Efficient Bayesian Social Learning on Trees

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October 29, 2013

Abstract

We consider a set of agents who are attempting to iteratively learn the 'state of the world' from their neighbors in a social network. Each agent initially receives a noisy observation of the true state of the world. The agents then repeatedly 'vote' and observe the votes of some of their peers, from which they gain more information. The agents' calculations are Bayesian and aim to myopically maximize the expected utility at each iteration.

This model, introduced by Gale and Kariv (2003), is a natural approach to learning on networks. However, it has been criticized, chiefly because the agents' decision rule appears to become computationally intractable as the number of iterations advances. For instance, a dynamic programming approach (part of this work) has running time that is exponentially large in min(n, (d − 1)t), where n is the number of agents.

We provide a new algorithm to perform the agents' computations on locally tree-like graphs. Our algorithm uses the dynamic cavity method to drastically reduce computational effort. Let d be the maximum degree and t be the iteration number. The computational effort needed per agent is exponential only in O(td) (note that the number of possible information sets of a neighbor at time t is itself exponential in td).

Under appropriate assumptions on the rate of convergence, we deduce that each agent is only required to spend polylogarithmic (in 1/ϵ) computational effort to approximately learn the true state of the world with error probability ϵ, on regular trees of degree at least five. We provide numerical and other evidence to justify our assumption on convergence rate.

We extend our results in various directions, including loopy graphs. Our results indicate efficiency of iterative Bayesian social learning in a wide range of situations, contrary to widely held beliefs.

1 Introduction

Consider a group of Facebook users who are each faced with the dilemma of whether to place an order for the new iGadget or the new Gadgetoid. Each boldly ventures to the wild and does independent research on the subject matter, discovering the “correct” answer with some probability p > 1/2. Then, over the next few weeks, before making the final decision, they daily share their current opinion on the matter with their Facebook contacts by posting either iGadget or Gadgetoid on their status line. Every day, after learning their friends’ opinions, they update their own by performing the Bayesian calculation that determines which of the two options is more likely to be true, given all they know. Eventually, they make a purchase based on this information. Such dynamics have become an integral part of electronic commerce, and understanding them is valuable to social media advertisers and vendors.

This model (or rather, a slightly more general version of it) was introduced by Gale and Kariv [7]. It is one in a long succession of social learning models. Already in 1785 Condorcet [4] considered how a group of individuals with weak private signals could reach a correct collective decision; he showed that a majority vote is likely to be correct when the group is large enough. Models such as those of Banerjee [2], Bikhchandani, Hirshleifer and Welch [3] and Smith and Sorensen [11] allow for each individual to make a single decision, learning from the decisions of her predecessors. The models of DeGroot [5] and Bala and Goyal [1] consider social networks and repeated interactions between agents.

The model of Gale and Kariv combines features from all of the above. It describes a group of individuals (or agents), each with a private signal that carries information on an unknown state of the world. The individuals form a social network, so that each observes the actions of some subset - her neighbors. The agents must choose between a set of possible actions, the relative merit of which depends on the state of the world. The agents iteratively learn by observing their neighbors’ actions, and picking an action that
is myopically optimal, given the information known to them.

Even in the simple case of two states of the world, two possible private signals and two actions, the required calculations appear to be very complicated. This has indeed been a recurring criticism of this model (see, e.g., [9][12]). One approach to this difficulty is the bounded rationality approach of Bala and Goyal [10], where agents ignore part of the information available to them and perform a Bayesian calculation on the rest.

While the bounded rationality approach has led to impressive results, it has two disadvantages, as compared with a fully Bayesian one: first, it is bound to involve a somewhat arbitrary decision of which heuristics the agents use. Second, a game theoretic analysis of strategic players is possible only if the players choose actions that are optimal by some criterion. Hence game-theoretic analyses of learning on networks (e.g. [12]) often opt for the more difficult but fully Bayesian model.

A different approach to the difficulty of computation in the Bayesian model is to show that the calculations are in fact not as difficult as they appear, at least in some cases. In this paper we show that when the graph of social ties is locally a tree, or close to one, then the computational outlook is not a bleak as previously thought.

We first give a simple dynamic programming algorithm for the Gale and Kariv model that is exponential in the number of individuals. Since at iteration $t$ one may consider only agents at distance $t$, then in graphs of maximum degree $d$ (on which we focus) the number of individuals to consider is $O((d-1)^t)$, and the time required of each individual to compute their action (or vote) at time $t$ is $2^{O((d-1)^t)}$. We then develop a sophisticated dynamic program for locally tree-like graphs that reduces the computational effort to $2^{O(td)}$.

We conjecture, and show supporting numerical evidence, that on infinite trees of degree at least five, the number of iterations needed to calculate the correct answer with probability $1 - \epsilon$ is $O(\log \log(1/\epsilon))$. In fact, we rigorously establish this for the ‘majority dynamics’ update rule, in which agents adopt the opinion of their neighbors in the previous round. Thus, our conjecture follows if iterative Bayesian learning learns at least as fast as majority, as suggested by intuition and numerical evidence, which we present. Assuming this conjecture, the computational effort required drops from quasi-polynomial in $1/\epsilon$ (using the naive dynamic program) to polylogarithmic in $(1/\epsilon)$.

An additional difficulty of the Gale and Kariv model is that it requires the individuals to exactly know the structure of the graph. A possible solution to this is a modification that allows the agents to know only their own neighborhoods and the distribution from which the rest of the graph was picked. We pursue this for the natural configuration model of random graphs (see below for full explanation) and show that the same computational upper bounds apply here.

We also introduce two further features into the model and show how to deal with them algorithmically. First, there may be a finite number of ‘hub’ nodes who are each observed by many nodes leading to several short loops in the connectivity graph. We show that our algorithm can be suitably modified for this case. Second, we consider that nodes may not all be ‘active’ in each round, and that nodes may observe only a random subset of active neighbors. We show that this can be handled when ‘inactive’ edges/nodes occur independently of each other and in time.

The key technique used in this paper is the dynamic cavity method, introduced by Kanoria and Montanari [10] in their study of “recursive majority” updates on trees, which was also motivated by social learning. A dynamical version of the cavity method of Statistical Physics, this technique was used to analyze majority dynamics on trees, and appears promising for the analysis of iterative tree processes in general. In this work, we use this technique for the first time to give an algorithm for efficient computation by nodes. This is in contrast to the case of majority updates, where the update rule is computationally trivial. Our algorithmic approach leveraging the dynamic cavity method may be applicable to a range of iterative update situations on locally treelike graphs.

## 2 Model

- There is a true state of the world $s \in \mathcal{S}$, where $\mathcal{S}$ is finite. The prior distribution $P[s]$ is common knowledge.

- Let $G = (V,E)$ be an undirected connected graph of agents and their social ties. Let $n \equiv |V|$.

- Denote by $\partial i$ the neighbors of agent $i$, not including $i$.

- Each agent $i$ receives a private signal $x_i \in \mathcal{X}$, where $\mathcal{X}$ is finite. Private signals are independent conditioned on $s$. The distribution $P[x_i | s]$ is common knowledge. We assume that the signal is informative, so that $P[x_i | s]$ is different for different values of $s$. 
• We identify the set of actions available to agents with the set $S$ of the states of the world (thus we call the actions ‘votes’). For each state of the world $s$, action $\sigma$ has utility one when the state of the world is $s = \sigma$, and zero otherwise. Thus the action that maximizes the expected utility corresponds to the maximum \textit{a posteriori} probability (MAP) estimator of the state of the world.

• At each time period $t \in \{0, 1, 2, \ldots\}$ each agent takes an action and then observes the actions taken by her neighbors.

• Denote by $F_t^i$ the information available to agent $i$ at time $t$. We do not include in this her neighbors’ votes at time $t$.

• At each time period, the agents’ goal is to maximize their expected utility. They are myopic and at each iteration cross out entries that are inconsistent with the signals that they’ve observed from their neighbors up to that point. Then, they calculate the probabilities of the different possible states of the world by summing over the entries that have yet to be crossed out.

This may not be as simple as it seems. To understand which initial configurations are ruled out by a signal coming from a neighbor, an agent must “simulate” that neighbor’s behavior, and so each agent must calculate the function $g_t^i$ for every other agent $i$ and every possible set of observations by $i$. We formalize this below.

Let $x \in \mathcal{X}^n$ be the vector of private signals $(x_i)_{i \in V}$. The trajectory of $i$, $\sigma_i$, is a deterministic function of $x_i$. Assume then that up to time $t - 1$ each agent has calculated the trajectory $\sigma_{t-1}^i(x)$ for all possible private signal vectors $x$ and all agents $i$. This is trivial for $t - 1 = 0$.

We say that $y$ is feasible for $i$ at time $t$ if $x_i = y_i$ and $\sigma_{t-1}^i = \sigma_{t-1}^j(y)$. We denote this set of feasible private signal vectors $I_t^i(x_i, \sigma_{t-1}^i)$. To calculate $\sigma_t^i(x)$, one need only note that

$$\mathbb{P}[s|F_t^i] \propto \mathbb{P}[s|x_i, \sigma_{t-1}^i|s] = \mathbb{P}[s] \sum_{y \in I_t^i} \mathbb{P}[y|s]$$

and

$$g_i(t, x_i, \sigma_{t-1}^i) = \arg\max_{s \in S} \mathbb{P}[s|F_t^i]$$

by definition. We use the standard abusive notation $\mathbb{P}[x_i]$ instead of $\mathbb{P}[x_i = y_i]$, $\mathbb{P}[\sigma_j]$ instead of $\mathbb{P}[\sigma_j = \omega_j]$, etc.

It is easy to verify that using this the calculation of each $\sigma_t^i(x)$ takes $O(tn|\mathcal{X}^n|)$. One can do better than perform each of these separately, but in any case the result is exponential in $n$, so we derive a rough upper bound of $2^{O(n)}$ for this method. Since we are in particular interested in graphs of maximum degree $d$, we note that up to time $t$ an agent need only perform this for agents at distance at most $t$, and so this bound becomes $2^{O((d-1)^t)}$ for large graphs, i.e., graphs for which $n > (d - 1)^{t}$ for relevant values of $t$.

3 A Simple Algorithm

A sign of the complexity of this Bayesian calculation is that even the brute-force solution for it is not trivial. We therefore describe it here.

One way of thinking of the agents’ calculation is to imagine that they keep a long list at the end of all the possible combinations of initial signals of all the other agents, and at each iteration cross out entries that are inconsistent with the signals that they’ve observed from their neighbors up to that point. Then, they calculate the probabilities of the different possible states of the world by summing over the entries that have yet to be crossed out.

4 The Dynamic Cavity Algorithm on Trees

Assume in this section that the graph $G$ is a tree with finite degree nodes. For $j \in \partial i$ let $G_{j-i} = (V_{j-i}, E_{j-i})$ denote $j$’s connected component in the graph $G$ with the edge $(i, j)$ removed. That is, $V_{j-i}$ is $j$’s subtree when $G$ is rooted in $i$. 

3
4.1 The Dynamic Cavity Method

We consider a modified process where agent \( i \) is replaced by a zombie which takes fixed actions \( \tau_i = (\tau_i(0), \tau_i(1), \ldots) \), and the true state of the world is assume to be some fixed \( s \). Furthermore, this ‘fixing’ goes unnoticed by the agents (except \( i \), who is a zombie anyway) who carry on their calculations, assuming \( i \) is her regular Bayesian self, and the state of the world is drawn randomly according to \( P[s] \). We denote by \( Q[A|\tau_i, s] \) the probability of event \( A \) in this modified process. This modified process is easier to analyze, as the processes on each of the subtrees \( V_{j \rightarrow i} \) are independent. This is formalized in the following claim, without proof:

**Claim 4.1.**

\[
Q[\sigma_{\partial i}^t|\tau_i, s] = \prod_{j \in \partial i} Q[\sigma_j^t|\tau_i^t, s]. \quad (1)
\]

(Since \( \sigma_j^t \) is unaffected by \( \tau_i(t') \) for all \( t' > t \), we only need to specify \( \tau_i^t \), and not the entire \( \tau_i \).

Now, it might so happen that for some number of steps the zombie behaves exactly as may be expected of a rational player. More precisely, given \( \sigma_{\partial i}^{t-1} \), it may be the case that \( \tau_i^t = g_i^t(x_i, \sigma_{\partial i}^{t-1}) \). This event provides the connection between the modified process and the original process, and is the inspiration for the following theorem.

**Theorem 4.2.** For all \( i, t \) and \( \tau_i \)

\[
P[\sigma_{\partial i}^{t-1}|s, x_i] 1 (\tau_i^t = g_i^t(x_i, \sigma_{\partial i}^{t-1})) = Q[\sigma_{\partial i}^{t-1}|\tau_i, s] 1 (\tau_i^t = g_i^t(x_i, \sigma_{\partial i}^{t-1})) . \quad (2)
\]

**Proof.** We couple the original process, after choosing \( s \), to the modified processes by setting the private signals to be identical in both.

Now, clearly if it so happens that \( \tau_i^t = g_i^t(x_i, \sigma_{\partial i}^{t-1}) \) then the two processes will be identical up to time \( t \). Hence the probabilities of events measurable up to time \( t \) will be identical when multiplied by \( 1 (\tau_i^t = g_i^t(x_i, \sigma_{\partial i}^{t-1})) \), and the theorem follows. \( \Box \)

Using Eqs. (1) and (2), we can easily write the posterior on \( s \) computed by node \( i \) at time \( t \), in terms of the probabilities \( Q[|s|] \):

\[
P[s|F_i^t] \propto P[s] P[x_i, \sigma_{\partial i}^{t-1}|s] = P[s] P[x_i|s] P[\sigma_{\partial i}^{t-1}|s, x_i] = P[s] P[x_i|s] \prod_{j \in \partial i} Q[\sigma_j^{t-1}|\sigma_i^{t-1}, s]. \quad (3)
\]

(Note that \( \sigma_i^{t-1} \) is a deterministic function of \( (x_i, \sigma_{\partial i}^{t-1}) \).)

Given that, the decision function is as before

\[
g_{i,t}(x_i, \sigma_{\partial i}^{t-1}) = \arg \max_{s \in S} P[s|F_i^t] \quad (4)
\]

As mentioned before, we assume there is a deterministic tie breaking rule that is common knowledge.

We are finally left with the task of calculating \( Q[|i|] \). The following theorem is the heart of the dynamic cavity method and allows us to perform this calculation:

**Theorem 4.3.** For \( j \in \partial i \) and \( t \in \mathbb{N} \)

\[
Q[\sigma_j^t||\tau_i, s] = \sum_{\sigma_j^{t-1}} \sum_{x_j} P[x_j|s] 1 \left[ \sigma_j^t = g_j^t(x_j, (\tau_i^{t-1}, \sigma_{\partial i}^{t-1})) \right] \cdot \prod_{l=1}^{d-1} Q[\sigma_l^{t-1}||\sigma_j^{t-1}, s]. \quad (5)
\]

where the neighbors of node \( j \) are \( \partial j = \{i, 1, 2, \ldots, d-1\} \).

We mention without proof that the recursion easily generalizes to the case of a random tie-breaking rule, provided the rule is common knowledge; it is a matter of replacing the expression \( 1[\sigma_j^t = \cdots] \) with \( P[\sigma_j^t = \cdots] \), where this probability is over the randomness of the rule. Eq. (3) continues to be valid in this case.

The following proof is similar to the proof of Lemma 2.1 in [10], where the dynamic cavity method is introduced and applied to a different process.

**Proof.** In the modified process the events in the different branches that \( i \) sees are independent. We therefore consider \( V_{j \rightarrow i} \) only, and view it as a tree rooted at \( j \). Also, for convenience we define \( \sigma_j^t \equiv \tau_i^t \); note that the random variable \( \sigma_j^t \) does not exist in the modified process, as \( i \)'s trajectory is fixed to \( \tau_i \).

Let \( x \) be the vector of private signals of \( j \) and all the vertices up to a distance \( t \) from \( j \) (call this set of vertices \( V^t_{j\rightarrow i} \)). For each \( l \in \{1, \ldots, d-1\} \), let \( x_l \) be the vector of private signals of \( V^t_{l \rightarrow j} \). Thus, \( x = (x_1, x_2, \ldots, x_{d-1}) \).

The trajectory \( \sigma_j^t \) is a function - deterministic, by our assumption - of \( x \) and \( \tau_i^t \). We shall denote this function by \( F_{j \rightarrow i} \) and write \( \sigma_j^t = F_{j \rightarrow i}(x, \tau_i^t) \). This function is uniquely determined by the update rules \( g_j^t(x_i, \sigma_{\partial i}^{t-1}) \) for \( l \in V^t_{j \rightarrow i} \).

We have therefore

\[
Q[\sigma_j^t = \lambda^t||\tau_i^t, s] = \sum_x P[x|s] 1(\lambda^t = F_{j \rightarrow i}(x, \tau_i^t)). \quad (6)
\]
We now analyze each of the terms appearing in this sum. Since the initialization is i.i.d., we have
\[ \mathbb{P}[\mathcal{E} \mid s] = \mathbb{P}[x_j|s] \mathbb{P}[x_1|s] \mathbb{P}[x_2|s] \cdots \mathbb{P}[x_{d-1}|s]. \] (7)

The function \( F_{j \to i}^{l} (\cdots) \) can be decomposed as follows:
\[ \mathbf{1} \left( \lambda^t = F_{j \to i}^{l} (\mathcal{E}, \tau^t_i) \right) = \sum_{\sigma_1^{t-1} \cdots \sigma_{d-1}^{t-1}} \mathbf{1} \left( \lambda^t = g_j^t (x_j, \sigma_1^{t-1}) \right) \cdot \prod_{l=1}^{d-1} \mathbf{1} \left( \sigma_l^{t-1} = F_{l \to j}^{t-1} (\mathcal{E}_l, \lambda^{t-1}) \right). \] (8)

Using Eqs. 7 and 8 in Eq. 9 and separating terms that depend only on \( \mathcal{E}_i \), we get
\[ \mathbb{Q} [\sigma_j^t = \lambda^t \mid \tau^t_i, s] = \sum_{\sigma_1^{t-1} \cdots \sigma_{d-1}^{t-1}} \sum_{x_j} \mathbb{P}[x_j|s] \mathbf{1} \left( \lambda^t = g_j^t (x_j, \sigma_1^{t-1}) \right) \cdot \prod_{l=1}^{d-1} \sum_{\mathcal{E}_l} \mathbb{P}[\mathcal{E}_l] \mathbf{1} \left( \sigma_l^{t-1} = F_{l \to j}^{t-1} (\mathcal{E}_l, \lambda^{t-1}) \right). \]
The recursion follows immediately by identifying that the product over \( l \) in fact has argument \( \mathbb{Q} [\sigma_j^{t-1} \mid \sigma_j^{t-1}, s] \).

4.2 The Agents’ Calculations

We now have in place all we need to perform the agent’s calculations. At time \( t = 0 \) these calculations are trivial. Assume then that up to time \( t \) each agent has calculated the following quantities:
1. \( \mathbb{Q} [\sigma_j^{t-1} \mid \tau_i^{t-1}, s] \) for all \( s \) and for all \( i, j \in V \) such that \( j \in \partial_i \), and for all \( \tau^{t-1} \) and \( \sigma_j^{t-1} \).
2. \( g_i^t (x_i, \sigma_1^{t-1}) \) for all \( i, x_i \) and \( \sigma_1^{t-1} \).

Note that these can be calculated without making any observations - only knowledge of the graph is needed.

At time \( t + 1 \) each agent makes the following calculations:
1. \( \mathbb{Q} [\sigma_j^t \mid \tau_i^t, s] \) for all \( s, i, j, \sigma_i^t, \tau_i^t \). These can be calculated using Eq. 9, given the quantities from the previous iteration.
2. \( g_i^{t+1} (x_i, \sigma_{\partial_i}^t) \) for all \( i, x_i \) and \( \sigma_{\partial_i}^t \). These can be calculated using Eqs. 7 and 8 and the newly calculated \( \mathbb{Q} [\sigma_j^t \mid \tau_i^t, s] \).

Since agent \( j \) calculates \( g_j^{t+1} \) for all \( i \), then she in particular calculates \( g_j^{t+1} \). Therefore, she can use this to calculate her next action, once she observes her neighbors’ actions. A simple calculation yields the following lemma.

**Lemma 4.4.** In a tree graph \( G \) with maximum degree \( d \), the agents can calculate their actions up to time \( t \) with computational effort \( n2^{O(td)} \).

In fact, each agent does not need to perform calculations for the entire graph. It suffices for node \( i \) to calculate quantities up to time \( t’ \) for nodes at distance \( t - t’ \) from node \( i \) (there are at most \((d - 1)^{t-t’}\) such nodes). A short calculation yields an improved bound on computational effort.

**Theorem 4.5.** In a tree graph \( G \) with maximum degree \( d \), each agent can calculate her action up to time \( t \) with computational effort \( 2^{O(td)} \).

4.3 Dynamic Cavity Algorithm: Extensions

Our algorithm admits several extensions that we explore in this section: Section 4.3.1 discusses graphs with loops and ‘hubs’, Section 4.3.2 discusses random graphs, Section 4.3.3 discusses graphs with loops and ‘hubs’, Section 4.3.4 relaxes the assumption that the entire graph is common knowledge and Section 4.3.5 allows nodes/edges to be inactive in some rounds.

First we mention some straightforward generalizations:

It is easy to see that dynamic cavity recursion (Theorem 4.3) does not depend on any special properties of the Bayesian update rule. The update rule \( g_i,t(\cdot) \) can be arbitrary. Thus, if agent \( i \) wants to perform a Bayesian update, he can do so (exactly) using our approach even if his neighbor, agent \( j \), is using some other update rule.

**Remark 4.6.** The dynamic cavity recursion can be used to enable computations of agents even if some of them are using arbitrary update rules (provided the rules are ‘well specified’ and common knowledge).

For instance, our approach should be applicable in ‘partial Bayesian’ settings.

Our algorithm is easily modified for the case of a general finite action set \( A \) that need not be the same as \( S \), associated with a payoff function \( u : A \times S \to \mathbb{R} \). Moreover the action set and payoff function can each be player dependent (\( A_i, u_i \) respectively).

We already mentioned that there is a simple generalization to the case of random tie breaking rules (that are common knowledge).

Instead of having only undirected edges (corresponding to bidirectional observations), we can allow a subset of the edges of the tree to be directed. In this case, the same algorithm works with suitably defined neighborhood sets \( \partial_i \). In other words, our result holds for the class of directed graphs lacking cycles of length greater than two (which correspond to undirected edges).
4.3.1 Loops and Hub nodes

A class of graphs that are not trees, but for which this dynamic cavity method can be easily extended is that of trees with ‘hub’ nodes in addition.

Consider then a graph that is not a tree, but can be transformed into a tree by the removal of some small set of nodes $V_{\text{loop}} \subset V$. Then the same calculations above can still be performed, with a time penalty of $|\mathcal{X}|^{\mid V_{\text{loop}}\mid}$; the calculation in Eq. (6) is simply repeated for each possible set of private signals of the hub nodes, and the probabilities in Eq. (6) are arrived at by averaging the $|\mathcal{X}|^{\mid V_{\text{loop}}\mid}$ different possible cases. In fact, one may not even need to average over all nodes in $V_{\text{loop}}$, since at iteration $t$ only those inside $B^t_i$ (the ball of radius $t$ around $i$) effect the outcome of $i$’s calculations. Hence the complexity of calculations up to iteration $t$ is now $|\mathcal{X}|^{n_t} 2^{O(td)}$, where $n_t = \max(|B^t_i \cap V_{\text{loop}}|) \in V$.

If we also allow directed edges in this model, then we can extend it to include nodes of unlimited indegree, i.e., some nodes may be observed by an unbounded number of others. These are agents who are observed by any number (perhaps an infinity) of other agents, in the spirit of Bala and Goyal’s “royal family” [1]. We call such nodes ‘hubs’ for obvious reasons. For instance, a popular blogger or a newspaper might constitute such a hub. Here too the same computational guarantees hold.

4.3.2 Random graphs

Consider a random graph on $n$ nodes drawn from the configuration model with a given degree distribution.$^1$ It is well known that such graphs are locally tree-like with high probability(see, e.g. [6]). More formally, for any $t < \infty$ we have

$$\lim_{n \to \infty} P [B^t_i \text{ is a tree.}] = 1. \quad (9)$$

Since node calculations up to time $t$ depend only on $B^t_i$, it follows that with high probability (w.h.p.), for an arbitrarily selected node, the tree calculations suffice for any constant number of iterations.$^2$ As we show in Section 3 just $O(\log \log 1/\epsilon)$ iterations (a small number independent of $n$) are enough to learn the true state of the world with probability at least $1 - \epsilon$ for any $\epsilon > 0$, provided private signals are not too noisy. Thus, our computational approach works for random graphs w.h.p.

4.3.3 Learning without Knowledge of the Graph

Here we consider the situation where nodes do not know the actual graph $G$, but know some distribution over possibilities for $G$. This is potentially a more realistic model. Moreover, the assumption that agents are assumed to know the entire graph structure is considered a weakness of the model of Gale and Kariv. We address this issue here, showing that our algorithm can be modified to allow Bayesian estimation in this case as well.

Let $G \equiv G_n$ be a random graph of $n$ nodes constructed according to the configuration model for a given (node perspective degree) distribution. Denote the degree distribution by $\rho_v$, so that $\rho_v(d) \equiv$ probability that a randomly selected node has degree $d$.

Now, in this ensemble, the local neighborhood up to distance $D$ of an arbitrary node $v$ with fixed degree $d_v$ converges in distribution as $n \to \infty$ to the following: Each of the neighbors of node $v$ has a degree drawn independently according to the ‘edge perspective’ degree distribution $\rho_E$, defined by:

$$\rho_E(d) = \frac{d \rho_v(d)}{\sum_{d' \in \mathbb{N}} d' \rho_v(d')}$$

Further, each of the neighbors of the neighbors (except $v$ itself) again have a degree drawn independently according to $\rho_E(d)$, and so on up to depth $D$. Call the resulting distribution over trees $T^D_{d_v}$.

Now suppose that agents are, in fact, connected in a graph drawn from the ensemble $G_n$ with degree distribution $\rho_v$. Suppose that each node $u$ knows the distribution $\rho_v$ and its own degree $d_u$, but does not know anything else about $G_n$.$^3$ Further, suppose that this is common knowledge. Now in the limit $n \to \infty$, an exact Bayesian calculation for a node $v$ up to time $t$ depends on $\rho_v$ via $T^t_{d_v}$. Since nodes know only their own degree, there are only $\Delta$ different ‘types’ of nodes, where $\Delta$ is the size of the support of $\rho_E(d)$. There is one type for each degree. This actually makes computations slightly simpler than in an arbitrary known graph.

Fix state $s$. Take an arbitrary node $i$. Make it a ‘zombie’ following the vote trajectory $\tau_i$. Now fix some $\delta i$ (ensure $\rho_v(|\delta i|) > 0$). Choose arbitrary $j \in \delta i$. Define $Q [\sigma_j^t = \omega_j^t | \tau_i^t, s]$ as the probability

1 In the configuration model, one first assigns a degree to each node, draws the appropriate number of ‘half-edges’ and then chooses a uniformly random pairing between them. One can further specify that a graph constructed thus is ‘rejected’ if it contains double edges or self-loops; this does not change any of the basic properties, e.g., the local description, of the ensemble.

2 In fact, as mentioned earlier, nodes with a small number of loops in the vicinity can also do their calculations without trouble.

3 Other ‘knowledge’ assumptions can be similarly handled, for instance where a node knows its own degree, the degree of its neighbors and $\rho_V$.
of seeing trajectory $\sigma_j^t = \omega_j^t$ at node $j$ in this setting. This probability is over the graph realization (given $\partial i$) and over the private signals. Note here that $Q [\sigma_j^t = \omega_j^t | \tau_i^t, s] \text{ is the same for any } i, \partial i \text{ and } j \in \partial i$.

Eqs. (2), (3) and (4) continue to hold w.h.p. for the same reasons as before.

The dynamic cavity recursion, earlier given by Eq. (5), becomes

$$Q [\sigma_j^t | \tau_i^t, s] = \sum_{d \in \mathbb{N}} \rho_E(d) \sum_{\sigma_i^{t-1} \ldots \sigma_{t-1}^{t-1}} \sum_{x_i} \mathbb{P} [x_j | s] \cdot$$

$$\cdot 1 \left[ \sigma_j^t = g_j^t \left( x_j, (\tau_i^{t-1}, \sigma_{t-1}^{t-1}) \right) \right] \cdot$$

$$\cdot \prod_{l=1}^{d-1} Q [\sigma_i^{t-1} | \sigma_j^{t-1}, s]. \tag{10}$$

We have written the recursion assuming the neighbors of $j$ are named according to $\partial j \setminus i = \{1, 2, \ldots, d-1\}$. Again, this holds w.h.p. with respect to $n$.

We comment that there is a straightforward generalization to the case of a multi-type configuration model with a finite number of types. Nodes may or may not be aware of the type of each of their neighbors (both cases can be handled). For instance, here is a simple example with two types: There are ‘red’ agents and ‘blue’ agents, and each ‘red’ agent is connected to 3 ‘blue’ agents, whereas each ‘blue’ agent is connected to either 5 or 6 ‘red’ agents with equal likelihood. In this case the degree distribution itself ensures that nodes know the type of their neighbors as being the opposite of their own type. Multi-type configuration models are of interest since they allow for a rich variety ‘social connection’ patterns.

### 4.3.4 Observing random subsets of neighbors

We may not interact with each of our friends every day. Suppose that for each edge $e$, there is a probability $p_e$ that the edge will be ‘active’ in any particular iteration, independent of everything else. Let $ae(t) \in \{*, a\}$, be an indicator variable for whether edge $e$ was active at time $t$ (a denotes ‘active’). Now, the observation by node $i$ of node $j$ belongs to an extended set that includes an additional symbol * corresponding to the edge being inactive. Thus, there are $(|S| + 1)^d + 1$ possible observed trajectories up to time $t$. Our algorithm can be easily adapted for this case. The modified ‘zombie’ process involves fixing state of the world $s$, trajectory $\tau_i$ and also $(a_{ij}(t))_{j \in \partial i}$ for all times $t$. The form of posterior on the state of the world, Eq. (3), remains unchanged. The cavity recursion Eq. (5) now includes a summation over the possibilities for $(a_1^{t-1}, \ldots, a_d^{t-1})$. The overall complexity remains $\mathcal{O}(d)$.

The case where node $v$ becomes inactive with some probability $p_v$ in an iteration, independent of everything else, can also be handled similarly. A suitable formulation can also be obtained when both the above situations are combined, so that both nodes and edges may be inactive in an iteration.

### 5 Rapid learning on trees

We say that there is **doubly exponential convergence** to the state of the world $s$ if the error probability $P[\sigma_i(t) \neq s]$ decays with round number $t$ as

$$- \log(P[\sigma_i(t) \neq s]) \in \Omega(b^t) \tag{11}$$

where $b > 1$ is some constant.

The following is an immediate corollary of Theorem 4.5

**Corollary 5.1.** Consider iterative Bayesian learning on a tree of with maximum degree $d$. If we have doubly exponential convergence to $s$, then computational effort that is polylogarithmic in $(1/e)$ suffices to achieve error probability $P[\sigma_i(t) \neq s] \leq e$.

We are handicapped by the fact that very little is known rigorously about convergence of iterative Bayesian learning. Nevertheless we provide the following evidence for doubly exponential convergence on trees:

In Section 5.1 we study a simple case with two possible states of the world and two possible private signal values on a regular directed tree. We show that except for the case of very noisy signals, we have doubly exponential convergence if the degree is at least five.

Next, in Section 5.2 we state a conjecture and show that it implies doubly exponential convergence of iterative Bayesian learning also on undirected trees. We provide numerical evidence in support of our conjecture.

#### 5.1 Directed trees

Consider an infinite directed $d$-ary tree. By this we mean a tree graph where each node $i$ has one ‘parent’ who observes $i$ and $d$ ‘children’ whom $i$ observes, but who do not observe $i$. Learning in such a tree is much easier to analyze (than an undirected tree) because the trajectories of the $d$ children are uncorrelated, given $s$.

We assume a binary state of the world $s$ and independent binary signals that are each incorrect with probability $\delta$.  

---

4We need the ball of radius $t$ around $i$ to be a tree.
Lemma 5.2. On an infinite directed $d$-ary tree, the error probability (at any node) at time $t$ is bounded above by $\delta_t$, where $\delta_0 \equiv \delta$ and we have a recursive definition

$$\delta_t \equiv \mathbb{P}[\text{Binomial}(d, \delta_{t-1}) \geq d/2]. \quad (12)$$

Proof. We proceed by induction on time $t$. Clearly, the error probability is bounded above by $\delta_0$ at $t = 0$. Suppose, $\mathbb{P}[\sigma_i(t) \neq s] \leq \delta_t$ by the induction hypothesis. Let the children of $j$ be $1, 2, \ldots, d$. Define $\bar{\sigma}_j$, the opinion of the majority of the children, by

$$\bar{\sigma}_j(t + 1) = \text{sgn} \left( \sum_{l=1}^{d} \sigma_l(t) \right),$$

where $\text{sgn}(0)$ is arbitrarily assigned the value $-1$ or $+1$. The ‘error-or-not’ variables $\{\sigma_l(t) \neq s\}$ are iid, with $\mathbb{P}[\sigma_i(t) \neq s] \leq \delta_t$. Consider a node $j$ making a decision at time $t + 1$. Using Eq. (13), it follows that $\mathbb{P}[\bar{\sigma}_j(t + 1) \neq s] \leq \delta_{t+1}$. Induction completes the proof.

It follows (by an argument similar to the one used in the proof of theorem 5.4 below) that we have doubly exponential convergence to the true state of the world, if the noise level is not too high. We obtain

$$-\log \mathbb{P}[\sigma_i(t) \neq s] \in \Omega \left( (d/2)^t \right). \quad (14)$$

implying that $O(\log \log((1/\epsilon)))$ rounds suffice to reduce the error probability to below $\epsilon$.

5.2 Bayesian vs. ‘majority’ updates

We conjecture that iterative Bayesian learning leads to lower error probabilities (in the weak sense) than a very simple alternative update rule we call “majority dynamics”. Under this rule the agents adopt the action taken by the majority of their neighbors in the previous iteration (this is made precise in Definition A.1). Our conjecture is natural since the iterative Bayesian update rule chooses the vote in each round that (myopically) minimizes the error probability.

Conjecture 5.3. On any regular tree with independent identically distributed private signals, the error probability under iterative Bayesian learning is no larger than the error probability under majority dynamics (cf. Definition A.1) after the same number of iterations.

We use $\bar{\sigma}_i(t)$ to denote votes under the majority dynamics.

In Appendix A we show doubly exponential convergence for majority dynamics on regular trees:

**Theorem 5.4.** Assume binary $s$ with uniform prior. Agents’ initial votes $\bar{\sigma}_i(0)$ are correct with probability $1 - \delta$, and independent conditioned on $s$. Let $i$ be any node in an (undirected) $d$ regular tree for $d \geq 5$. Then, under the majority dynamics,

$$-\log \mathbb{P}[\bar{\sigma}_i(t) \neq s] \in \Omega \left( \left( \frac{1}{2} (d - 2) \right)^t \right). \quad (15)$$

Thus, if Conjecture 5.3 holds:

- We also have doubly exponential convergence for iterative Bayesian learning on regular trees with $d \geq 5$, implying that for any $\epsilon > 0$, an error probability $\epsilon$ can be achieved in $O(\log \log(1/\epsilon))$ iterations under iterative Bayesian learning.

- Combining with Theorem 4.5 (cf. Corollary 5.1), we see that the computational effort that is polylogarithmic in $(1/\epsilon)$ suffices to achieve error probability $1/\epsilon$.

This compares favorably with the quasi-poly$(1/\epsilon)$ bound on computational effort that we can derive by combining Conjecture 5.3 and the simple dynamic program described in Section 3.

In Table 1 we provide numerical evidence on regular undirected trees in support of our conjecture. Further numerical results are presented in Appendix B. All computations are exact, and were performed using the dynamic cavity equations. The results are all consistent with our conjecture over different values of $d$ and $\mathbb{P}[x_i \neq s]$.

<table>
<thead>
<tr>
<th>Round</th>
<th>Bayesian</th>
<th>Majority</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>$0.15$</td>
<td>0.15</td>
</tr>
<tr>
<td>1</td>
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<td>$2.7 \cdot 10^{-2}$</td>
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<td>2</td>
<td>$7.6 \cdot 10^{-4}$</td>
<td>$1.7 \cdot 10^{-3}$</td>
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<tr>
<td>3</td>
<td>$2.8 \cdot 10^{-7}$</td>
<td>$8.4 \cdot 10^{-6}$</td>
</tr>
<tr>
<td>4</td>
<td>$1.4 \cdot 10^{-12}$</td>
<td>$2.5 \cdot 10^{-10}$</td>
</tr>
</tbody>
</table>

Table 1: Error probability on regular tree with $d = 5$ and $\mathbb{P}[x_i \neq s] = 0.15$, for (i) Bayesian and (ii) majority updates. The agents break ties by picking their original private signals.
We would like to thank Andrea Montanari, Elchanan Mossel and Allan Sly for valuable discussions.

References


A Majority dynamics: Proof of Theorem 5.4

In this section we study a very simple update rule, ‘majority dynamics’. We use \( \hat{\sigma}_i(t) \) to denote votes under the majority dynamics.

**Definition A.1.** Under the majority dynamics, each node \( i \in V \) chooses his vote in round \( t+1 \) according to the majority of the votes of his neighbors in round \( t \), i.e.

\[
\hat{\sigma}_i(t+1) = \text{sign} \left( \sum_{j \in \partial i} \hat{\sigma}_j(t) \right)
\]
Ties are broken by flipping an unbiased coin.

As before, \( s \in \{-1, +1\} \) is drawn from a 50-50 prior and nodes receive ‘private signals’ \( \hat{\sigma}_i(0) \) that are correct with probability \( 1 - \delta \), and independent conditioned on \( s \).

Consider an undirected \( d \) regular tree. The analysis in this case is complicated (relative to the case of a directed tree) by dependencies which have to be carefully handled.

**Lemma A.2.** Let \( i \) and \( j \) be adjacent nodes in the tree. Then for all \( (\hat{\sigma}_i^t, \hat{\sigma}_j^t) \in \{-1, +1\}^{2t} \)

\[
\Pr \left[ \hat{\sigma}_i(t+1) = -1 \mid \hat{\sigma}_i^t, \hat{\sigma}_j^t, s = +1 \right] \leq \delta_t
\]

where \( \delta_t \) is defined recursively by \( \delta_0 \equiv \delta \), and

\[
\delta_t \equiv \Pr \left[ \text{Binomial}(d-1, \delta_{t-1}) \geq d/2 - 1 \right]
\]

**Proof.** We proceed by induction. Clearly Eq. (16) holds for \( t = 0 \). Suppose Eq. (16) holds for some \( t \). We want to show

\[
\Pr \left[ \hat{\sigma}_i(t+1) = -1 \mid \hat{\sigma}_i^t, \hat{\sigma}_j^t, \hat{\sigma}_i^t, ... \hat{\sigma}_{d-1}^t, s = +1 \right] \leq \delta_{t+1}
\]

for all possible \( \xi \equiv (\hat{\sigma}_i^t, \hat{\sigma}_j^t, \hat{\sigma}_i^t, ... \hat{\sigma}_{d-1}^t) \).

We reason as follows. Fix the state of the world \( s \) and the trajectories \( \hat{\sigma}_i^t \) and \( \hat{\sigma}_j^t \). Now this induces correlations between the trajectories of the neighbors \( l_1, l_2, ... l_{d-1} \), caused by the requirement of consistency, but only up to time \( t - 1 \). If we further fix \( \hat{\sigma}_{l_m}^t \), then \( \hat{\sigma}_{l_m}^t \) (and \( \hat{\sigma}_{l_m} \) at all future times) is conditionally independent of \( \hat{\sigma}_{l_m}^t \) for \( m \neq m \). Thus, we have

\[
\Pr \left[ \hat{\sigma}_{l_m}^t = \pm 1 \mid \xi, s = +1 \right] = \Pr \left[ \hat{\sigma}_{l_m}^t = \pm 1 \right] \]

and therefore, using the induction hypothesis

\[
\Pr \left[ \hat{\sigma}_{l_m}^t = -1 \mid \xi, s = +1 \right] \leq \delta_t
\]

for all \( m \in \{1, 2, ... d - 1\} \). Also, the actions \( \hat{\sigma}_{l_1}(t), ... \hat{\sigma}_{l_{d-1}}(t) \) are conditionally independent of each other given \( \xi, s = +1 \). We have

\[
\hat{\sigma}_i(t+1) = \text{sgn}(\hat{\sigma}_j(t) + \hat{\sigma}_{l_1}(t) + ... + \hat{\sigma}_{l_{d-1}}(t)),
\]

with \( \text{sgn}(0) \) being assigned value \(-1\) or \(+1\) with equal probability. We have

\[
\Pr \left[ \hat{\sigma}_i(t+1) = -1 \mid \xi, s = +1 \right] \leq \Pr \left[ \text{Binomial}(d-1, \delta_t) \geq d/2 - 1 \right]
\]

from Eq. (20) and conditional independence of \( \hat{\sigma}_i(t), ... \hat{\sigma}_{l_{d-1}}(t) \). This yields Eq. (19).

**Proof of Theorem 5.4.** By applying the multiplicative version of the Chernoff bound\(^6\) to Eq. (17) we have that

\[
\delta_{t+1} \leq e^{(d-2)/2 - (d-1)\delta_t(2\delta_t(d-1)/(d-2))^{(d-2)/2}}
\]

Dropping the term \( e^{-1/d}\delta_t \), we obtain

\[
\delta_{t+1} \leq (2e\delta_t(d-1)/(d-2))^{1/2(d-2)}.
\]

This is a first order non-homogeneous linear recursion in \( \log \delta_t \). If it was an equality it would yield

\[
\log \delta_t = \left( \log \delta + \frac{d - 2}{d - 4} \log \left( 2e(d-1)/(d-2) \right) \right) \left( \frac{1}{2} \right)^t (d-2) / 2
\]

and so

\[
- \log \delta_t \leq \Omega \left( \left( \frac{1}{2} \right) (d-2) \right)^t,
\]

as long as

\[
\log \delta < \frac{d - 2}{d - 4} \log \left( 2e(d-1)/(d-2) \right).
\]

Theorem 5.4 is non-trivial for \( d \geq 5 \). The upper limit of the noise for which it establishes rapid convergence approaches \( (2e)^{-1} \) as \( d \) grows large (see also the discussion below for large \( d \)).

### A.1 Convergence for large \( d \)

We present now a short informal discussion on the limit \( d \to \infty \). We can, in fact, use Lemma 5.2 to show convergence is doubly exponential for \( \delta = 1/2 - c/d \) for some \( c < \infty \) that does not depend on \( d \).

Here is a sketch of the argument. Suppose \( \delta = 1/2 - c_1/d \). Then, for all \( d > d_1 \) where \( d_1 < \infty \), there exists \( c_2 < \infty \) such that

\[
\Pr[X \geq (1 + \eta)E[X]] \leq \left( \frac{\text{sgn}(\delta t \alpha)}{\beta \delta t \alpha + \eta} \right)^{E[X]}. \quad \text{We substitute } E[X] = \delta_t(d-1) + 1 + \eta = (d/2 - 1)/\delta_t(d-1).
\]
\(\mathbb{P}\{\text{Binomial}(d - 1, \delta) \geq d/2 - 1\} < 1/2 - c_2/\sqrt{d}\). This can be seen, for instance, by coupling with the Binomial\((d - 1, 1/2)\) process and using an appropriate local central limit theorem (e.g., see [10, Theorem 4.4]). Thus, \(\delta_1 < 1/2 - c_2/\sqrt{d}\). Further, \(c_2\) can be made arbitrarily large by choosing large enough \(c_1\). Next, with a simple application of the Azuma’s inequality, we arrive at \(\delta_2 < c_3\) (where \(c_3 \to 0\) as \(c_2 \to \infty\)). Now, for small enough \(c_3\), we use the Chernoff bound analysis in the proof of Theorem 5.4 and obtain doubly exponential convergence.

### B Further numerical results

Table 2, together with table 1 above, contrast the error probabilities of Bayesian updates with those of majority updates. All cases exhibit lower error probabilities (in the weak sense) for the Bayesian update, consistent with Conjecture 5.3. Table 3 contains the data plotted in Figure 1. Also for these parameters, we found that the Bayesian updates showed lower error probabilities than the majority updates (though we omit to present the majority results here).

The running time to generate these tables, on a standard desktop machine was less than a minute. We did not proceed with more rounds because of numerical instability issues which begin to appear as error probabilities decrease.

<table>
<thead>
<tr>
<th>Round</th>
<th>Bayesian</th>
<th>Majority</th>
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</thead>
<tbody>
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<td>0.15</td>
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<tr>
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</tr>
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</table>

Table 2: \(d = 3, \mathbb{P}\{x_i \neq s\} = 0.15\)

<table>
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<th>(d = 5)</th>
<th>(d = 7)</th>
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<td>7</td>
<td>1.5 \cdot 10^{-3}</td>
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</table>

Table 3: Error probabilities with \(\mathbb{P}\{x_i \neq s\} = 0.3\), for regular tree of different degrees \(d\). This data is displayed in Figure 1.