene	0.238	0.254	-6.7	141	-0.08
68		Atrazine	0.263	0.270	-2.7	131	-0.08
69	C	Pentachlorophenol	0.128	0.112	12.5	111	-0.08
70		Phenanthrene	1.115	1.094	1.9	130	-0.08
71		Anthracene	1.114	1.085	2.6	130	-0.08
72		Carbazole	1.100	1.010	8.2	122	-0.08
73		Di-n-butylphthalate	1.409	1.270	9.9	118	-0.08
74	C	Fluoranthene	1.458	1.500	-2.9	137	-0.08
75	I	Chrysene-d12	1.000	1.000	0.0	133	-0.08
76		Benzidine	0.687	0.550	19.9	99	-0.08
77		Pyrene	1.232	1.261	-2.4	134	-0.08
78	S	Terphenyl-d14	0.968	1.054	-8.9	143	-0.08
79		Butylbenzylphthalate	0.533	0.500	6.2	121	-0.08
80		Benzo(a)anthracene	1.230	1.336	-8.6	142	-0.08
81		3,3'-Dichlorobenzidine	0.460	0.465	-1.1	130	-0.08
82		Chrysene	1.114	1.225	-10.0	144	-0.08
83		Bis(2-ethylhexyl)phthalate	0.769	0.726	5.6	123	-0.08
84	c	Di-n-octyl phthalate	1.289	1.152	10.6	118	-0.09
85		Indeno(1,2,3-cd)pyrene	1.399	1.373	1.9	128	-0.18
86	I	Perylene-d12	1.000	1.000	0.0	124	-0.12
87		Benzo(b)fluoranthene	1.267	1.355	-6.9	129	-0.10

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Instrument : BNA_G
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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)
88	Benzo(k)fluoranthene	1.213	1.310	-8.0	137	-0.11
89 C	Benzo(a)pyrene	1.162	1.235	-6.3	130	-0.12
90	Dibenzo(a,h)anthracene	1.208	1.244	-3.0	126	-0.18
91	Benzo(g,h,i)perylene	1.150	1.192	-3.7	129	-0.20

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