BROADER PERCEPTION FOR LOCAL COMMUNITY IDENTIFICATION

F. T. W. (Frank) Koopmans and Th. P. (Theo) van der Weide
Institute for Computing and Information Sciences, Radboud University, Nijmegen, The Netherlands

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Abstract: A local community identification algorithm can identify the network community of a given start node without knowledge of the entire network. Such algorithms only consider nodes within or directly adjacent to the local community. Therefore a local algorithm is more effective than an algorithm that partitions the entire network when only a small portion of a large network is of interest or when it is difficult to obtain information about the network (such as the world wide web). However, local algorithms cannot deliver the same quality as their global counterparts that use the entire network. We propose an improvement to local community identification algorithms that will decrease the gap between relevant network knowledge of global and local methods. Benchmarks on synthetic networks show our approach increases the quality of locally identified communities in general and a decrease of the dependency on specific source nodes.

1 INTRODUCTION

The science of complex systems is a popular interdisciplinary field of research. Using a network to represent the elements and interactions in a complex network is a commonly used tool to study complex system phenomena (Strogatz, 2001; Albert and Barabasi, 2002). Such analysis is done (among others) on protein networks, social networks and (parts of) the internet (Newman, 2003; Palla et al., 2007).

Complex systems, and thus their representation as a network, often exhibit an a-priori unknown structure of building blocks with varying functions. These blocks are represented in the network as sets of nodes that are densely connected and have relatively few connections to the rest of the network. Such building blocks are referred to as modules, communities or clusters. Identifying and observing such communities may lead to a greater understanding of the structure and functioning of a complex system. Many different techniques for identifying all communities within a network, with varying computational costs and accuracy, have been developed (Danon et al., 2005; Lancichinetti and Fortunato, 2009).

In some cases one would rather identify a specific community instead of all communities in the entire network in order to reduce the amount of information that needs to be processed or acquired. For example, finding the community around a given website on the internet or finding all friends of a specific person in a social network. In such cases there should be no need to process the entire network. A local community is a community in the network identified by starting from a start node and using only information from the context of the local community to identify local community members. Thus no knowledge of the entire network is required.

There are many different approaches to local community identification (Clauset, 2005; Schaeffer, 2005; Chen et al., 2009). But local approaches have a lacking quality compared to approaches that cluster entire networks and they are overly dependant on a well chosen start node (Bagrow, 2008).

In this paper we focus on improving local community identification algorithms in order to increase their overall quality and reduce the dependency on the start node. First we discuss local community structure, then we propose an improvement to local algorithms and finally we present benchmark results and discussion thereof.
2 LOCAL COMMUNITY IDENTIFICATION

A local community is considered from the perspective of a start node from which that community can be generated. As a consequence, each node in the local community has a special relation with this start node and the local community as a whole is a regular community in the network. Local communities are a soft clustering, where each node can be a member of multiple communities, of a part of a network thus overlap may occur. The ideal local community for some start node $v$ would be the community with the highest quality that contains $v$. Let the quality of a community $C \subseteq N$ be quantified by community quality measure $\mu(C)$.

A community is denoted as the set of nodes it consists of. We define the universe of a community $C$, denoted as $U(C)$, as all nodes that are outside of $C$ and adjacent to any node within $C$.

\[ U(C) = \{ v \in N - C \mid \exists_{x \in C} [u \longrightarrow x] \} \quad (1) \]

The boundary of $C$, denoted as $B(C)$, consists of the nodes within the local community that have a relation with nodes outside that community.

\[ B(C) = \{ u \in C \mid \exists_{v \in U(C)} [u \longrightarrow v] \} \quad (2) \]

We introduce node relation $\varphi(C, u, v)$ to indicate that in community $C$ node $u$ is related to node $v$. This relation is expressed by a path from $u$ to $v$ within that community:

\[ \varphi(C, u, v) \equiv u \longrightarrow v \lor \exists_{x \in C} [u \longrightarrow x \land \varphi(C, x, v)] \quad (3) \]

where $u \longrightarrow v$ denotes that there is an edge from $u$ to $v$.

A local community $C$ from the perspective of start node $v$ is a proper community, denoted as $\Phi(C, v)$, if all its nodes are related to the start node:

\[ \Phi(C, v) = \{ v \in C \land \forall_{x \in C - \{v\}} [\varphi(C, v, x)] \} \quad (4) \]

In that case, we refer to $v$ as a proper generator of $C$. A local community $C$ may have different generators. Furthermore, a node may be the generator of different communities.

We will assume that the communities that can be properly generated from the same start node can be ordered hierarchically according to the subset relation:

\[ \Phi(C_1, v) \land \Phi(C_2, v) \implies (5) \exists_{C \in [C_1, C_2]} [\Phi(C, v)] \]

We will call $C$ a root community from node $v$, denoted as $\Phi^+(C, v)$, if $C$ is a minimal community that can be properly generated from $v$:

\[ \Phi^+(C, v) \equiv \forall_{X} [\Phi(X, v) \land X \subseteq C \implies X = C] \quad (6) \]

**Lemma 1.** Let $\Phi^+(C, v)$, then $C = \bigcap \{ X \mid \Phi(X, v) \}$. Note that $\Phi$ is not related to community quality. The ideal local community for some start node $v$ will be a local community properly generated by $v$ with a maximal $\mu$-score. We introduce $\arg\max_{x \in X} f(x)$ to indicate that $x \in X$ yields a result for $f(x)$ that is not surpassed by any value in $X$:

\[ \arg\max_{x \in X} f(x) = \{ x \in X \mid \forall_{y \in X} [f(y) \leq f(x)] \} \quad (7) \]

Let the set of communities where $v$ is a proper generator, denoted as $\zeta(v)$, be defined as:

\[ \zeta(v) = \{ C \subseteq N \mid \Phi(C, v) \} \quad (8) \]

then the set of ideal communities for start node $v$, denoted as $I(v)$, is defined as:

\[ I(v) = \arg\max_{C \in \zeta(v)} \mu(C) \quad (9) \]

where multiple ideal local communities imply there are different local community compositions for start node $v$ that yield the same $\mu$-score, which is the maximum. The computation of these ideal communities is not feasible in practice because of the time complexity of this measure. While we cannot use this concept in practice it does provide insight in the desired result from a local algorithm. It sets a goal, find the best local community around a given start node.

3 BROADENING THE LOCAL SCOPE

A strong advantage of global community identification algorithms over their local counterparts is knowledge of the entire network. For a local algorithm there is no way of telling what lies beyond the universe of the community it is building. There may be knowledge about general network characteristics, but node specific information is not available. This shortsightedness has a negative impact on the quality of the resulting local community because local algorithms tend to halt at every sign of community quality decline, they cannot see possible improvement beyond. A related and common problem for local algorithms is their dependence on a well chosen start node. Direct neighbors of a start node may not seem like good building blocks for a community. A commonly applied band-aid solution to the start node dependency is to make the local algorithm ignore its stopping criterion until hitting a predetermined size or quality to discourage preemptive stopping. But this method relies on a preset threshold and thus priori knowledge of the expected community size and structure. Also, after this preset threshold has been reached
the algorithm will still fail to see beyond minor barriers.

We suggest the improvement of local community identification algorithms in general by adding more contextual information to their selection criteria. By making a local algorithm look ahead further than one edge from the community we decrease the shortsightedness of this approach and allow for a more informed and balanced judgement on community membership, at the cost of higher computational and situational crawling complexity. This improvement can be applied to any local algorithm.

We propose the following example application where we extend the algorithm its network knowledge by offering nodes one step beyond the universe. Besides the universe node itself, we also investigate the addition of that a node together with any combination of its neighbors. This will provide insight into possible local community structure, and quality, that lies beyond the community universe. The set of possible addition sets for community $C$ at distance 2, denoted as $A(C)$, is defined as follows:

$$A(C) = \bigcup_{u \in U(C)} \{X \cup \{u\} | X \subseteq \varphi(b(u))\}$$  \hspace{1cm} (10)

where $b(u) = \{v | u \rightarrow v\}$ is the set of nodes that can be reached from $u$ in one step. Then we are interested in:

$$\operatorname{argmax}_{C \in A(C)} \mu(C \cup C')$$  \hspace{1cm} (11)

Note that the computational complexity will increase rapidly as the lookahead distance $k$ is increased, especially in dense networks. Further research should consider different distances for the lookahead approach, consider the quality and computational complexity tradeoff and determine an optimum (if any).

4 VALIDATION

Our improvement to local community identification algorithms as proposed in section three is validated by a series of tests on synthetic networks. Since we aim to verify the improvement of local identification algorithms in the area where these often struggle we will generate networks with a low average degree (4, 5 and 8). This will result in networks that contain a lot of potentially troublesome start nodes. We will run four tests on every node in the network, varying the algorithm (regular or improved) and the community measure. We measure community quality by the widely known local modularity (Clauset, 2005) and relative density (Schaeffer, 2005; Lancichinetti et al., 2009) definitions.

For our test we generate an undirected network according to the Barabasi–Albert model (Albert and Barabasi, 2002) and rewire it to create a flat community structure as proposed by Bagrow (Bagrow, 2008). The rewiring is done by creating $k$ sets of nodes (representing the communities) in the network and then rewiring inter-community edges to intra-community edges while preserving the degree distribution. Our benchmark networks contain 128 nodes equally divided amongst 4 communities.

The quality of a community identification algorithm is evaluated by quantifying the similarity between the algorithm output and the synthetic community structure. We will adopt the Jaccard Similarity Coefficient (JSC) which is defined as the commonality of both sets divided by their generality:

$$JSC(X, Y) = \frac{|X \cap Y|}{|X \cup Y|}$$  \hspace{1cm} (12)

4.1 Results

Running these tests on 50 generated networks yields the plot of the JSC score frequency shown in Figure 1. In the networks with an average degree of 4 and 5 we observe a significant increase in high similarity and decrease of outliers for both quality measures when our algorithm improvement is applied. There is a relatively low gain for the more dense networks with average degree 8. The plots also show that the result, while strongly improved, is not perfect yet. We observe a couple of reasons why even the improved algorithm is struggling for some start nodes.

First of all, when the boundaries of two communities are not very sharp and the start node is a boundary node (as defined by the synthetic graph structure) the algorithm may start of in the wrong direction and identify the wrong community. Suppose we start with node $v$ that is a member of community $C$ according to the synthetic structure. If the algorithm identifies $C' \cup \{v\}$ where $C'$ is another community defined by the synthetic structure, then the result of the algorithm may be a quite strong community. But the similarity measure will yield a bad result because there is very little overlap between the reference community and the found community.

Also, the local algorithm may find a strong community that is a subset of the community it is supposed to find. The gap between the found community and agglomerating until the algorithm identifies a larger and stronger community may be too large for the lookahead algorithm to recognize. There lies a tradeoff between the time complexity of the lookahead algorithm and the effectiveness of identifying improvement beyond the universe candidates.
Finally, the use of the synthetic network structure while measuring quality is not ideal. When the goal of a local algorithm is to identify the strongest community around a start node (as is the case in this paper) that synthetic reference community may not be the strongest community thus we may not compare the algorithm output with the ideal community.

5 CONCLUSIONS

Local community identification is particularly useful in large networks where only a small part of the network is of interest. However, the quality of local identification algorithms is lacking when compared to their global counterparts. In this paper we suggest the addition of contextual information beyond the direct neighbors of a local community when evaluating local community candidates. By decreasing the gap between relevant network knowledge of global and local methods we improve the quality of local community identification algorithms in general.

Our experiments on synthetic networks have shown promising results. Extending generic local community identification algorithms such that they consider nodes one step beyond their universe will lead to a significant increase in quality and decrease on start node dependency.

After deriving this approach from theory and showing a proof of concept in this paper we are interested in further research on several aspects. Firstly, as suggested in section three, one should consider several distances $k$ where $k > 1$. Secondly, an analysis of the computational complexity of this approach is required (while also considering varying network densities). This should lead to an improved experiment using both synthetic and real world datasets to determine the appropriate tradeoff between computational complexity and quality.

REFERENCES


