Detection of the dominant direction of information flow in densely interconnected regulatory networks

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ABSTRACT

Motivation: Finding the dominant direction of flow of information in densely interconnected regulatory or signaling networks is required in many applications in computational biology and neuroscience. This is achieved by first identifying and removing links which close feedback loops in the original network and hierarchically arranging nodes in the remaining network. In mathematical language this corresponds to a problem of making a graph acyclic by removing as few links as possible and thus altering the original graph in the least possible way. Practically all applications the exact solution of this problem requires an enumeration of all combinations of removed links, which is computationally intractable.

Results: We introduce and compare two algorithms: the deterministic, ‘greedy’ algorithm that preferentially cuts the links that participate in the largest number of feedback cycles, and the probabilistic one based on a simulated annealing of a hierarchical layout of the network which minimizes the number of “backward” links going from lower to higher hierarchical levels. We find that the annealing algorithm outperforms the deterministic one in terms of speed, memory requirement, and the actual number of removed links. Implications for system biology and directions for further research are discussed.

Availability: Source codes of F90 and Matlab implementation of these two algorithms are available from the authors upon request.

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1 INTRODUCTION

During the last several years, a substantial amount of information on large-scale structure of intracellular regulatory networks has been accumulated. However, the growth in our understanding of how these networks manage to function in a robust and specific manner was lagging behind the sheer rate of data acquisition. The fact these networks are frequently visualized as a giant “hairball” (Fig. 1) consisting of a multitude of edges, linking most constituent protein-nodes to each other serves as a striking illustration of the complexity of the issue at hand.

To understand the functioning or even to efficiently visualize a densely interconnected directed network it is desirable to determine the dominant direction of information flow and to identify links that go against this flow and thus close feedback loops. Ordering a network with respect to the dominant direction of information flow can help to determine its previously unknown inputs and outputs, to track back hidden sources of perturbations based on their observable downstream effects, etc. A simple-minded hierarchical layout of a densely interconnected network is often impossible due to a ubiquitous presence of feedback loops. Indeed, all nodes in a strongly connected component of a network by definition are simultaneously upstream and downstream of each other. However, if most feedback loops are closed by relatively few feedback signaling links, the dominant direction of information flow could still be reconstructed based on a network topology alone. An identification and removal of these relatively infrequent feedback links would enable one to perform a hierarchical layout of the remaining acyclic network which still sufficiently resembles the original one.

In this work we consider the problem of identifying the minimum set of links, removal of which would render a graph acyclic. In the next section we introduce two rather different algorithms allowing one to approximately accomplish this goal, a deterministic ‘greedy’ algorithm and a probabilistic Metropolis annealing, and compare their performance. We find that the probabilistic algorithm outperforms the deterministic one in better minimizing the number of removed links, and memory requirements, while maximizing the speed. A simple visual example is provided for the situation when the deterministic algorithm is non-optimal. Following that, we discuss biological implications and applications of our findings as well as how additional constraints such as a priori knowledge of the function and therefore hierarchical position of certain nodes may affect the algorithm performance.

2 APPROACH

Consider a graph of $N$ vertices labeled as 1, 2, 3, . . . , $N$ and $L$ directed links labeled by pairs of vertices they connect, $l_i \equiv (n_i, m_i)$. The goal is to remove as few as possible of the links to make the graph acyclic, or feedback-free.

An exact way to solve this problem is to sample all possible combinations of links to be removed, starting with enumerating individual links, then pairs of links, etc, until the first acyclic graph is obtained. Evidently, if a removal of $l$ links finally yields an acyclic graph, such sampling would require checking the $\sum_{i=1}^{L} \binom{l}{i}$ networks for cycles. For the biologically relevant values of $L \sim 10^3 - 10^4$ and $l \sim 10 - 10^2$ this approach is clearly unfeasible.  

\footnote{From an obvious identity, $\sum_{i=1}^{L} \binom{l}{i} = 2^L - 1$, it follows that even for fairly modest $L = 10^2$ and $l = L/2$ the number of such attempts is $\sim 10^{15}$.}
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Fig. 1. Caption, A part of the post-translational regulatory network in human shown here includes 1671 automatically and manually curated protein modification interactions (phosphorylation, proteolytic cleavage, etc.) between 732 proteins from our ResNet database Nikitin et al., 2003. Panel A contains the “hairball” visualization of the network structure emphasizing interconnections between individual pathways. Red edges lie within the strongly connected component of this network consisting of 107 proteins that could all be linked to each other by a path in both directions. This makes any two of these proteins to be simultaneously upstream and downstream from each other. In Panel B we optimally distribute proteins over a number of hierarchical levels. Red arrows represent 208 putative feedback links going from lower levels of the hierarchy to higher ones, while yellow ones – 512 feed-forward links jumping over one or more hierarchical levels. Only proteins and links reachable from one of the 71 receptors placed at the top hierarchical level were included.

Fig. 2. Caption, Removal of a single (3, 1) link makes this 3-vertex graph acyclic.

2.1 Greedy algorithm

A natural reduction of such exact enumeration approach is a “greedy” algorithm which performs the “steepest descent” in the number of cycles. We implemented the following realization of such link removal algorithm:

- By enumerating all cycles in a graph, each link is assigned a score equal to the number of cycles it is a member of.
- The link with the highest score is removed. When more than one link have the same highest score, a link to be removed is selected among the highest-scored ones by random.
- This procedure of cycle enumeration and link removal is repeated until no cycles are found.

The cycle enumeration can be implemented by following all paths that originate from a given vertex and recording only the cycles that come back to this vertex. The procedure is repeated for each of the N graph vertices; evidently, each cycle of length C is counted C times and a proper normalization is performed.

An example of network where the greedy algorithm performs flawlessly is shown in Figure 2. Here the link (3, 1) carries the maximum score 2. A removal of this link indeed makes the graph acyclic, while a removal of any other than (3, 1) link would require a subsequent removal of the second link to achieve the same goal. However, one would suspect that as any “steepest descent” method, the proposed greedy algorithm, performing a sometimes near-sighted local one-step optimization, may miss the globally optimal solution. This is indeed often the case for bigger and more complex graphs; a fairly simple example is given in Fig. 3.

2.2 Simulated annealing network ordering

The task of finding the minimum number of links, cutting which makes the graph acyclic, can be interpreted as an optimization problem and tackled by probabilistic methods such as simulated annealing. Evidently, there exist more than one way to define the optimization function, and after exploring several possibilities we converged to the following one:

- For a given network, a set of M levels is introduced ($M \leq N$), in reality, $M \ll N$ and is of the order of the graph diameter. Initially, all nodes are distributed on the levels randomly.
- For a particular distribution of nodes on levels, the number of links that go opposite to the hierarchy, that is, from a lower level to the same or a higher one, is declared to be the energy $E$ of the distribution, or the optimization function.
In the previous section we introduced two algorithms intended to make a network acyclic by removing the least number of links. The stochastic stimulated annealing level-ordering algorithm outperforms the deterministic greedy algorithm in all respects. Indeed, the greedy algorithm requires tracking along all paths originating from a given vertex, which uses a lot of memory and slows the performance significantly. We found it impractical to apply the greedy algorithm to networks with more than 100–200 vertices. This rules out its use for all-organism network ordering and limits its utility in the ability to determine how many cycles pass through a given link. Otherwise, $M$ could be determined self-consistently, by observing when the number of counter-hierarchical links stops to decrease upon the increase in the number of levels. This is illustrated in Fig. 5 where a plot of the number of non-hierarchical links vs number of levels is presented for the human protein phosphorylation network.

3 DISCUSSION

A useful property of this algorithm is that in addition to making a network acyclic, it also produces a hierarchical layout. The number of levels $M$ could be fixed by the requirements for such layout. Otherwise, $M$ could be determined self-consistently, by observing when the number of counter-hierarchical links stops to decrease upon the increase in the number of levels. This is illustrated in Fig. 5 where a plot of the number of non-hierarchical links vs number of levels is presented for the human protein phosphorylation network.
Often there exist some *a priori* knowledge on the hierarchical positions of certain network nodes. For example, many of the receptor proteins localized in the membrane upon activation pass the signals downstream signaling cascades made of proteins localized in the cytoplasm and ultimately in the cell’s nucleus. Thus receptor proteins might have to be forcefully put on the upper levels of the hierarchical layout of such signaling network. Contrary to receptors, many transcription factors serve the role of effectors of signaling pathways and thus must occupy the lowest levels of the hierarchy. Initial, or possibly permanent, position of such nodes on the hierarchical levels often helps to converge to the better in terms of fewer feedback links, or more biologically relevant solution.

In a similar way, the orientation of certain links (or equivalently, pairs of nodes) could be quenched if they are known to be of the feed forward of feed back nature. Based on the initial knowledge of network functioning, it is also possible to assign a certain weight to a link, so that the energy $E$ of a particular assignment of nodes to layers is a sum of weights of the counter-hierarchical links. Thus the *a priori* known plausibility of a link to be (or not to be) a feedback can be introduced into the layering algorithm.

It is also possible to improve the visual perception of the layout by shortening the hierarchical links. In its present edition, a “good” or hierarchical link may be arbitrary long, i.e. go down many levels, without carrying any energetic penalty. This interferes with identifying the hierarchical levels as certain stages of network flow. Introduction of a small energetic penalty for particularly long links may alleviate this shortcoming.

We leave these questions as well as those of particular application of ordering algorithms to catalytic signaling and transcription regulation cellular networks for future studies and publications.

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