Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060615\
Data File : BG017537.D
Acq On : 7 Jun 2015 16:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 05:29:36 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	143	0.00
2		1,4-Dioxane	40.000	37.265	6.8	140	-0.01
3		Pyridine	40.000	35.359	11.6	121	-0.01
4		n-Nitrosodimethylamine	40.000	45.247	-13.1	154	-0.01
5	S	2-Fluorophenol	80.000	80.365	-0.5	142	0.00
6		Aniline	40.000	38.727	3.2	138	0.00
7	S	Phenol-d6	80.000	79.907	0.1	139	0.00
8		2-Chlorophenol	40.000	39.502	1.2	142	0.00
9		Benzaldehyde	40.000	38.425	3.9	131	0.00
10	C	Phenol	40.000	39.944	0.1	137	0.00
11		bis(2-Chloroethyl)ether	40.000	38.784	3.0	136	0.00
12		1,3-Dichlorobenzene	40.000	41.400	-3.5	149	0.00
13	C	1,4-Dichlorobenzene	40.000	41.504	-3.8	149	0.00
14		1,2-Dichlorobenzene	40.000	40.582	-1.5	145	0.00
15		Benzyl Alcohol	40.000	38.183	4.5	130	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.389	1.5	139	0.00
17		2-Methylphenol	40.000	38.730	3.2	135	0.00
18		Hexachloroethane	40.000	43.644	-9.1	153	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.238	9.4	128	0.00
20		3+4-Methylphenols	40.000	39.046	2.4	135	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	141	0.00
22		Acetophenone	40.000	39.007	2.5	135	0.00
23	S	Nitrobenzene-d5	80.000	83.379	-4.2	142	0.00
24		Nitrobenzene	40.000	40.845	-2.1	141	0.00
25		Isophorone	40.000	34.999	12.5	123	0.00
26	C	2-Nitrophenol	40.000	39.432	1.4	133	0.00
27		2,4-Dimethylphenol	40.000	38.633	3.4	134	0.00
28		bis(2-Chloroethoxy)methane	40.000	36.898	7.8	125	0.00
29	C	2,4-Dichlorophenol	40.000	40.175	-0.4	136	0.00
30		1,2,4-Trichlorobenzene	40.000	40.926	-2.3	144	0.00
31		Naphthalene	40.000	39.027	2.4	136	0.00
32		Benzoic acid	40.000	34.127	14.7	115	0.04
33		4-Chloroaniline	40.000	35.935	10.2	124	0.00
34	C	Hexachlorobutadiene	40.000	45.199	-13.0	157	0.00
35		Caprolactam	40.000	29.071	27.3#	95	0.00
36	C	4-Chloro-3-methylphenol	40.000	35.713	10.7	120	0.01
37		2-Methylnaphthalene	40.000	37.977	5.1	130	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	125	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	45.730	-14.3	145	0.00
40	P	Hexachlorocyclopentadiene	40.000	15.110	62.2#	31	0.00
41	S	2,4,6-Tribromophenol	80.000	74.607	6.7	113	0.00
42	C	2,4,6-Trichlorophenol	40.000	41.700	-4.3	121	0.00
43		2,4,5-Trichlorophenol	40.000	40.570	-1.4	126	0.00
44	S	2-Fluorobiphenyl	80.000	86.939	-8.7	134	0.00

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Quant Time: Jun 08 05:29:36 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	42.922	-7.3	130	0.00
46		2-Chloronaphthalene	40.000	43.169	-7.9	128	0.00
47		2-Nitroaniline	40.000	40.429	-1.1	120	0.00
48		Acenaphthylene	40.000	40.216	-0.5	122	0.00
49		Dimethylphthalate	40.000	35.773	10.6	108	0.00
50		2,6-Dinitrotoluene	40.000	34.889	12.8	107	0.00
51	C	Acenaphthene	40.000	40.167	-0.4	124	0.00
52		3-Nitroaniline	40.000	34.699	13.3	105	0.00
53	P	2,4-Dinitrophenol	40.000	11.635	70.9#	18	0.00
54		Dibenzofuran	40.000	39.119	2.2	118	0.00
55	P	4-Nitrophenol	40.000	30.791	23.0	92	0.02
56		2,4-Dinitrotoluene	40.000	34.837	12.9	101	0.00
57		Fluorene	40.000	39.964	0.1	121	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.292	1.8	114	0.00
59		Diethylphthalate	40.000	34.804	13.0	106	0.00
60		4-Chlorophenyl-phenylether	40.000	39.960	0.1	122	0.00
61		4-Nitroaniline	40.000	29.701	25.7#	85	0.00
62		Azobenzene	40.000	39.557	1.1	119	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	40.000	10.377	74.1#	27	0.00
65	c	n-Nitrosodiphenylamine	40.000	42.998	-7.5	110	0.00
66		4-Bromophenyl-phenylether	40.000	43.970	-9.9	117	0.00
67		Hexachlorobenzene	40.000	42.774	-6.9	113	0.00
68		Atrazine	40.000	37.924	5.2	96	0.00
69	C	Pentachlorophenol	40.000	38.941	2.6	103	0.00
70		Phenanthrene	40.000	40.929	-2.3	108	0.00
71		Anthracene	40.000	40.975	-2.4	109	0.00
72		Carbazole	40.000	38.323	4.2	101	0.00
73		Di-n-butylphthalate	40.000	38.871	2.8	101	0.00
74	C	Fluoranthene	40.000	39.027	2.4	103	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
76		Benzidine	40.000	38.051	4.9	94	0.00
77		Pyrene	40.000	39.398	1.5	103	0.00
78	S	Terphenyl-d14	80.000	81.376	-1.7	107	0.00
79		Butylbenzylphthalate	40.000	40.384	-1.0	104	0.00
80		Benzo(a)anthracene	40.000	39.811	0.5	104	0.00
81		3,3'-Dichlorobenzidine	40.000	40.652	-1.6	104	0.00
82		Chrysene	40.000	39.530	1.2	103	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	41.584	-4.0	108	0.00
84	c	Di-n-octyl phthalate	40.000	40.912	-2.3	108	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	35.216	12.0	92	0.02
86	I	Perylene-d12	20.000	20.000	0.0	107	0.01
87		Benzo(b)fluoranthene	40.000	39.342	1.6	102	0.00

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

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	40.731	-1.8	111	0.01
89 C	Benzo(a)pyrene	40.000	38.807	3.0	102	0.01
90	Dibenzo(a,h)anthracene	40.000	34.387	14.0	90	0.01
91	Benzo(g,h,i)perylene	40.000	28.563	28.6#	76	0.01

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