

Online Appendix – Objective Bayesian Edge Screening and Structure Selection for Ising Networks

This is the online appendix for the paper "Objective Bayesian Edge Screening and Structure Selection for Ising Networks."

A Comparison of Bayesian Approaches to Structure Selection for Networks of Binary Data

We have introduced novel edge screening and structure selection procedures for estimating the topology of Ising networks. As discussed in our paper, other Bayesian approaches exist for structure selection for the Ising model (Pensar et al., 2017)—implemented in the R-package `BDgraph` (Mohammadi and Wit, 2019)—and the multivariate probit model (Talhok et al., 2012)—implemented in the R-package `BGGM` (Williams and Mulder, 2020b). We wish to compare the performance of these two approaches to that of our procedures in terms of edge-selection sensitivity and specificity.

Similar to the other simulations in our paper, we focus on simulations in which and Erdős and Rényi (1960) model is used to generate the topologies. The parameters corresponding to present edges are simulated uniformly between 0.5 and 2, and set to zero otherwise. In the simulations, we vary π , the probability of generating an edge between two variables, p , the number of variables, and n , the number of observations and generate 100 datasets for each combination of values for π , p , and n .

We analyze the simulated datasets using the default settings of `rbinnet` (`precision = .975`), `BDgraph`, and `BGGM`, except that for the latter we varied the specification of the `delta` parameter. We have used `delta` values of 5, and 24, which translate into a prior standard deviation for the partial correlation between the latent Gaussian variables of approximately .41, and .2, respectively. These `delta` values were suggested in Williams and Mulder (2020a). We ran each of the sampling-based methods—structure selection, `BDgraph` and `BGGM`—for 10,000 iterations. Sensitivity and Specificity were computed as defined in our paper. Table 1 shows the results for a generating probability $\pi = 0.1$ and Table 2 shows the results for $\pi = 0.5$.

With increasing sample size, the sensitivity increases for all methods. In Table 1, sensitivity starts around 13% for most methods (36% for `BGGM`, `delta = 24`), it increases to around 80% (86% for `rbinnet`, edge screening). In Table 2, which considers denser networks, sensitivity starts around 85% and increases to around 100%. The same observation could be made for specificity. In Table 1, specificity starts around 95% for most methods (87% for `BGGM`, `delta = 24`), it increases to 95% and above for all methods. In Table 2, specificity already starts above 95% and slowly increases to over 99% for all methods, except that

		<i>n</i>			
		250	500	1,000	10,000
Screening	SEN	.133	.193	.438	.859
	SPE	.954	.959	.966	.953
Selection	SEN	.141	.167	.394	.807
	SPE	.941	.964	.978	.987
Screening & Selection	SEN	.133	.172	.394	.807
	SPE	.954	.963	.975	.987
BDgraph	SEN	.117	.151	.306	.808
	SPE	.949	.976	.988	.988
BGGM (<code>delta</code> = 5)	SEN	.133	.163	.317	.802
	SPE	.941	.972	.985	.995
BGGM (<code>delta</code> = 24)	SEN	.362	.297	.446	.802
	SPE	.874	.921	.959	.996

Table 1

Sensitivity and specificity for $\pi = 0.1$, as a measure of performance of the methods implemented in `rbinnet`, `BDgraph`, and `BGGM`.

for edge screening it appears to hover around 97%. There are no major differences in the performance of the different methods, except that for smaller sample sizes BGGM—using a `delta` value equal to 24—showed an increased sensitivity and reduced specificity.

Computing Time Differences Between the Methods and Their Implementations

The methods are currently implemented in the `rbinnet` package using native R code. We therefore expect it to run considerably slower than other packages, such as `BGGM` and `BDgraph`, which run code in a compiled language. To get a sense of the differences in running time between the different methods and implementation, we simulated data for $n = 200, 500$, and $1,000$ cases for a $p = 10$, and 15 variable network and ran the different procedures. The samplers each ran for 10,000 iterations. Simulations ran on a MacBook pro with an Intel i5 CPU and clock speed 2.0 Ghz and 16Gb of memory.

The results are shown in Table 3. Observe that edge screening and eLasso are both really fast. But these only do optimization. When comparing the simulation based methods, we observe that structure selection is much slower than both `BGGM` and `BDgraph`. Whereas `BGGM` is roughly four to six times faster than structure selection for the $p = 10$ variable network, it is seven to eight times faster for the $p = 15$ variable network. Thus, `BGGM` becomes relatively faster than structure selection for increasing network sizes. Interestingly, the relative speed difference shrinks with increasing sample sizes. `BDgraph`, on the other hand, is roughly 30 to 50 times faster than structure selection. In contrast to the implementation of `BGGM`, `BDgraph` does not become relatively faster than structure selection when the network’s size increases, although it does appear to become relatively faster for increasing sample sizes.

		<i>n</i>			
		250	500	1,000	10,000
Screening	SEN	.846	.966	.994	1.000
	SPE	.978	.959	.971	.964
Selection	SEN	.859	.966	.991	1.000
	SPE	.970	.957	.983	.995
Screening & Selection	SEN	.846	.960	.994	1.000
	SPE	.978	.962	.983	.995
BDgraph	SEN	.815	.895	.948	1.000
	SPE	.977	.992	.995	.999
BGGM ($\delta = 5$)	SEN	.816	.934	.977	1.000
	SPE	.984	.994	.989	.992
BGGM ($\delta = 24$)	SEN	.882	.957	.986	1.000
	SPE	.956	.988	.975	.989

Table 2

Sensitivity and specificity for $\pi = 0.5$, as a measure of performance of the methods implemented in `rbinnet`, `BDgraph`, and `BGGM`.

<i>p</i>	<i>n</i>	Screening	Selection	BGGM	BDgraph	IsingFit
10	200	0.4	55.4	10.1	1.4	0.3
	500	0.5	84.1	19.7	2.0	0.3
	1,000	0.7	166.4	38.9	4.9	0.4
15	200	1.1	150.8	19.0	4.7	0.5
	500	1.3	342.2	44.7	7.7	0.6
	1,000	2.2	640.0	88.9	13.5	0.8

Table 3

Running time in seconds for the different methods and implementations in `rbinnet`, `BDgraph`, `BGGM`, and `IsingFit`. The simulation based methods ran for 10,000 iterations each.

References

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