

## UNIVERSITÀ DI PISA

Facoltà di Scienze Matematiche, Fisiche e Naturali Corso di Laurea Triennale in Informatica

#### Tesi di Laurea

# SUBGRAPH SIMILARITY IN COMPLEX NETWORKS

SIMILARITÀ DI SOTTOGRAFI NELLE RETI COMPLESSE

Relatori:	Candidato:
Prof. <i>Roberto Grossi</i> Prof. <i>Andrea Marino</i>	Gaspare Ferraro

ANNO ACCADENICO 0016 0017

ANNO ACCADEMICO 2016-2017

Ai miei genitori per non avermi tagliato i viveri

# Contents

1	Intr	oduction 1
	1.1	Basic definitions
	1.2	The problem
	1.3	Pratical applications
<b>2</b>	Bas	ic tools 4
	2.1	Similarity indices
	2.2	Documents similarity
	2.3	Graphs similarity
	2.4	Subgraphs similarity
	2.5	Sketches
		2.5.1 min-wise permutation
		2.5.2 bottom-k sketches
	2.6	Color Coding
3	Con	nputation of subgraph similarity 6
_	3.1	Naive approach
	3.2	Efficient computation
	3.3	Baseline algorithm
4	Pro	ject development 8
_	4.1	Implementation choices
	4.2	Dataset
	4.3	Experimental results
Δ	Cod	le snippets 10
А		Color Coding
		<u> </u>
		Colorful sampling
		Frequency count
	A.4	Frequency sampling
	$\rightarrow$	AUDURIUM INCUCES 1.3

## Introduction

With the spread of Internet and more importantly of the social networks, efficient data analysis becomes increasingly important. Graphs are a powerful data structure that model in a natural way information .

#### 1.1 Basic definitions

**Definition 1.1.** A graph is a pair of sets G = (V, E), where V is the set of vertices (or nodes) and  $E \subset V \times V$  is the set of edges.

If two vertices  $u, v \in V$  are connected by an edge they are called extreme of the edge, in this case we denote the edge with the pair  $(u, v) \in E$ 

If  $(u, v) \in E \Leftrightarrow (v, u) \in E$  the graph is called undirected, where not specified we will only deal with undirected graphs.

A sequence of nodes  $v_1, v_2, \ldots, v_k$  is called path if  $(v_i, v_{i+1}) \in E \ \forall i = 1, \ldots, k-1$ ; a path is called simple if  $v_i \neq v_j \ \forall i, j \ 1 \leq i < j \leq k$ . A cycle is a path where  $(v_k, v_1) \in E$ .

We denote by  $N(u) = \{v : (u, v) \in E\}$  the set of neighbors of the vertex u, the cardinality of this set is called degree of u (deg u = |N(u)|).

With  $N^{< k}(u)$  we indicate the set of vertex connected to u by a simple path of length less than k (note that  $N(u) = N^{<2}(u)$ ).

**Definition 1.2.** A graph G' = (V', E') is called subgraph of G = (V, E) if  $V' \subset V$  and  $E' \subset E$ . A subgraph is called induced if  $E' = (V' \times V') \cap E$ .

We use  $G' \subset G$  to indicate that the graph G' is a subgraph of G and G' < G to indicate that the graph G' is a induced subgraph of G.

Note that an induced subgraph G' = (V', E') can be uniquely identified by the set of its vertex V'.

2 Introduction

**Definition 1.3.** A labeled graph is a triple (V, E, L) where (V, E) is a graph and  $L: V \to \Sigma$  is a function that assign for every node v a symbol of the alphabet  $\Sigma$ . We call  $L(u) \in \Sigma$  label of the node u.

Given a path  $\pi = v_1, v_2, \dots, v_k$  we extend the function L and we indicate with  $L(\pi) = L(v_1)L(v_2) \dots L(v_k) \in \Sigma^k$  the string obtained by the concatenation of the labels of the nodes in the path.

In this thesis we mainly focus to analyze complex network: special graph with a non-trivial topology like random graph. Complex network occur in graphs modelling real system like social networks or computer networks and are characterized by a specific structural features:

**Definition 1.4.** We define as power-law degree distribution a networks where the degree of a node u follow, for some  $\gamma$  (usually  $2 < \gamma < 3$ ), the probability:

$$P(deg(u) = k) \sim k^{-\gamma} \tag{1.1}$$



Figure 1.1: Degree distribution of a random network



Figure 1.2: Degree distribution of a scale-free complex network



Figure 1.3: Random network with |N| = 100 and |E| = 1000



Figure 1.4: Complex network with |N| = 100 and |E| = 1000

#### 1.2 The problem

**Problem 1.5.** Given an undirected labeled graph G = (V, E, L) over an alphabet  $\Sigma$ , an integer q and two set of nodes  $V_1, V_2 \subset V$ , we want to estimate the similarity between the two induced subgraphs  $V_1, V_2 < G$  based on the labels frequency of simple paths with nodes in  $V_1 \cup N^{< q}(V_1)$  and  $V_2 \cup N^{< q}(V_2)$ .

Will discuss about a more formal and rigorous definition of subgraphs similarity in chapter 2.

In the definition we use  $V_1 \cup N^{< q}(V_1)$  and  $V_2 \cup N^{< q}(V_2)$  instead of simply  $V_1$  and  $V_2$  because in a complex graph we also want to keep in mind of the interaction between the subgraph and the external graph.

The difficulty we must face is that, in a complex network, the labels can exponentially explode for increasing values of q and  $|\Sigma|$  to  $|\Sigma|^q \gg |V|$  and, even worse, the number of simple paths can exponentially explode to  $|V|^q$ . For the simple reason that in complex networks the average separation is very low (the famous idea of *six degrees of separation*).

In this thesis we exploit the problem using randomized techniques and parallelization, which makes the problem suitable even for big network.

#### 1.3 Pratical applications

The problem can be applied to a lot of context. That is why it is very important to choose the right domains for the values of the  $V, E, L, \Sigma, q$ :

- $\bullet$  V are
- E represent the set of interactions, two vertices are connected if exists a relation among them.
- L and  $\Sigma$  are the category that partition V, note that if  $|\Sigma| = 1$  the labeling is useless.
- q should be low as  $N^{< q}(u)$  could be a large portion of G, (e.g. in Facebook for  $q \simeq 4$  we have  $N^{< q}(u) \simeq G$ ).

Furthermore, we have to choose G1 and G2, like ego networks or connected components.

## Basic tools

In this chapter we first give a definition of subgraphs similarity

#### 2.1 Similarity indices

**Definition 2.1.** Given two set A and B we define the **Jaccard index** as the ratio between the size of intersection and of the union between the two sets:

$$J(A,B) = \frac{|A \cap B|}{|A \cup B|} \tag{2.1}$$

**Definition 2.2.** Given two set A and B we define the **Bray-Curtis index** as:

$$BC(A,B) = \frac{2 \times |A \cap B|}{|A| + |B|} \tag{2.2}$$

We can easily extended the two previous definition to multiset:

**Definition 2.3.** Given two multiset  $A = (a_1, \ldots, a_n)$  and  $B = (b_1, \ldots, b_n)$  we define the **weighted Frequency Jaccard index** as:

$$FJ(A,B) = \frac{\sum_{i=1}^{n} \min(a_i, b_i)}{\sum_{i=1}^{n} \max(a_i, b_i)}$$
(2.3)

**Definition 2.4.** Given two multiset  $A = (a_1, \ldots, a_n)$  and  $B = (b_1, \ldots, b_n)$  we define the **Bray-Curtis index** on multiset as:

$$BC(A,B) = \frac{2 \times \sum_{i=1}^{n} \min(a_i, b_i)}{\sum_{i=1}^{n} a_i + b_i}$$
(2.4)

2.2 Documents similarity

- 2.3 Graphs similarity
- 2.4 Subgraphs similarity

...

- 2.5 Sketches
- 2.5.1 min-wise permutation
- 2.5.2 bottom-k sketches
- 2.6 Color Coding

# Computation of subgraph similarity

#### **Algorithm 1:** Bray-Curtis

**Input**: W a dictionary of strings

 $f_A[W], f_B[W]$ 

Output: BC(A, B) Bray-Curtis index

ı return M

## 3.1 Naive approach

Algorithm 2: preprocess: BRUTE-FORCE

**Input**: G = (V, E) undirected graph with q random colors.

**Output:**  $(f_A[x], f_B[x])$  dynamic programming table for color coding.

 $_1$  return M

```
Algorithm 3: preprocess: COLOR-CODING

Input: G = (V, E) undirected graph with q random colors.

Output: M = dynamic programming table for color coding.

1 parallel foreach u \in V do M_{1,u} = \langle \chi(u), 1 \rangle

2 for i \in \{2, 3, \dots, q\} do

3 parallel foreach u \in V do

4 foreach v \in N(u) do

5 foreach \langle C, f \rangle \in M_{i-1,v} such that \chi(u) \notin C do

6 f' \leftarrow M_{i,u} (C \cup \{\chi(u)\})

7 M_{i,u} \leftarrow \langle C \cup \{\chi(u)\}, f' + f \rangle

8 return M
```

#### 3.2 Efficient computation

Color coding

Colorful sampling

Frequency count

Frequency sampling

Estimating similarity indices

#### 3.3 Baseline algorithm

## Project development

In this chapter we describe

#### 4.1 Implementation choices

#### 4.2 Dataset

For the experiments we use two differents kind of dataset, a small one so we can easily brute-force the real indices and compare the relative error, and a big one in order to

**NetInf** This graph represents the flow of information on the web among blogs and news websites. The graph was computed by the *NetInf* approach, as part of the SNAP project [?], by tracking cascades of information diffusion to reconstruct "who copies who" relationships. Each node represents a blog or news website, and a website is connected to those who frequently copy their content. The graph contains 854 nodes and 3824 edges. We labelled websites according to their importance, using Amazon's Alexa ranking [?]: the labels correspond to respectively the websites ranked in the top 4%, the following 15%, the following 30%, and the remaining 51% (i.e.  $|\Sigma| = 4$ ).

Considered query: compute the similarity of two websites a and b or two sets of websites.

**IMDb** In this graph, taken from the *Internet Movie Database* [?], nodes correspond to movies, and there is a link between two movies if their casts share at least one actor. The graph contains  $1\,060\,209$  movies (nodes) and  $288\,008\,472$  edges. Each movie is labeled with one of  $|\Sigma| = 36$  genres.

Considered query: similarity of actors' ego networks. Given two actors a and b, let A and B be their ego-networks, i.e., the sets of nodes corresponding to movies in which respectively a and b starred. Compute the similarity of A and B.

### 4.3 Experimental results

We describe the experimental evaluation for our approach. Our computing platform is a machine with Intel(R) Xeon(R) CPU E5-2620 v3 at 2.40GHz, 24 virtual cores, 128 Gb RAM, running Ubuntu Linux version 4.4.0-22-generic. Code written in C++ and compiled with g++ version 5.4.1 with OpenMP.

# Appendix A

# Code snippets

All the code written for this thesis can be found in the personal GitHub page<sup>1</sup>

#### A.1 Color Coding

```
map < COLORSET, long long > M[Q][V];
void ColorCoding() {
  #pragma omp parallel for schedule(static, 1)
  for (int u = 0; u < N; u++)
    M[0][u][setBit(0, color[u])] = 1;
  for (int i = 1; i < q; i++) {
    #pragma omp parallel for schedule(static, 1)
    for (int u = 0; u < V; u++) {</pre>
      for (int v : G[u]) {
        for (auto d : M[i-1][v]) {
          COLORSET s = d.first;
          long long f = d.second;
          if ( !getBit(s, color[u]))
            M[i][u][setBit(s, color[u])] += f;
        }
      }
    }
  }
```

<sup>1</sup>https://github.com/GaspareG/ColorCoding

#### A.2 Colorful sampling

```
vector < int > randomPathTo(int u) {
  list<int> P;
  P.push_front(u);
  COLORSET D = getCompl(setBit(01, color[u]));
  for (int i = q - 1; i > 0; i--) {
    vector<ll> freq;
    for (int v : G[u]) freq.push_back(M[i][v][D]);
    disc_distr<int> dist(freq.begin(), freq.end());
    u = G[u][dist(eng)];
    P.push_front(u);
    D = clearBit(D, color[u]);
  }
// reverse(P.begin(), P.end());
  return vector < int > ( begin(P), end(P));
}
set < string > colorfulSample(vector < int > X, int r) {
  set < string > W;
  set < vector < int >> R;
  vector<ll> freqX;
  for (int x : X) freqX.push_back(M[q][x][getCompl(0)]);
  disc_distr<int> dist(freqX.begin(), freqX.end());
  while (R.size() < (size_t)r) {</pre>
    int u = X[dist(eng)];
    vector < int > P = randomPathTo(u);
    if (R.find(P) == R.end()) R.insert(P);
  }
  for (auto r : R) {
    // reverse(r.begin(), r.end());
    W.insert(L(r));
 return W;
}
```

#### A.3 Frequency count

```
map<string, ll> frequencyCount(set<string> W, multiset<int> X) {
  set < string > WR;
  for (string w : W) {
    reverse(w.begin(), w.end());
    WR.insert(w);
  }
  vector<tuple<int, string, COLORSET>> old;
  for (int x : X)
    if (isPrefix(WR, string(&label[x], 1)))
      old.push_back(make_tuple(x, string(&label[x], 1), setBit(011,
  for (int i = q - 1; i > 0; i--) {
    vector<tuple<int, string, COLORSET>> current;
    #pragma omp parallel for schedule(static, 1)
    for (int j = 0; j < (int)old.size(); j++) {</pre>
      auto o = old[j];
      int u = get < 0 > (o);
      string LP = get<1>(o);
      COLORSET CP = get <2>(o);
      for (int v : G[u]) {
        if (getBit(CP, color[v])) continue;
        COLORSET CPv = setBit(CP, color[v]);
        string LPv = LP + label[v];
        if (!isPrefix(WR, LPv)) continue;
        #pragma omp critical
        { current.push_back(make_tuple(v, LPv, CPv)); }
    }
    old = current;
  map < string, 11 > frequency;
  for (auto c : old) {
    string s = get < 1 > (c);
    reverse(s.begin(), s.end());
    frequency[s]++;
  }
  return frequency;
```

}

Code snippets

## A.4 Frequency sampling

```
map<pair<int, string>, ll> randomColorfulSamplePlus(vector<int> X,
  map<pair<int, string>, ll> W;
  set < vector < int >> R;
  vector<ll> freqX;
  freqX.clear();
  for (int x : X) freqX.push_back(M[q][x][getCompl(011)]);
  discrete_distribution < int > distribution (freq X . begin (), freq X . end (
  while (R.size() < (size_t)r) {</pre>
    int u = X[distribution(eng)];
    vector < int > P = randomPathTo(u);
    if (R.find(P) == R.end()) R.insert(P);
  }
  for (auto r : R) {
    reverse(r.begin(), r.end());
    W[make_pair(*r.begin(), L(r))]++;
  }
  return W;
}
```

## A.5 Similarity indices

```
double BCW(set<string> W,
           map<string, 11> freqA,
           map<string, 11> freqB) {
  11 \text{ num} = 011;
  11 \ den = 011;
  for (string x : W) {
    ll fax = freqA[x];
    11 fbx = freqB[x];
    num += 2 * min(fax, fbx);
    den += fax + fbx;
 return (double)num / (double)den;
}
double FJW(set<string> W, map<string, ll> freqA, map<string, ll> freq
           long long R) {
  11 \text{ num} = 011;
  for (string x : W) {
    ll fax = freqA[x];
    11 fbx = freqB[x];
    num += min(fax, fbx);
  return (double)num / (double)R;
}
```