Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
Data File : BM019260.D
Acq On : 20 Mar 2019 09:07
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 20 11:08:51 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 11 17:30:46 2019
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	121	-0.02
2		1,4-Dioxane	40.000	33.067	17.3	106	-0.02
3		Pyridine	40.000	39.309	1.7	114	-0.02
4		n-Nitrosodimethylamine	40.000	35.453	11.4	107	-0.02
5	S	2-Fluorophenol	80.000	74.040	7.4	113	-0.02
6		Aniline	40.000	38.798	3.0	118	-0.02
7	S	Phenol-d6	80.000	78.892	1.4	119	-0.02
8		2-Chlorophenol	40.000	38.833	2.9	120	-0.03
9		Benzaldehyde	40.000	36.609	8.5	111	-0.02
10	C	Phenol	40.000	40.015	-0.0	121	-0.02
11		bis(2-Chloroethyl)ether	40.000	38.276	4.3	116	-0.02
12		1,3-Dichlorobenzene	40.000	37.390	6.5	116	-0.03
13	C	1,4-Dichlorobenzene	40.000	37.411	6.5	117	-0.02
14		1,2-Dichlorobenzene	40.000	37.632	5.9	117	-0.02
15		Benzyl Alcohol	40.000	39.750	0.6	119	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	37.308	6.7	116	-0.03
17		2-Methylphenol	40.000	39.820	0.4	121	-0.02
18		Hexachloroethane	40.000	37.432	6.4	114	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	39.380	1.5	118	-0.02
20		3+4-Methylphenols	40.000	40.928	-2.3	122	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	126	-0.03
22		Acetophenone	40.000	38.048	4.9	119	-0.03
23	S	Nitrobenzene-d5	80.000	75.483	5.6	119	-0.02
24		Nitrobenzene	40.000	37.495	6.3	118	-0.02
25		Isophorone	40.000	38.354	4.1	120	-0.02
26	C	2-Nitrophenol	40.000	40.304	-0.8	122	-0.03
27		2,4-Dimethylphenol	40.000	38.991	2.5	122	-0.02
28		bis(2-Chloroethoxy)methane	40.000	37.968	5.1	120	-0.02
29	C	2,4-Dichlorophenol	40.000	40.971	-2.4	126	-0.02
30		1,2,4-Trichlorobenzene	40.000	37.940	5.2	122	-0.03
31		Naphthalene	40.000	37.382	6.5	120	-0.02
32		Benzoic acid	40.000	38.186	4.5	119	0.00
33		4-Chloroaniline	40.000	39.813	0.5	124	-0.02
34	C	Hexachlorobutadiene	40.000	37.575	6.1	121	-0.03
35		Caprolactam	40.000	42.386	-6.0	128	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.113	-0.3	125	-0.02
37		2-Methylnaphthalene	40.000	38.342	4.1	124	-0.03
38		1-Methylnaphthalene	40.000	38.196	4.5	122	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	132	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	37.265	6.8	124	-0.02
41	P	Hexachlorocyclopentadiene	40.000	29.426	26.4#	96	-0.02
42	S	2,4,6-Tribromophenol	80.000	80.334	-0.4	131	-0.02
43	C	2,4,6-Trichlorophenol	40.000	39.448	1.4	128	-0.02
44		2,4,5-Trichlorophenol	40.000	39.768	0.6	128	-0.02

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Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

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

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	74.351	7.1	125	-0.02
46		1,1'-Biphenyl	40.000	36.914	7.7	124	-0.02
47		2-Chloronaphthalene	40.000	37.176	7.1	125	-0.02
48		2-Nitroaniline	40.000	38.868	2.8	122	-0.02
49		Acenaphthylene	40.000	37.641	5.9	125	-0.02
50		Dimethylphthalate	40.000	36.736	8.2	124	-0.02
51		2,6-Dinitrotoluene	40.000	38.469	3.8	124	-0.02
52	C	Acenaphthene	40.000	36.770	8.1	124	-0.02
53		3-Nitroaniline	40.000	37.337	6.7	124	-0.02
54	P	2,4-Dinitrophenol	40.000	30.401	24.0	99	-0.02
55		Dibenzofuran	40.000	37.198	7.0	126	-0.02
56	P	4-Nitrophenol	40.000	36.708	8.2	120	-0.02
57		2,4-Dinitrotoluene	40.000	38.056	4.9	128	-0.02
58		Fluorene	40.000	37.922	5.2	127	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	38.215	4.5	125	-0.02
60		Diethylphthalate	40.000	37.232	6.9	124	-0.02
61		4-Chlorophenyl-phenylether	40.000	38.014	5.0	128	-0.02
62		4-Nitroaniline	40.000	37.215	7.0	124	-0.02
63		Azobenzene	40.000	37.722	5.7	123	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	132	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	33.840	15.4	111	-0.02
66	c	n-Nitrosodiphenylamine	40.000	37.941	5.1	125	-0.02
67		4-Bromophenyl-phenylether	40.000	38.552	3.6	127	-0.02
68		Hexachlorobenzene	40.000	38.655	3.4	129	-0.02
69		Atrazine	40.000	37.391	6.5	122	-0.02
70	C	Pentachlorophenol	40.000	37.171	7.1	123	-0.02
71		Phenanthrene	40.000	37.088	7.3	125	-0.02
72		Anthracene	40.000	38.225	4.4	126	-0.02
73		Carbazole	40.000	37.302	6.7	123	-0.02
74		Di-n-butylphthalate	40.000	39.278	1.8	127	-0.02
75	C	Fluoranthene	40.000	36.262	9.3	122	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	117	-0.02
77		Benzidine	40.000	34.776	13.1	101	-0.02
78		Pyrene	40.000	42.284	-5.7	121	-0.02
79	S	Terphenyl-d14	80.000	86.847	-8.6	125	-0.02
80		Butylbenzylphthalate	40.000	44.886	-12.2	124	-0.02
81		Benzo(a)anthracene	40.000	37.617	6.0	111	-0.02
82		3,3'-Dichlorobenzidine	40.000	38.581	3.5	111	-0.02
83		Chrysene	40.000	37.136	7.2	111	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	45.623	-14.1	128	-0.01
85	c	Di-n-octyl phthalate	40.000	42.125	-5.3	123	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	33.431	16.4	112	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	110	-0.02

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Instrument : BNA_M
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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	37.534	6.2	104	-0.02
89	Benzo(k)fluoranthene	40.000	39.208	2.0	105	-0.02
90 C	Benzo(a)pyrene	40.000	37.520	6.2	104	-0.03
91	Dibenzo(a,h)anthracene	40.000	38.633	3.4	112	-0.04
92	Benzo(g,h,i)perylene	40.000	39.099	2.3	113	-0.04

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