Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120315\
Data File : BG019891.D
Acq On : 4 Dec 2015 4:25
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 04 05:02:12 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	143	-0.02
2		1,4-Dioxane	40.000	42.864	-7.2	155	0.00
3		Pyridine	40.000	43.918	-9.8	157	-0.02
4		n-Nitrosodimethylamine	40.000	39.940	0.2	135	-0.01
5	S	2-Fluorophenol	80.000	84.386	-5.5	152	-0.01
6		Aniline	40.000	42.149	-5.4	148	-0.01
7	S	Phenol-d6	80.000	85.087	-6.4	152	-0.01
8		2-Chlorophenol	40.000	41.693	-4.2	148	-0.01
9		Benzaldehyde	40.000	40.026	-0.1	144	-0.01
10	C	Phenol	40.000	41.975	-4.9	148	0.00
11		bis(2-Chloroethyl)ether	40.000	42.200	-5.5	152	0.00
12		1,3-Dichlorobenzene	40.000	41.143	-2.9	148	-0.01
13	C	1,4-Dichlorobenzene	40.000	39.909	0.2	145	0.00
14		1,2-Dichlorobenzene	40.000	41.562	-3.9	150	-0.01
15		Benzyl Alcohol	40.000	40.602	-1.5	148	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	41.632	-4.1	150	0.00
17		2-Methylphenol	40.000	42.970	-7.4	151	-0.01
18		Hexachloroethane	40.000	41.930	-4.8	148	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.551	-1.4	146	-0.01
20		3+4-Methylphenols	40.000	42.147	-5.4	146	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	147	0.00
22		Acetophenone	40.000	40.589	-1.5	147	-0.01
23	S	Nitrobenzene-d5	80.000	84.581	-5.7	153	-0.01
24		Nitrobenzene	40.000	42.300	-5.7	151	-0.01
25		Isophorone	40.000	40.296	-0.7	147	-0.01
26	C	2-Nitrophenol	40.000	44.784	-12.0	160	0.00
27		2,4-Dimethylphenol	40.000	40.265	-0.7	148	-0.01
28		bis(2-Chloroethoxy)methane	40.000	41.971	-4.9	154	-0.01
29	C	2,4-Dichlorophenol	40.000	41.372	-3.4	149	-0.01
30		1,2,4-Trichlorobenzene	40.000	39.533	1.2	143	-0.01
31		Naphthalene	40.000	40.986	-2.5	149	-0.01
32		Benzoic acid	40.000	47.627	-19.1	191	0.00
33		4-Chloroaniline	40.000	40.267	-0.7	148	-0.01
34	C	Hexachlorobutadiene	40.000	38.519	3.7	141	-0.01
35		Caprolactam	40.000	40.442	-1.1	152	-0.02
36	C	4-Chloro-3-methylphenol	40.000	40.693	-1.7	150	0.00
37		2-Methylnaphthalene	40.000	40.769	-1.9	149	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	147	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	39.626	0.9	147	-0.01
40	P	Hexachlorocyclopentadiene	40.000	36.439	8.9	130	-0.01
41	S	2,4,6-Tribromophenol	80.000	76.594	4.3	141	-0.01
42	C	2,4,6-Trichlorophenol	40.000	38.910	2.7	146	-0.01
43		2,4,5-Trichlorophenol	40.000	39.239	1.9	148	-0.01
44	S	2-Fluorobiphenyl	80.000	76.553	4.3	143	-0.01

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

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	38.713	3.2	142	-0.01
46		2-Chloronaphthalene	40.000	38.985	2.5	144	-0.01
47		2-Nitroaniline	40.000	42.665	-6.7	154	0.00
48		Acenaphthylene	40.000	38.733	3.2	144	-0.01
49		Dimethylphthalate	40.000	37.483	6.3	141	-0.01
50		2,6-Dinitrotoluene	40.000	41.815	-4.5	148	-0.01
51	C	Acenaphthene	40.000	39.291	1.8	145	-0.01
52		3-Nitroaniline	40.000	42.181	-5.5	152	-0.01
53	P	2,4-Dinitrophenol	40.000	39.657	0.9	163	-0.01
54		Dibenzofuran	40.000	37.828	5.4	141	-0.01
55	P	4-Nitrophenol	40.000	39.155	2.1	141	-0.01
56		2,4-Dinitrotoluene	40.000	43.074	-7.7	150	-0.01
57		Fluorene	40.000	37.729	5.7	141	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	38.000	5.0	137	-0.01
59		Diethylphthalate	40.000	36.116	9.7	137	-0.01
60		4-Chlorophenyl-phenylether	40.000	37.699	5.8	140	-0.01
61		4-Nitroaniline	40.000	39.578	1.1	142	-0.01
62		Azobenzene	40.000	37.644	5.9	140	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	136	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	42.147	-5.4	152	-0.01
65	c	n-Nitrosodiphenylamine	40.000	40.430	-1.1	139	-0.01
66		4-Bromophenyl-phenylether	40.000	40.306	-0.8	137	-0.02
67		Hexachlorobenzene	40.000	40.611	-1.5	139	-0.01
68		Atrazine	40.000	39.190	2.0	132	-0.01
69	C	Pentachlorophenol	40.000	43.780	-9.5	148	-0.01
70		Phenanthrene	40.000	39.541	1.1	135	-0.01
71		Anthracene	40.000	39.829	0.4	136	-0.01
72		Carbazole	40.000	38.788	3.0	132	-0.01
73		Di-n-butylphthalate	40.000	39.122	2.2	133	-0.02
74	C	Fluoranthene	40.000	38.736	3.2	131	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	132	-0.02
76		Benzidine	40.000	38.164	4.6	122	-0.02
77		Pyrene	40.000	40.201	-0.5	130	-0.01
78	S	Terphenyl-d14	80.000	78.073	2.4	126	-0.02
79		Butylbenzylphthalate	40.000	41.077	-2.7	133	-0.02
80		Benzo(a)anthracene	40.000	39.655	0.9	131	-0.02
81		3,3'-Dichlorobenzidine	40.000	38.914	2.7	126	-0.02
82		Chrysene	40.000	40.205	-0.5	132	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	40.790	-2.0	135	-0.02
84	c	Di-n-octyl phthalate	40.000	42.282	-5.7	137	-0.04
85		Indeno(1,2,3-cd)pyrene	40.000	42.784	-7.0	141	-0.07
86	I	Perylene-d12	20.000	20.000	0.0	137	-0.03
87		Benzo(b)fluoranthene	40.000	38.899	2.8	134	-0.04

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.432	1.4	137	-0.03
89 C	Benzo(a)pyrene	40.000	39.760	0.6	137	-0.04
90	Dibenzo(a,h)anthracene	40.000	40.312	-0.8	141	-0.05
91	Benzo(g,h,i)perylene	40.000	40.920	-2.3	141	-0.07

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