Preserving Privacy in Social Networks Against Neighborhood Attacks

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Abstract-Recently, as more and more social network data has been published in one way or another, preserving privacy in publishing social network data becomes an important concern. With some local knowledge about individuals in a social network, an adversary may attack the privacy of some victims easily. Unfortunately, most of the previous studies on privacy preservation can deal with relational data only, and cannot be applied to social network data. In this paper, we take an initiative towards preserving privacy in social network data. We identify an essential type of privacy attacks: neighborhood attacks. If an adversary has some knowledge about the neighbors of a target victim and the relationship among the neighbors, the victim may be re-identified from a social network even if the victim's identity is preserved using the conventional anonymization techniques. We show that the problem is challenging, and present a practical solution to battle neighborhood attacks. The empirical study indicates that anonymized social networks generated by our method can still be used to answer aggregate network queries with high accuracy.

I. INTRODUCTION

Recently, as more and more social network data has been made publicly available [1], [2], [3], [4], preserving privacy in publishing social network data becomes an important concern. Is it possible that releasing social network data, even with individuals in the network aonymized, still intrudes privacy?

A. Motivation Example

With some local knowledge about individual vertices in a social network, an adversary may attack the privacy of some victims. As a concrete example, consider a synthesized social network of "close-friends" shown in Figure 1(a). Each vertex in the network represents a person. An edge links two persons who are close friends.

Suppose the network is to be published. To preserve privacy, is it sufficient to remove all identities as shown in Figure 1(b)? Unfortunately, if an adversary has some knowledge about the neighbors of an individual, the privacy may still be leaked.

If an adversary knows that Ada has two close friends who know each other, and has another two close friends who do not know each other, i.e., the 1-neighborhood graph of Ada as shown in Figure 1(c), then the vertex representing Ada can be identified uniquely in the network since no other vertices have the same 1-neighborhood graph. Similarly, Bob can be identified in Figure 1(b) if the adversary knows the 1-neighborhood graph of Bob.

Identifying individuals from released social networks intrudes privacy immediately. In this example, by identifying



Fig. 1. Neighborhood attacks in a social network.

Ada and Bob, an adversary can even know from the released social network (Figure 1(b)) that Ada and Bob are close friends, and they share one common close friend. Other private information can be further derived such as how well a victim is connected to the rest of the network and the relative position of the victim to the center of the network.

To protect the privacy satisfactorily, one way is to guarantee that any individual cannot be identified correctly in the anonymized social network with a probability higher than $\frac{1}{k}$, where k is a user-specified parameter carrying the same spirit in the k-anonymity model [5]. By adding a noise edge linking Harry and Irene, the 1-neighborhood graph of every vertex in Figure 1(d) is not unique. An adversary with the knowledge of 1-neighborhood cannot identify any individual from this anonymous graph with a confidence higher than $\frac{1}{2}$.

B. Challenges

Although privacy preservation in data publishing has been studied extensively and several important models such as *k*anonymity [5] and *l*-diversity [6] as well as many efficient algorithms have been proposed, most of the existing studies can deal with relational data only. Those methods cannot be applied to social network data straightforwardly.

As elaborated in Section I-A, privacy may be leaked if a social network is released improperly to public. In practice, we need a systematic method to anonymize social network data before it is released. However, anonymizing social network data is much more challenging than anonymizing relational data on which most of the previous work focuses.

First, it is much more challenging to model the background knowledge of adversaries and attacks about social network data than that about relational data. On relational data, it is often assumed that a set of attributes serving a quasi-identifier is used to associate data from multiple tables, and attacks mainly come from identifying individuals from the quasi-identifier. However, in a social network many pieces of information can be used to identify individuals, such as labels of vertices and edges, neighborhood graphs, induced subgraphs, and their combinations. It is much more complicated and much more difficult than the relational case.

Second, it is much more challenging to measure the information loss in anonymizing social network data than that in anonymizing relational data. Typically, the information loss in an anonymized table can be measured using the sum of information loss in individual tuples. Given one tuple in the original table and the corresponding anonymized tuple in the released table, we can calculate the distance between the two tuples to measure the information loss at the tuple level. However, a social network consists of a set of vertices and a set of edges. It is hard to compare two social networks by comparing the vertices and edges individually. Two social networks having the same number of vertices and the same number of edges may have very different network-wise properties such as connectivity, betweenness, and diameter. Thus, there can be many different ways to define the measures of information loss and anonymization quality.

Last but not least, it is much more challenging to devise anonymization methods for social network data than for relational data. Divide-and-conquer methods are extensively applied to anonymization of relational data due to the fact that tuples in a relational table are separable in anonymization. In other words, anonymizing a group of tuples does not affect other tuples in the table. However, anonymizing a social network is much more difficult since changing labels of vertices and edges may affect the neighborhoods of other vertices, and removing or adding vertices and edges may affect other vertices and edges as well as the properties of the network.

C. Contributions and Organization

Privacy preservation in publishing social networks is a new challenging problem that cannot be solved by one shot. In this paper, we take the initiative to tackle the problem. We identify an essential type of privacy attacks in social networks: neighborhood attacks, which are illustrated in Section I-A. We model the attacks and the anonymization problem systematically, show its difficulty, and develop a practical solution. We conduct an empirical study which indicates that anonymized social networks generated by our method can still be used to answer aggregate network queries with satisfactory accuracy.

The rest of the paper is organized as follows. We model neighborhood attacks and review related work in Section II. A practical solution is developed in Section III, and examined empirically using both a real data set and a series of synthetic data sets in Section IV. In Section V, we discuss some other related privacy attacks in social networks as the further challenges for future work, and also conclude the paper.

II. PROBLEM DEFINITION AND RELATED WORK

In this section, we first formulate the problem of privacy preserving in social networks by anonymization. Then, we review the related work briefly.

A. Preliminaries

In this paper, we model a *social network* as a simple graph $G = (V, E, L, \mathcal{L})$, where V is a set of vertices, $E \subseteq V \times V$ is a set of edges, L is a set of labels, and a labeling function $\mathcal{L} : V \to L$ assigns each vertex a label. For a graph G, V(G), E(G), L_G , and \mathcal{L}_G are the set of vertices, the set of edges, the set of labels, and the labeling function in G, respectively. To keep our discussion simple, we assume that edges do not carry labels. However, our methods can be straightforwardly extended to remove this assumption.

The items in the label set L form a hierarchy. For example, if occupations are used as labels of vertices in a social network, L contains not only the specific occupations such as dentist, general physician, optometrist, high school teacher, and primary school teacher, but also general categories like medical doctor, teacher, and professional. We assume that there exists a meta symbol $* \in L$ which is the most general category generalizing all labels. For two labels $l_1, l_2 \in L$, if l_1 is more general than l_2 , we write $l_1 \prec l_2$. For example, medical doctor \prec optometrist. Moreover, $l_1 \preceq l_2$ if and only if $l_1 \prec l_2$ or $l_1 = l_2$. \preceq is a partial order on L.

For a graph G and a set of vertices $S \subseteq V(G)$, the *induced* subgraph of G on S is $G(S) = (S, E_S, L_G, \mathcal{L}_G)$ where $E_S = \{(u, v) | (u, v) \in E(G) \land u, v \in S\}.$

In a social network G, the *neighborhood* of $u \in V(G)$ is the induced subgraph of the neighbors of u, denoted by $Neighbor_G(u) = G(N_u)$ where $N_u = \{v | (u, v) \in E(G)\}.$

Given a graph $H = (V_H, E_H, L, \mathcal{L})$ and a social network $G = (V, E, L, \mathcal{L})$, an *instance* of H in G is a tuple (H', f) where $H' = (V_{H'}, E_{H'}, L, \mathcal{L})$ is a subgraph in G and $f : V_H \to V_{H'}$ is a bijection function such that (1) for any $u \in V_H$, $\mathcal{L}(f(u)) \leq \mathcal{L}(u)$, and (2) $(u, v) \in E_H$ if and only if $(f(u), f(v)) \in E_{H'}$. Literally, the first condition states that the corresponding labels in H' can be more general than those in H.

B. Problem Formulation

To define the problem of privacy preservation in publishing social network data, we need to formulate the following issues. First, we need to identify the privacy information to be preserved. Second, we need to model the background knowledge that an adversary may use to attack the privacy. Last, we need to specify the usage of the published social network data so that an anonymization method can try to retain the utility as much as possible while the privacy information is fully preserved.

Different formulations of the above issues may lead to different versions of privacy preservation in social networks.

Here, we propose a version which we believe is useful in many applications.

1) Privacy in Social Networks and Anonymization: In this paper, we are interested in preserving the privacy of individuals which are represented as vertices in a social network. Specifically, how a small subset of vertices are connected in a social network is considered as the privacy of those vertices.

Consider a social network $G = (V, E, L, \mathcal{L})$ and the anonymization $G' = (V', E', L', \mathcal{L}')$ for publishing. We assume that no fake vertices are added in the anonymization. That is, there is a bijection function $\mathcal{A} : V \to V'$. This assumption is often desirable in applications since introducing fake vertices may often change the global structure of a social network. Moreover, we assume that for $(u, v) \in E$, $(\mathcal{A}(u), \mathcal{A}(v)) \in E'$. That is, the connections between vertices in G are retained in G'.

For a vertex $u \in V$, if an adversary can identify a vertex $u' \in V'$ such that how u connects to other vertices in G is very similar to how u' connects to other vertices in G', and is substantially different from how any other vertices connect to others, then the privacy of u is leaked.

Therefore, privacy preservation in publishing social network data is to prevent any vertex $u \in V(G)$ from being reidentified in G' with high confidence. Technically, given a positive integer k, G' preserves the privacy in G if every vertex $u \in V(G)$ cannot be re-identified in G' with a confidence larger than $\frac{1}{k}$.

2) Adversary Background Knowledge: In order to attack the privacy of a target individual in the original network, i.e., analyze the released anonymization network and re-identify the vertex, an adversary needs some background knowledge. Equipped with different background knowledge, an adversary may conduct different types of attacks against privacy. Therefore, the assumptions of adversaries' background knowledge play a critical role in both modeling privacy attacks on social networks and developing anonymization strategies to protect privacy in social network data.

In this paper, we assume that an adversary may have the background knowledge about the neighborhood of some target individuals. This assumption is realistic in many applications. Among many types of information about a target victim that an adversary may collect to attack the victim's privacy, one essential piece of information easy to be collected is the neighborhood, i.e., what the neighbors of the victim are and how the neighbors are connected.

Generally, we can consider the *d*-neighbors of the target vertex, i.e., the vertices within distance *d* to the target vertex in the network where *d* is a positive integer. However, when *d* is large, collecting information about the *d*-neighbors of a target vertex may often be impractical for an adversary since the adversary may often have a limited access to a large social network. Moreover, as found in many social networks, the network diameter is often small. In other words, when d > 1, an adversary may have to collect information about many vertices. Therefore, we confine our discussion in this paper to the essential case where only the immediate neighbors, i.e., vertices in $Neighbor_G(u)$, are considered. The case of d > 1

is interesting for future work.

An adversary may attack the privacy using the neighborhoods. For a social network G, suppose an adversary knows $Neighbor_G(u)$ for a vertex $u \in V(G)$. If $Neighbor_G(u)$ has k instances in G' where G' is an anonymization of G, then u can be re-identified in G' with confidence $\frac{1}{k}$.

Similar to the philosophy in the k-anonymity model [5], to protect the privacy of vertices sufficiently, we want to keep the re-identification confidence lower than a threshold. Let k be a positive integer. For a vertex $u \in V(G)$, u is k-anonymous in anonymization G' if there are at least (k - 1) other vertices $v_1, \ldots, v_{k-1} \in V(G)$ such that $Neighbor_{G'}(\mathcal{A}(u))$, $Neighbor_{G'}(\mathcal{A}(v_1)), \ldots, Neighbor_{G'}(\mathcal{A}(v_{k-1}))$ are isomorphic. G' is k-anonymous if every vertex in G is k-anonymous in G'. Analogous to the correctness of k-anonymity model [5] on relational data, we have the following claim.

Theorem 1 (K-anonymity): Let G be a social network and G' an anonymization of G. If G' is k-anonymous, then with the neighborhood background knowledge, any vertex in G cannot be re-identified in G' with confidence larger than $\frac{1}{k}$.

An adversary knowing the neighborhood of a target vertex is a strong assumption. Provided privacy is preserved under this assumption, privacy is also preserved when an adversary knows only part of the neighborhood (i.e., only some neighbors and some connections among neighbors) of a target vertex.

3) Usage of Anonymized Social Networks: An important aspect of anonymizing social network data is how the anonymized networks are expected to be used. Different applications may have different expectations. In some situations, anonymized networks may be used to analyze the global structures. In some other situations, anonymized networks may be used to analyze the micro-structures. Clearly, different usage expectations may lead to different anonymization schemes.

In this paper, we focus on using anonymized social networks to answer aggregate network queries. An aggregate network query computes the aggregate on some paths or subgraphs satisfying some given conditions. As an example, suppose a user is interested in the average distance from a medical doctor vertex to a teacher vertex in a network. For each doctor vertex, we can find the nearest neighbor vertex that is a teacher. Then, the aggregate network query returns the average of the distance between a doctor vertex to its nearest teacher neighbor.

Aggregate network queries are useful in many applications, such as customer relationship management. While many types of queries on social networks are interesting, we are particularly interested in aggregate network queries in this paper since typically detail data is needed to answer such queries accurately. Using aggregate network queries we can examine the effectiveness of social network anonymization in a meaningful way.

4) Problem Definition and Complexity: Given a social network G, we want to compute an anonymization G' such that (1) G' is k-anonymous; (2) each vertex in G is anonymized to a vertex in G' and G' does not contain any fake vertex; (3) every edge in G is retained in G'; and (4) G' can be used to answer aggregate network queries as accurately as possible.

To understand the difficulty of the problem, we consider a simple case where all vertices in G carry the same label, or equivalently, G is not labeled.

Theorem 2 (Complexity): The following *k*-anonymity problem in social network is NP-hard.

Instance: a social network $G = (V, E, L, \mathcal{L})$ where L = * and $\forall u \in V, \mathcal{L}(u) = *$, positive integers k and n.

Question: is there an anonymized social network $G' = (V, E', L, \mathcal{L})$ such that $E \subset E'$, |E' - E| = n, and G' is k-anonymous?

Proof sketch. The proof is constructed by reducing the NPhard *k*-DIMENSIONAL PERFECT MATCHING problem [7] to the *k*-anonymity problem in social networks. Limited by space, we omit the details here.

C. Related Work

Privacy becomes a more and more serious concern in many applications. The development of techniques that incorporate privacy concerns has become a fruitful direction for database and data mining research.

One of the privacy concerned problems is publishing microdata for public use [8], which has been extensively studied recently. A large category of privacy attacks is to re-identify individuals by joining the published table with some external tables modeling the background knowledge of users. To battle this type of attacks, the mechanism of k-anonymity was proposed in [9], [5]. Specifically, a data set is said to be k-anonymous ($k \ge 1$) if, on the quasi-identifier attributes (i.e., the minimal set of attributes in the table that can be joined with external information to re-identify individual records), each record is indistinguishable from at least k - 1 other records within the same data set. The larger the value of k, the better the privacy is protected.

Machanavajjhala *et al.* [6] showed that a k-anonymized dataset has some subtle but severe privacy problems due to the lack of diversity in the sensitive attributes. In particular, they showed that, the degree of privacy protection does not really depend on the size of the quasi-identifier attribute set. Instead, it is determined by the number of distinct sensitive values associated with each quasi-identifier attribute set. The observation leads to the notion of l-diversity [6]. [10] proved that l-diversity always guarantees stronger privacy preservation than k-anonymity.

In this paper, we focus on k-anonymity since k-anonymity is the most essential and most applicable privacy model, which can be used even when sensitive attributes are not defined. A few governmental privacy regulations including HIPAA and European Union Data Directive adopted k-anonymity.

Beyond microdata, some other data sources such as social network data also have privacy concerns when they are published for public use. Typically, social network data can be represented as a graph, in which vertices correspond to people or other social entities, and edges correspond to social links between them [11]. As a first step to hide information about social entities while preserving the global network properties, the released social network data has to go through the anonymization procedure which replaces social entity names with meaningless unique identifiers [12]. Although this kind of anonymization can exactly preserve the unannotated structure of the social network, it may still leak a lot of information.

Attacks in social network data can be regarded as one kind of link mining [13]. Specifically, as a pioneer work about privacy in social network data, Backstrom *et al.* [12] described a family of attacks based on random graph theory. For example, an attacker may plant some well constructed substructures associated with the target entities in advance. Once the social network data is collected and published, the attacker can first try to identify the planted structures and thus peek the linkage between the target vertices. However, there is no practical solution proposed in [12] to counter those attacks.

The attacks proposed in [12] are different from the neighborhood attacks addressed in this paper. The attacks in [12] need to plant a set of deliberative structures before the social network data is anonymized, which is a task hard to achieve in some situations. As shown before, even without planting deliberative structures, the released social network data is still in danger, as neighborhood attacks are still possible.

Wang *et al.* [14] adopted description logic as the underlying knowledge representation formalism, and proposed some metrics of anonymity for assessing the risk of breaching confidentiality by disclosing social network data. However, they did not give any anonymization algorithms for social network data.

Simultaneous to our study, Hay *et al.* [15] presented a framework for assessing the privacy risk of sharing anonymized network data. They modeled the adversaries' background knowledge as vertex requirement structural queries and subgraph knowledge structural queries, and proposed a privacy requirement k-candidate anonymity which is similar to k-anonymity in tabular data. They developed a random graph perturbation method by randomly deleting or adding edges to anonymize a social network. Their model assumes that the nodes and the edges in a social network are not labeled.

Recently, Zheleva *et al.* [16] proposed a model different from ours. They focused on social networks where nodes are not labeled but edges are labeled. Some types of edges are sensitive and should be hidden. They provided the edge anonymization methods based on edge clustering and removal to prevent link re-identification.

III. AN ANONYMIZATION METHOD

In this section, we introduce a practical method to anonymize a social network to satisfy the k-anonymity requirement. The method is in two steps.

First, we extract the neighborhoods of all vertices in the network. To facilitate the comparisons among neighborhoods of different vertices including the isomorphism tests which will be conducted frequently in anonymization, we propose a simple yet effective *neighborhood component coding technique* to represent the neighborhoods in a concise way.

In the second step, we greedily organize vertices into groups and anonymize the neighborhoods of vertices in the same group. Due to the well recognized power law distribution of



Fig. 2. Neighborhood and neighborhood components (the dashed edges are just for illustration and are not in the neighborhood subgraph).



the degrees of vertices in large social networks, we start with those vertices of high degrees.

A. Neighborhood Extraction and Coding

In order to meet the k-anonymity requirement, we need to put vertices into groups and anonymize the neighborhoods of vertices in a group. Ideally, vertices having similar neighborhoods should be grouped together. As the first step, we extract neighborhoods of vertices and represent them in a concise way.

Extracting the neighborhood of a vertex is straightforward. The challenge is how we can represent the neighborhood information to facilitate the later operations in anonymization. Since we need to anonymize all neighborhoods of the vertices in one group to the same, isomorphism tests are frequently conducted.

The general graph isomorphism problem which determines whether two graphs are isomorphic is NP-hard [17]. Here, we propose a coding technique for neighborhood subgraphs so that whether two neighborhoods are isomorphic can be determined by the corresponding coding.

In a social network G, a subgraph C of G is a *neighborhood* component of $u \in V(G)$ if C is a maximal connected subgraph in $Neighbor_G(u)$.

Example 1 (Neighborhood component): Figure 2 shows $Neighbor_G(u)$, the neighborhood of a vertex u. $Neighbor_G(u)$ contains three neighborhood components, C_1 , C_2 , and C_3 as shown in the figure.

Clearly, the neighborhood of a vertex can be divided into neighborhood components. To code the whole neighborhood, we need to first code each component.

The *depth-first search tree* (*DFS-tree* for short) [18] is popularly used for navigating connected graphs. Thus, it is natural to encode the edges and vertices in a graph based on its DFS-tree. All the vertices in G can be encoded in the preorder of T. However, the DFS-tree is generally not unique for a graph. That is, there can be multiple DFS-trees corresponding to a given graph. For example, Figures 3(b) and 3(c) show two DFS-trees of the graph G in Figure 3(a). The thick edges in Figures 3(b) and 3(c) are those in the DFS-trees, and are called the *forward edges*, while the thin edges are those not in the DFS-trees, and are called the *backward edges*. The vertices in the graph are encoded v_0 to v_3 according to the pre-order of the corresponding DFS-trees.

To solve the uniqueness problem, a minimum DFS code notation is proposed in [19]. For any connected graph G, let T be a DFS-tree of G. Then, an edge is always listed as (v_i, v_j) such that i < j. A linear order \prec on the edges in G can be defined as follows. Given edges $e = (v_i, v_j)$ and $e' = (v_{i'}, v_{j'})$. $e \prec e'$ if (1) when both e and e' are forward edges (i.e., in DFS-tree T), j < j' or $(i > i' \land j = j')$; (2) when both e and e' are backward edges (i.e., edges not in DFS-tree T), i < i' or $(i = i' \land j < j')$; (3) when e is a forward edge and e' is a backward edge, $j \leq i'$; or (4) when e is a backward edge and e' is a forward edge, i < j'.

For a graph G and a DFS-tree T, a list of all edges in E(G) in order \prec is called the DFS code of G with respect to T, denoted by code(G,T). For example, the DFS code with respect to the DFS-tree T_1 in Figure 3(b) is $code(G,T_1) = \langle (v_0, v_1, x, x) - (v_1, v_2, x, z) - (v_2, v_0, z, x) - (v_1, v_3, x, y) \rangle$, where an edge (v_i, v_j) is written as $(v_i, v_j, \mathcal{L}(v_i), \mathcal{L}(v_j))$, i.e., the labels are included. Similarly, the DFS code with respect to the DFS-tree T_2 in Figure 3(c) is $code(G,T_2) = \langle (v_0, v_1, y, x) - (v_1, v_2, x, x) - (v_2, v_3, x, z) - (v_3, v_1, z, x) \rangle$.

Suppose there is a linear order over the label set L. Then, for DFS-trees T_1 and T_2 on the same graph G, their DFS codes can be compared lexically according to the vertex pairs as labels of edges. For example, we have $code(G, T_1) < code(G, T_2)$ in Figures 3(b) and 3(c).

The lexically *minimum DFS code* is selected as the representation of the graph, denoted by DFS(G). In our example in Figure 3, $DFS(G) = code(G, T_1)$.

Minimum DFS code has a nice property [19]: two graphs G and G' are isomorphic if and only if DFS(G) = DFS(G'). Using minimum DFS code, we can code every component of the neighborhood of a vertex. Now, the problem becomes how we can combine the minimum DFS codes of all components in a neighborhood into one code.

For two neighborhood components C_i and C_j in $Neighbor_G(u)$, we define $C_i \prec C_j$ if (1) $|V(C_i)| < |V(C_j)|$; or (2) $|V(C_i)| = |V(C_j)|$ and $|E(C_i) < |E(C_j)|$; or (3) $|V(C_i)| = |V(C_j)|$, $|E(C_i) = |E(C_j)|$, and $DFS(C_i)$ is smaller than $DFS(C_j)$.

Based on the neighborhood component order, we can assign a canonical label for each neighborhood. In a social network G, for vertex $u \in V(G)$, the neighborhood component code of Neighbor_G(u) is a vector $NCC(u) = (DFS(C_1), \ldots, DFS(C_m))$ where C_1, \ldots, C_m are the neighborhood components of Neighbor_G(u), i.e., Neighbor_G(u) = $\bigcup_{i=1}^{m} C_i, C_i \leq C_j$ for $1 \leq i < j \leq m$.

Example 2 (Neighborhood component code): In Figure 2, the neighborhood component code of $Neighbor_G(u)$ is $NCC(u) = (DFS(C_1), DFS(C_2), DFS(C_3)).$

Using neighborhood component code, we can easily identify isomorphic neighborhoods.

Theorem 3 (Neighborhood component code): For two vertices $u, v \in V(G)$ where G is a social network, $Neighbor_G(u)$ and $Neighbor_G(v)$ are isomorphic if and only if NCC(u) = NCC(v).

Using neighborhood component code to label and index neighborhoods has some advantages. First, it is easy to examine whether a group of neighborhoods are isomorphic. Second, since each neighborhood is decomposed into several neighborhood components, it is easy to calculate the structure similarity between two neighborhoods. Third, we can find similar components between two neighborhoods. Anonymizing similar components can lead to low information loss and high similarity between the anonymization and the original social network.

B. Social Network Anonymization

One major challenge in anonymizing a social network is that changing labels of vertices and adding edges may affect the neighborhoods of some other vertices as well as the properties of the network. It has been well recognized that the following two properties often hold in practical social networks. The properties help us in designing anonymization methods.

Property 1: vertex degree in power law distribution. The degrees of vertices in a large social network often follow the power law distribution [20]. Such degree distributions have been identified in various social networks including Internet, biological networks, and co-authorship networks.

Property 2: the "small-world phenomenon" [11]. It is also popularly known as "six degrees of separation", which states that large social networks in practice often have surprisingly small average diameters.

Our social network anonymization method processes vertices in the degree descending order, and utilizes the above two properties of large social networks in practice.

The k-anonymity requires that each vertex $u \in V(G)$ is grouped with at least (k - 1) other vertices such that their anonymized neighborhoods are isomorphic. For a group S of vertices having the isomorphic anonymized neighborhoods, all vertices in S have the same degree. Since the degrees of vertices in a large social network follow a power law distribution, only a small number of vertices have a high degree. Processing those vertices of high degrees first can keep the information loss about those vertices low. There are often many vertices of a low degree. It is relatively easy to anonymize those low degree vertices and retain high quality. Moreover, as will be shown soon, low degree vertices can be used to anonymize those high degree vertices and do not affect the diameters of the network too much.

1) Anonymization Quality Measure: In our social network anonymization model, there are two ways to anonymize the neighborhoods of vertices: generalizing vertex labels and adding edges. Each of the two methods leads to some information loss.

The information loss due to generalization of vertex labels can be measured by the normalized certainty penalty [21]. Consider a vertex u of label l_1 , where l_1 is at the leaf level of the label hierarchy, i.e., l_1 does not have any descendant. Suppose l_1 is generalized to l_2 for u where $l_2 \prec l_1$. Let $size(l_2)$ be the number of descendants of l_2 that are leafs in the label hierarchy, and size(*) be the total number of leafs in the label hierarchy. Then, the normalized certainty penalty of l_2 is $NCP(l_2) = \frac{size(l_2)}{size(*)}$.

The information loss due to adding edges can be measured by the total number of edges added and the number of vertices that are not in the neighborhood of the target vertex and are linked to the anonymized neighborhood for the purpose of anonymization.

Consider two vertices $u_1, u_2 \in V(G)$ where G is a social network. Suppose $Neighbor_G(u_1)$ and $Neighbor_G(u_2)$ are generalized to $Neighbor_{G'}(\mathcal{A}(u_1))$ and $Neighbor_{G'}(\mathcal{A}(u_2))$ such that $Neighbor_{G'}(\mathcal{A}(u_1))$ and $Neighbor_{G'}(\mathcal{A}(u_2))$ are isomorphic. Let $H = Neighbor_G(u_1) \cup Neighbor_G(u_2)$ and $H' = Neighbor_{G'}(\mathcal{A}(u_1)) \cup Neighbor_{G'}(\mathcal{A}(u_2))$. The anonymization cost is

$$Cost(u, v) = \alpha \cdot \sum_{v' \in H'} NCP(v') +\beta \cdot |\{(v_1, v_2) | (v_1, v_2) \notin E(H), (\mathcal{A}(v_1), \mathcal{A}(v_2)) \in E(H')\}| +\gamma \cdot (|V(H')| - |V(H)|)$$

where α , β and γ are the weights specified by users. Literally, the cost consists of three parts. The first part is the normalized certainty penalty measuring the information loss of generalizing labels of vertices. The second part measures the information loss due to adding edges. The last part counts the number of vertices that are linked to the anonymized neighborhoods to achieve k-anonymity.

The anonymization cost of two vertices u and v measures the similarity between $Neighbor_G(u)$ and $Neighbor_G(v)$. The smaller the anonymization cost, the more similar the two neighborhoods.

2) Anonymizing Two Neighborhoods: Now, let us consider a greedy method to anonymize two neighborhoods $Neighbor_G(u)$ and $Neighbor_G(v)$.

We first find all perfect matches of neighborhood components in $Neighbor_G(u)$ and $Neighbor_G(v)$. Two components perfectly match each other if they have the same minimum DFS code. Those perfect matches are marked as "matched" and pass over for further consideration.

For example, consider two vertices u and v whose neighborhoods are shown in Figure 4. Each vertex is shown in the form of (id, label). The neighborhood component $C_2(u) \in Neighbor_G(u)$ perfectly matches $C_3(v) \in Neighbor_G(v)$.

For those unmatched components, the anonymization algorithm tries to pair similar components and anonymize them. The similarity between two components is based on the anonymization cost. To calculate the similarity between two components, we try to match similar vertices in the two components as much as possible. This is a traditional substructure similarity search problem, which has been proved NP-hard [22]. Instead of computing the optimal matching, we conduct a greedy match.

To start with, we first try to find two vertices with the same degree and the same label in the two components to be



Fig. 4. Anonymizing two neighborhoods.

matched. If there are multiple matching vertex pairs, the pair with the highest vertex degree is chosen. If there is no such a pair of matching vertices, we relax the matching requirement (vertex degree and label), calculate the difference of degrees and the normalized certainty penalty of generalizing the labels in the label hierarchy, and choose the one with the minimum anonymization cost. Then, we conduct a breadth-first search to match vertices one by one, until all possible vertex matchings are found. The anonymization cost is calculated according to the matching, and is used to measure the similarity of the two components.

Consider components $C_1(u)$ and $C_1(v)$ in Figure 4. Vertices u_1 and v_1 match. We start from these two vertices and conduct a breadth-first search. Vertex v_2 partially matches vertex u_2 . Vertex v_3 partially matches vertex u_3 . The vertex matching stops since all possible vertex matchings are found. However, vertex u_4 does not find any vertex matching in $C_1(v)$. Thus we have to find a vertex $w_1 \in V(G)$ that is neither in $C_1(v)$ nor in $C_1(u)$, and add it into $C_1(v)$, so that $C_1(u)$ and $C_1(v)$ can be anonymized to the same.

When a vertex has to be introduced into the neighborhood for the sake of anonymization, the following rules are used: we first consider those vertices in V(G) that are unanonymized. The vertex with smallest degree has the highest priority. If there are more than one candidate with the same smallest degree, we choose the one having the closest label in terms of normalized certainty penalty. If we cannot find any other vertex that is unanonymized, we select one anonymized vertex wwith the smallest degree and satisfying the label requirement, and mark w and its (k - 1) other vertices anonymized in the same group as "unanonymized".

In our example, suppose we can find an unanonymized vertex (w_1, l_4) to be added to $C_1(u)$, the anonymization cost of $C_1(u)$ and $C_1(v)$ is $\alpha \cdot \sum_{v' \in V(C_1(u)) \cup V(C_1(v))} NCP(\mathcal{L}(\mathcal{A}(v')) + \beta \cdot 1 + \gamma \cdot 1 = \alpha \cdot \frac{4}{5} + \beta + \gamma.$

Input: a social network
$$G = (V, E)$$
, the anonymization requirement parameter k, the cost function parameters α , β and γ ;

Output: an anonymized graph G';

Method:

- 1: initialize G' = G;
- 2: mark $v_i \in V(G)$ as "unanonymized";
- 3: sort $v_i \in V(G)$ as *VertexList* in neighborhood size descending order;
- 4: WHILE (VertexList $\neq \emptyset$) DO
- 5: let *SeedVertex = VertexList.head()* and remove it from *VertexList*;
- 6: FOR each $v_i \in VertexList$ DO
- 7: calculate *Cost(SeedVertex, v_i)* using the anonymization method for two vertices; END FOR
- 8: IF (*VertexList.size*() $\geq 2k 1$) DO let *CandidateSet* contain the top k - 1 vertices with the smallest *Cost*;

- let CandidateSet contain the remaining unanonymized vertices;
- 11: suppose *CandidateSet*= $\{u_1, \ldots, u_m\}$, anonymize *Neighbor(SeedVertex)* and *Neighbor(u_1)* as discussed in Section III-B.2;
- 12: FOR j = 2 to m DO

13: anonymize
$$Neighbor(u_j)$$
 and $\{Neighbor(SeedVertex), Neighbor(u_1), \ldots, Neighbor(u_{j-1})\}$ as discussed in Section III-B.2, mark them as "anonymized";

14: update VertexList; END FOR END WHILE

Fig. 5. Anonymizing a Social Network.

Based on the component similarity, we can pair similar components. We start with the component with the largest number of vertices. This component is paired with the most similar component in the other neighborhood. The two paired components are anonymized to the same, marked "matched", and removed from consideration. The matching continues until all components in one neighborhood are marked "matched".

If there are some components left in the other neighborhood say $Neighbor_G(u)$, we use some other vertices in V(G) that are not in $Neighbor_G(u)$ to construct a component and add it to $Neighbor_G(u)$ to construct the matching and anonymization. The vertices are selected using the same criteria as selecting vertices to match two components.

We anonymize each pair of matched neighborhood components to the same. The two neighborhoods then are anonymized. For example, in Figure 4, the algorithm matches components $C_1(u)$ and $C_1(v)$, and $C_2(v)$ and $C_3(u)$ in turn. As a result, two vertices w_1 and w_2 from V(G) have to be added into components $C_1(v)$ and $C_3(u)$, respectively.

3) Anonymizing a Social Network: We propose a greedy method to anonymize a social network as shown in Figure 5.

First, we mark all vertices in the network as "unanonymized". We maintain a list *VertexList* of "unanonymized" vertices in the neighborhood size descending order: for vertices $u, v \in V(G)$, if $|V(Neighbor_G(u))|$ $< |V(Neighbor_G(v))|$, or $|V(Neighbor_G(u))| =$ $|V(Neighbor_G(v))|$ and $|E(Neighbor_G(u))| <$

^{9:} ELSE

 $|E(Neighbor_G(v))|$, then v precedes u in the list. If their neighborhoods have the same numbers of vertices and edges, they can be ordered arbitrarily.

Iteratively, we pick the first vertex SeedVertex in the list VertexList. The anonymization cost of SeedVertex and any other vertices in VertexList is calculated using the anonymization method for two vertices discussed in Section III-B.2. If the number of unanonymized vertices in VertexList is at least 2k - 1, we select a set CandidateSet of the top k - 1 vertices in VertexList with the smallest anonymization cost. We can easily give a lower bound of the anonymization cost based on the number of vertices and the number of edges in two neighborhoods. Since all vertices in VertexList have a neighborhood size smaller than or equal to that of SeedVertex, we scan VertexList in the neighborhood size descending order, and stop once the lower bound of the anonymization cost exceeds the cost of the current (k - 1)-th most similar vertex.

The SeedVertex and the vertices in CandidateSet= $\{u_1, \ldots, u_m\}$ are anonymized in turn using the anonymization method for two vertices discussed in Section III-B.2. The anonymization of SeedVertex and u_1 is straightforward. After these two vertices are anonymized, their neighborhoods are identical. When we anonymize them with respect to u_2 , any change (e.g., adding an edge or a neighbor node) to the neighborhood of SeedVertex will be applied to u_1 as well, so that the neighborhoods of SeedVertex, u_1 and u_2 are anonymized to the same. The process continues until the neighborhoods of SeedVertex and u_1, \ldots, u_m are anonymized.

During the anonymization of a group of vertices, some changes may occur to some other vertices v that have been marked as "anonymized" in another group (e.g., adding edges between an anonymized vertex and a vertex being anonymized based on vertex matching). In order to maintain the k-anonymity for those vertices, we apply the same changes to every other k-1 vertices having the isomorphic neighborhoods as v. Once those k vertices are changed, they are marked as "unanonymized" and inserted into the VertexList again.

When the number of unanonymized vertices in *VertexList* is less than 2k, to satisfy the *k*-anonymity, the remaining vertices in *VertexList* have to be considered together in anonymization. They are added to *CandidateSet* in a batch.

The social network anonymization algorithm continues until all the vertices in the graph are marked as "anonymized".

Theorem 4 (Termination): The algorithm in Figure 5 terminates for a finite social network of at least k vertices.

Proof sketch. Clearly, a clique of n vertices where each vertex is labeled * is k-anonymous provided $n \ge k$. In each iteration such that *VertexList* is not shortened, the algorithm either adds some edges into the network, or generalizes the labels of some vertices towards *. In the worst case, the network will be anonymized to a clique.

Surprisingly, as shown in our experiments, the algorithm can stop very quickly in practice, and the anonymization cost is relatively small. This is due to the two important properties of real social networks (vertex degree in power law distribution and small world phenomenon). Using different synthetic data sets, we find that most of the time we do not

k	Removing labels	Generalizing to affiliations
5	1.3%	12.7%
10	3.9%	16.1%
15	7.1%	19.4%
20	12.0%	23.2%

TABLE I

THE PERCENTAGES OF VERTICES VIOLATING *k*-ANONYMITY IN THE CO-AUTHORSHIP DATA.

need to anonymize the remaining vertices that are less than 2k, since they have the same labels, very low degree (1 in many sparse social networks), and isomorphic neighborhoods.

IV. EMPIRICAL EVALUATION

In this section, we report a systematic empirical study to evaluate our anonymization method using both real data sets and synthetic data sets. All the experiments were conducted on a PC computer running the Microsoft Windows XP SP2 Professional Edition operating system, with a 3.0 GHz Pentium 4 CPU, 1.0 GB main memory, and a 160 GB hard disk. The program was implemented in C/C++ and was compiled using Microsoft Visual Studio .NET 2005.

A. Neighborhood Attacks in Real Data

We used a real co-authorship data set from KDD Cup 2003 to examine whether neighborhood attacks may happen in practice. The data set was from the e-print arXiv (arXiv.org), and contains a subset of papers in the high-energy physics section of the arXiv. The LATEX sources of all papers are provided. We extracted author names from the data sources and constructed a co-authorship graph. Each vertex in the graph represents an author, and two vertices are linked by an edge if the two corresponding authors co-authored at least one paper in the data set. There are 57, 448 vertices and 120, 640 edges in the co-authorship graph and the average number of vertex degrees is about 4.

We tested two ways to preserve the privacy of authors by generalizing labels. In the first method, we removed all labels, i.e., author names. In the second method, we use the author's affiliation as the label, i.e., all authors from the same institution have the same label. After generalization, for each vertex, we extracted its neighborhood and counted the number of other vertices in the graph that have the isomorphic neighborhoods. Table I shows the percentages of vertices whose neighborhoods violate the k-anonymity.

Table I clearly shows that the neighborhood attacks are a real issue for social network data publishing. When the value of k increases, the number of vertices violating k-anonymity increases. Moreover, the more specific are the vertex labels, the more vertices fail the k-anonymity. Anonymizing labels only cannot prevent neighborhood attacks effectively.

B. Anonymization Performance

We use the R-MAT graph model [23] to generate synthetic data sets. R-MAT can generate graphs with power law vertex degree distribution and small-world characteristic, which are



Fig. 8. The runtime on various synthetic data sets.

the two most important properties for many real-world social networks.

R-MAT takes 4 probability parameters, namely a, b, c and d. Consider the adjacency matrix representation A of the graph. Assume that each element a_{ij} in A is non-zero if there exists an edge between vertices i and j. Given the number of vertices n and the number of edges m, R-MAT starts with an empty $n \times n$ adjacency matrix A. It recursively divides the adjacency matrix into four equal-size partitions, and distributes edges within these partitions with a set of unequal probabilities. Each edge chooses one of the four partitions with probabilities a, b, c and d, respectively, such that a + b + c + d = 1. The parameters a, b, c and d can be determined based on the required community structure and vertex degree distribution, as detailed in [23]. The study [23] conjectures that the ratios a/b and a/c are approximately 3/1 in many real world graphs, and $a \ge d$.

We used the default values of 0.45, 0.15, 0.15 and 0.25 for those four parameters, respectively. We generated a series of synthetic data sets by varying the number of vertices from 5,000 to 25,000 and the average vertex degree from 3 to 7. Hereafter, we refer to a synthetic data set as D(n,d), where nand d are the number of vertices and the average vertex degree, respectively. The edge weight was set in the range [0,100] by default. We assigned each vertex a label based on its average edge weight. A simple two level hierarchy structure for those labels was generated.

We first examined the effect of parameters α , β and γ in the anonymization quality measure. β was set to 1 as the base. We changed the values of α and γ , and measured the number of edges added and the normalized certainty penalty incurred in the anonymized graphs. The results for data set D(15K, 5)and k = 10 are shown in Figure 6. The results show that we can trade off between adding edges and generalizing labels by tuning the three parameters. Often adding less edges is more desirable in anonymizing a social network since the network structures can be preserved better. We observed that, on the synthetic data sets, when $\alpha = 100$ and $\gamma = 1.1$, the number of edges added is small and the normalized certainty penalty is moderate. Hereafter, we use 100, 1, and 1.1 as the default values for α , β and γ .

Figure 7 reports the anonymization quality on various synthetic data sets with respect to different k values, and shows the anonymization cost in both the number of edges added and the normalized certainty penalty. First, when the number of vertices increases, the anonymization cost increases. However, the increase of the number of edges added is sublinear, since in a larger network it is more likely to find similar neighborhoods. Second, when k increases, the anonymization cost also increases, because more neighborhoods need to be anonymized in a group. Last, when the average number of vertex degree increases, the anonymization cost increases, too. In a denser network, the neighborhoods are more diverse and more edges are needed to anonymize different neighborhoods.

Figure 8 shows the runtime on various synthetic data sets with respect to different k values. The runtime increases when the average vertex degree increases, since the network becomes denser. Moreover, the larger the k, the longer the runtime since more neighborhoods in a group need to be anonymized.

C. Anonymizing the Co-authorship Data Set

We built a 3-level label hierarchy for the KDD cup 2003 co-authorship data set. The leaf vertices are those labels with author affiliations. The middle level contains the countries of the authors' affiliations. The root is the most general label *. We anonymized the co-authorship data set using our anonymization method. The number of edges added, the normalized certainty penalty and the runtime with different values of k are shown in Figure 9. Comparing to the total number of edges in the data set (120, 640), the number of edges added is less than 6% even when k = 20. Moreover, the runtime is scalable with respect to k.

D. Aggregate Query Answering

To test the utility of the anonymized social networks, we conducted aggregate network queries on the KDD Cup 2003 co-authorship data set, and the anonymized networks. For two labels l_1 and l_2 in the data set, we calculated the average



Fig. 10. Query answering on the KDD Cup 2003 co-authorship data set.

Fig. 9. Anonymizing the KDD cup 2003 co-authorship data set.

distance from a vertex with label l_1 to its nearest vertex with label l_2 . Since labels are organized in a hierarchy structure, when we calculated the average distance, we also considered distance from vertices with labels l'_1 to vertices with labels l'_2 such that $l_1 \leq l'_1$ and $l_2 \leq l'_2$. The error rate is $\frac{d-d'}{d}$, where dand d' are the average distances in the original network and in the anonymized network, respectively. We randomly picked 10 label pairs from the label hierarchy, and calculated the average error rate of them. The results are shown in Figure 10. After the anonymization, with some edges added, the average distance decreases. Therefore, the error rate is always positive. However, the error rate is small even when k is up to 20, since the number of edges added is small, as shown in Figure 9.

V. DISCUSSION AND CONCLUSIONS

In this paper, we tackled the novel and important problem of preserving privacy in social network data, and took an initiative to combat neighborhood attacks. We modeled the problem systematically and developed a practically feasible approach. An extensive empirical study using both a real data set and a series of synthetic data sets strongly indicated that neighborhood attacks are real in practice, and our method is highly feasible. Moreover, anonymized social networks can still be used to answer aggregate queries accurately.

As social network data is much more complicated than relational data, privacy preserving in social networks is much more challenging and needs many serious efforts in the future. Particularly, modeling adversarial attacks and developing privacy preservation strategies are critical. For future work, we believe that the following two types of attacks should be addressed systematically.

We only handle 1-neighborhoods in this paper. It is very interesting and could be desirable in some applications that d-neighborhoods (d > 1) are protected. However, this will introduce a serious challenge in computation. The neighborhood size increases exponentially as d increases. Isomorphism tests and anonymization of large neighborhoods are very challenging.

A k-anonymous social network still may leak privacy. If an adversary can identify a victim in a group of vertices anonymized in a group, but all are associated with some sensitive information, then the adversary still can know that sensitive attribute of the victim. Some mechanism analogous to l-diversity [6] should be introduced.

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