Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
Data File : BG018020.D
Acq On : 21 Jul 2015 13:03
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 22 01:49:58 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 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	92	-0.02
2		1,4-Dioxane	40.000	31.622	20.9	78	-0.03
3		Pyridine	40.000	33.394	16.5	77	-0.02
4		n-Nitrosodimethylamine	40.000	30.701	23.2	74	-0.02
5	S	2-Fluorophenol	80.000	74.418	7.0	88	-0.02
6		Aniline	40.000	35.793	10.5	85	-0.02
7	S	Phenol-d6	80.000	72.585	9.3	86	-0.01
8		2-Chlorophenol	40.000	38.071	4.8	91	-0.02
9		Benzaldehyde	40.000	34.661	13.3	83	-0.02
10	C	Phenol	40.000	35.584	11.0	85	-0.01
11		bis(2-Chloroethyl)ether	40.000	35.292	11.8	85	-0.02
12		1,3-Dichlorobenzene	40.000	37.569	6.1	90	-0.02
13	C	1,4-Dichlorobenzene	40.000	37.465	6.3	90	-0.02
14		1,2-Dichlorobenzene	40.000	37.104	7.2	90	-0.02
15		Benzyl Alcohol	40.000	36.100	9.7	86	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	30.246	24.4	75	-0.02
17		2-Methylphenol	40.000	36.398	9.0	87	-0.01
18		Hexachloroethane	40.000	36.334	9.2	87	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	33.764	15.6	81	-0.02
20		3+4-Methylphenols	40.000	37.455	6.4	88	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	88	-0.02
22		Acetophenone	40.000	37.328	6.7	86	-0.02
23	S	Nitrobenzene-d5	80.000	74.446	6.9	84	-0.02
24		Nitrobenzene	40.000	37.054	7.4	83	-0.02
25		Isophorone	40.000	36.972	7.6	83	-0.02
26	C	2-Nitrophenol	40.000	47.770	-19.4	110	-0.02
27		2,4-Dimethylphenol	40.000	40.247	-0.6	92	-0.02
28		bis(2-Chloroethoxy)methane	40.000	36.505	8.7	82	-0.02
29	C	2,4-Dichlorophenol	40.000	44.208	-10.5	97	-0.01
30		1,2,4-Trichlorobenzene	40.000	40.777	-1.9	95	-0.02
31		Naphthalene	40.000	38.102	4.7	87	-0.02
32		Benzoic acid	40.000	55.935	-39.8#	166	0.00
33		4-Chloroaniline	40.000	38.950	2.6	87	-0.02
34	C	Hexachlorobutadiene	40.000	42.755	-6.9	96	-0.02
35		Caprolactam	40.000	40.701	-1.8	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.116	-0.3	89	0.00
37		2-Methylnaphthalene	40.000	38.347	4.1	88	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	89	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	40.888	-2.2	94	-0.02
40	P	Hexachlorocyclopentadiene	40.000	46.439	-16.1	108	-0.02
41	S	2,4,6-Tribromophenol	80.000	93.230	-16.5	109	-0.01
42	C	2,4,6-Trichlorophenol	40.000	44.710	-11.8	104	-0.02
43		2,4,5-Trichlorophenol	40.000	45.045	-12.6	102	0.00
44	S	2-Fluorobiphenyl	80.000	78.525	1.8	90	-0.01

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 2015
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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	38.864	2.8	89	-0.01
46		2-Chloronaphthalene	40.000	38.497	3.8	89	-0.02
47		2-Nitroaniline	40.000	37.842	5.4	87	-0.01
48		Acenaphthylene	40.000	38.769	3.1	88	-0.01
49		Dimethylphthalate	40.000	39.739	0.7	90	-0.01
50		2,6-Dinitrotoluene	40.000	42.323	-5.8	97	-0.01
51	C	Acenaphthene	40.000	37.923	5.2	88	-0.01
52		3-Nitroaniline	40.000	41.938	-4.8	93	-0.01
53	P	2,4-Dinitrophenol	40.000	54.030	-35.1#	146	0.00
54		Dibenzofuran	40.000	38.208	4.5	88	-0.01
55	P	4-Nitrophenol	40.000	41.821	-4.6	96	0.00
56		2,4-Dinitrotoluene	40.000	43.068	-7.7	99	-0.01
57		Fluorene	40.000	38.472	3.8	89	-0.01
58		2,3,4,6-Tetrachlorophenol	40.000	44.293	-10.7	102	0.00
59		Diethylphthalate	40.000	38.962	2.6	89	-0.01
60		4-Chlorophenyl-phenylether	40.000	39.763	0.6	91	-0.01
61		4-Nitroaniline	40.000	38.238	4.4	89	0.00
62		Azobenzene	40.000	32.920	17.7	75	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	89	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	51.456	-28.6#	129	-0.01
65	c	n-Nitrosodiphenylamine	40.000	39.261	1.8	88	-0.01
66		4-Bromophenyl-phenylether	40.000	41.613	-4.0	95	-0.02
67		Hexachlorobenzene	40.000	41.030	-2.6	96	-0.01
68		Atrazine	40.000	41.857	-4.6	93	0.00
69	C	Pentachlorophenol	40.000	48.435	-21.1#	113	-0.01
70		Phenanthrene	40.000	37.750	5.6	87	-0.02
71		Anthracene	40.000	39.062	2.3	87	-0.01
72		Carbazole	40.000	38.462	3.8	87	-0.01
73		Di-n-butylphthalate	40.000	40.219	-0.5	89	-0.01
74	C	Fluoranthene	40.000	39.117	2.2	89	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	89	-0.01
76		Benzidine	40.000	40.357	-0.9	91	-0.01
77		Pyrene	40.000	38.150	4.6	88	-0.01
78	S	Terphenyl-d14	80.000	79.411	0.7	92	-0.01
79		Butylbenzylphthalate	40.000	41.466	-3.7	91	-0.01
80		Benzo(a)anthracene	40.000	39.122	2.2	89	-0.01
81		3,3'-Dichlorobenzidine	40.000	44.174	-10.4	99	-0.01
82		Chrysene	40.000	38.774	3.1	89	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	41.543	-3.9	91	-0.02
84	c	Di-n-octyl phthalate	40.000	43.631	-9.1	94	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	39.818	0.5	91	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	93	-0.02
87		Benzo(b)fluoranthene	40.000	40.071	-0.2	96	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	36.204	9.5	85	-0.01
89 C	Benzo(a)pyrene	40.000	38.833	2.9	91	-0.02
90	Dibenzo(a,h)anthracene	40.000	39.103	2.2	94	-0.02
91	Benzo(g,h,i)perylene	40.000	38.910	2.7	93	-0.02

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