Unsupervised Exceptional Attributed Sub-graph Mining in Urban Data
Anes Bendimerad, Marc Plantevit, Céline Robardet

To cite this version:

HAL Id: hal-01430622
https://hal.archives-ouvertes.fr/hal-01430622
Submitted on 10 Jan 2017

HAL is a multi-disciplinary open access archive for the deposit and dissemination of scientific research documents, whether they are published or not. The documents may come from teaching and research institutions in France or abroad, or from public or private research centers.

L’archive ouverte pluridisciplinaire HAL, est destinée au dépôt et à la diffusion de documents scientifiques de niveau recherche, publiés ou non, émanant des établissements d’enseignement et de recherche français ou étrangers, des laboratoires publics ou privés.
Unsupervised Exceptional Attributed Sub-graph Mining in Urban Data

Ahmed Anes Bendimerad*, Marc Plantevit†, Céline Robardet*

*INSA de Lyon, CNRS, LIRIS UMR5205, F-69621 France
†Université Lyon 1, CNRS, LIRIS UMR5205, F-69622 France
{firstname.name}@liris.cnrs.fr

Abstract—Geo-located social media provide a wealth of information that describes urban areas based on user descriptions and comments. Such data makes possible to identify meaningful city neighborhoods on the basis of the footprints left by a large and diverse population that uses this type of media. In this paper, we present some methods to exhibit the predominant activities and their associated urban areas to automatically describe a whole city. Based on a suitable attributed graph model, our approach identifies neighborhoods with homogeneous and exceptional characteristics. We introduce the novel problem of exceptional sub-graph mining in attributed graphs and propose a complete algorithm that takes benefits from new upper bounds and pruning properties. We also propose an approach to sample the space of exceptional sub-graphs within a given time-budget. Experiments performed on 10 real datasets are reported and demonstrate the relevancy and the limits of both approaches.

I. INTRODUCTION

In today’s increasingly global and interconnected world, people have opportunities to live abroad of their country, generally in urban areas. They face the challenge of making decisions about where to live, how to find appropriate areas to go out or a place to visit. Thanks to the current numerical development, numerous sources of collected data can help to make better decisions. Nevertheless, such geo-enabled social data must be processed with efficient methods to take into account the heterogeneity and the complexity of urban areas by the discovery of useful and understandable insights. Such questions have recently raised the interests of researchers such as discovering similar neighborhoods across several cities [5] or matching social attributes with geographic spaces [19].

Using social and urban data of a city (such as the ones provided by social networks as FOURSQUARE or GOOGLE PLACE), we aim to identify neighborhoods with homogeneous and exceptional characteristics: Areas are described by their associated characteristics that distinguish them from the rest of the city. To this end, we propose a suitable attributed graph model (as illustrated in Fig. 1) that results from the combination of social and urban data, and we achieve the task by applying a constraint-based graph pattern mining approach. The devised algorithm identifies connected sub-graphs associated to some characteristics that discriminate the subgraphs from the rest of the graph.

Attributed graph analysis has received much attention in the past decade. For example, [14] designed a method to find dense homogeneous sub-graphs, where vertices are described by categorical attributes and [7] proposes subspace clustering approach using numerical vertex attributes. However, all these works focus their attention on the similarity inside the sub-graphs, while underestimate exceptionality of the sub-graph characteristics with respect to the whole graph.

We design two algorithms to solve this problem. The first one is an exact algorithm that uses original and efficient upper bounds and some other techniques to reduce the search space. We also propose a method that reduces the number of output patterns by providing a concise summary of the complete result set while limiting the overlap between patterns. This summary represents the whole set of solutions, while being faster to compute. The second algorithm mines exceptional sub-graphs by sampling the space of patterns in a similar way as [2], [8], [17].

Our main contributions are manifold: First, we propose a new kind of graph analysis that exploits both of the contrasts of vertices attributes and the graph structure with a connectivity constraint. Second, to solve the defined problem, we present an efficient algorithm based on new upper bounds and pruning properties. We also propose a method to summarize the output set. Third, we design a probabilistic approach that samples the output space of patterns within a time budget specified by end-users. Forth, we provide a thorough empirical study that includes (1) a demonstration of the efficiency of the used pruning techniques, (2) the impact of the parameters and the input graph dimensions on the performance of the algorithms, and (3) the relevancy of the discovered results.

The rest of the paper is organized as follows. Section II formally introduces the problem. The proposed solutions are presented in Section III. Section IV reports on the empirical study on both synthetic and real-world datasets. Section V discusses related work and section VI concludes.
II. PROBLEM FORMULATION

Data describing geographic venues are numerous, ranging from census data to collaborative data produced through social-media platforms. To describe a city, nearby venues are grouped into small areas (geographers generally use tiles of 200 meters) over which venue characteristics are aggregated into count data. These areas are hereafter considered as the vertices \( V \) of a graph \( G = (V, E, C, D) \) whose edges \( E \) connect adjacent areas (that share a part of their borders), \( C = \{ c_i, i \in \mathbb{I} \} \) is a set of \( p \) categories and the vertices of \( V \) are described by \( D = \{ c_i(v) \in \mathbb{N}, \text{with } c_i \in C \text{ and } v \in V \} \), the counts of venues of each category in the area associated to each vertex.

The values of \( D \) can be aggregated over a set of vertices \( K \subseteq V \) and a set of categories \( L \subseteq C \): \( \text{sum}(L, K) = \sum_{v \in K} \sum_{c_i \in L} c_i(v) \). To simplify the notation, we use \( \text{sum}(K) \) to denote \( \text{sum}(C, K) \).

As an example, Fig. 1 presents a graph derived from the division of a city into 6 areas (from \( v_1 \) to \( v_6 \)). The area represented by \( v_1 \) is adjacent to the ones represented by \( v_2 \) and \( v_4 \), and consequently an edge connects \( v_1 \) to \( v_2 \) and another one \( v_1 \) to \( v_4 \). The number of venues of each category in a given area composed a vector associated to the corresponding vertex. The distribution of venue categories \( C = \{ \text{Health, Tourism, Store, Food} \} \) is detailed in Fig. 2. \( \text{sum}(\text{health}, \{ v_1 \}) = 1 \) as there is one venue with the category \text{health} in the area associated to \( v_1 \). We can also observe that \( \text{sum}(\{ \text{Health, Tourism, Store, Food} \}, \{ v_1 \}) = 22 \), and for the set \( K = \{ v_2, v_5 \} \), \( \text{sum}(K) = 49 \).

\[ \text{Fig. 2: Example of the distribution of venues in areas.} \]

Our objective is to identify neighborhoods whose characteristics distinguish them from the rest of the city. Therefore, we propose to discover connected sub-graphs associated to exceptional categories. A category is exceptional for a sub-graph if it is more frequent in its vertices than in the remaining of the graph. The scarcity of a category can also be a relevant descriptive element of a neighborhood. In Fig. 2, vertices \( v_2 \) and \( v_5 \) have a surplus on the category \text{Health} compared to the rest of the graph, while having a loss on category \text{Tourism}. Thus, we associate to a sub-graph a characteristic defined as a pair \( S = (S^+, S^-) \), with \( S^+ \) and \( S^- \) two disjoint subsets of \( C \). In order to assess the relevancy of the characteristics \( S \) with respect to the sub-graph induced by \( K \subseteq V \), noted \( G(K) \), we define the measure \( WRAcc(S, K) \), an adaptation of the weighted relative accuracy measure widely used in Subgroup Discovery [9].

A set of categories \( L \) is discriminant to \( G(K) \) if it is more or, on the contrary, less frequent in \( G(K) \) than in \( G \). This is evaluated by the gain function:

\[ \text{gain}(L, K) = \frac{\text{sum}(L, K)}{\text{sum}(K)} - \frac{\text{sum}(L, V)}{\text{sum}(V)} \]

The validity of a characteristic \( S = (S^+, S^-) \) with respect to \( G(K) \) is given by

\[ \text{valid}(S, K) = \bigwedge_{v \in K} \left( \bigwedge_{c_i \in S^+} \delta_{\text{gain}(c_i, v) > 0} \right) \land \left( \bigwedge_{c_i \in S^-} \delta_{\text{gain}(c_i, v) < 0} \right) \]

\( \text{valid}(S, K) \) means that each vertex \( v \in K \) has a positive gain for each category \( c_i \in S^+ \), and a negative gain for each category \( c_i \in S^- \). The quality of characteristic \( S \) can be globally measured by the numerical function \( A \):

\[ A(S, K) = \text{gain}(S^+, K) - \text{gain}(S^-, K) \]

However, a major drawback of the gain is that it is easy to obtain high value with highly specific characteristics [9], more precisely characteristics associated to a small set of vertices. Weighted relative accuracy makes a trade-off between generality and gain by considering the relative size of the sub-graph:

\[ WRAcc(S, K) = \left\{ \begin{array}{ll} A(S, K) \times \frac{|K|}{|V|} & \text{if } \text{valid}(S, K) \\ 0 & \text{otherwise} \end{array} \right. \]

The main differences with the \( WRAcc \) used in Subgroup Discovery [9] are (1) our adapted \( WRAcc \) considers both the positive and the negative contrasts in an unsupervised setting, (2) it takes into account the homogeneity of elements of \( K \), using the predicate \( \text{valid}(S, K) \). We can now define the pattern domain we consider:

**Definition 1** (Exceptional sub-graph), Given a graph \( G = (V, E, C, D) \) and two thresholds \( \sigma \) and \( \delta \), an exceptional sub-graph \( (S, K) \) is such that (1) \( |K| \geq \sigma \), (2) \( G(K) \) is connected, and (3) \( WRAcc(S, K) \geq \delta \). \( \mathcal{E}(G) \) denotes the whole set of exceptional sub-graphs in \( G \).

The computation of the complete set of exceptional sub-graphs requires to search into two combinatorial search spaces, with constraints that cannot be used according to the usual techniques of search space pruning. Thus, a naive approach cannot achieve this task for large graphs or large number of categories. In the following we propose several enumeration strategies that takes benefit from computation of tight upper-bounds on \( WRAcc(S, K) \).

III. COMPUTING EXCEPTIONAL SUB-GRAPHS

This section introduces three distinct approaches to extract exceptional sub-graphs. First, we present an exact algorithm that aims at discovering the complete set of exceptional sub-graphs. Second, we propose to only compute a summary of this set by controlling the overlap between patterns. Third, we devise a heuristic algorithm that samples the space of
computes the exceptional sub-graphs. This approach makes possible to obtain instant results and to successfully scales up to a large number of attributes.

A. The complete approach

To compute all exceptional sub-graphs, two search spaces have to be explored: The space of characteristics $S = (S^{+}, S^{-})$ is first traveled, and then, for each promising characteristic, the sub-graphs $G(K), K \subseteq V$, are considered.

1) Characteristic enumeration: We explore the space of characteristics recursively. We start from the empty characteristic $(S^{+}, S^{-}) = (\emptyset, \emptyset)$ and consider the candidate categories that can be used to expand $S: (X^{+}, X^{-}) = (C, C)$. $X^{+}$ contains the categories that can be added to $S^{+}$, and $X^{-}$ the ones that can be added to $S^{-}$.

However, if the size of the category set $C$ is large, it is unrealistic to explore all the combinations. To early discard unpromising candidates, we exploit two properties: The anti-monotony of the predicate valid and an upper-bound on the WRAcc measure.

The predicate valid is anti-monotone with respect to the inclusion of characteristics: Considering two characteristics $S_{1}, S_{2}$ such that $S_{1}^{+} \subseteq S_{2}^{+}$ and $S_{1}^{-} \subseteq S_{2}^{-}$, denoted $S_{1} \subseteq S_{2}$, we have $\forall K \subseteq V$, valid$(S_{2}, K) \Rightarrow valid(S_{1}, K)$. Let $V(S) = \{v \in V \mid valid(S, v)\}$, if $V(S)$ has a size smaller than $\sigma$, then it is unnecessary to explore the children of $S$. To quickly obtain a small-sized set $V(S)$, we expand $S^{+}$ with the candidate $c \in X^{+}$ that has the smallest number of vertices $v \in V$ with $gain(c, v) > 0$ (a.k.a. fail-first principle), and to expand $S^{-}$ with the candidate $c \in X^{-}$ that has similar property for $gain(c, v) < 0$ (lines 1 and 2 of Algorithm 1).

Let us now define an upper bound on the WRAcc measure. Let $\Omega(S, X)$ be the exploration tree rooted in $(S, X)$ that contains all the characteristics $T$ obtained from the characteristic $S$ extended with some categories of $X$. An upper bound $UB_{1}$ on WRAcc with respect to a connected component $CC \subseteq V(S)$ is defined as: $UB_{1}(S, X, CC) = \max_{S \in \Omega(S, X), v \in CC} A(S, v) \times |CC|/|V|$.

Property 1. $WRAcc(T, K) \leq UB_{1}(S, X, CC), \forall T \in \Omega(S, X)$ and $K \subseteq CC$

Proof. From the definition of the function $A$, we have $A(S, K) = \sum_{v \in K} \frac{\text{sum}(v)}{\text{sum}(K)} \times A(S, v)$. Thus, $A(S, K)$ is simply a weighted average of $A(S, v)$. Furthermore, as $A(S, v) \leq A(S^{*}, v)$ with $S^{*} = \text{argmax}_{S \in \Omega(S, X)} A(S, v)$, from the properties of the average, we can infer that $A(S, K) \leq A(S^{*}, v)$. Thus, we can conclude that $A(S, K) \leq A(S^{*}, v) \times \frac{|CC|}{|V|}$.

The computation of $UB_{1}$ can be done in $\Theta(|CC| \times |X|)$. Algorithm 1 presents the pseudo-code of ENERGETICS\(^{1}\) that computes the exceptional sub-graphs. Line 3 stops the algorithm when there is no more candidate. The loop in line 4 discards unpromising connected components using the aforementioned upper bound. The four recursive calls consider the four possible extensions of $S$ by a category $c$. The function SUB-GRAPH is detailed below.

Algorithm 1: ENERGETICS$(S, X, Y, R, \delta, \sigma)$

Input: $S = (S^{+}, S^{-})$ the current explored characteristic, $X = (X^{+}, X^{-})$ the candidate sets, $Y = \{v \in V \mid valid(S, v)\}$

Output: $R$ the result set under construction

1. Sort $c \in X^{+}$ in ascending order of $|\{v \in V \mid gain(c, v) > 0\}|$
2. Sort $c \in X^{-}$ in ascending order of $|\{v \in V \mid gain(c, v) < 0\}|$
3. if $X \neq (\emptyset, \emptyset)$ then
   for $CC \subseteq G(Y)$ do
      if $UB_{1}(S, X, CC) < \delta$ or $|CC| < \sigma$ then
         $Y \leftarrow Y \setminus CC$
      if $X^{+} \neq \emptyset$ then
         $c \leftarrow \text{pop}(X^{+})$
         $Y' \leftarrow \{v \in Y \mid valid((S^{+} \cup \{c\}, S^{-}))\}$
         if $Y' \neq \emptyset$ then
            ENERGETICS$((S^{+} \cup \{c\}, S^{-}), (X^{+} \setminus \{c\}, X^{-} \setminus \{c\}), Y', R, \delta, \sigma)$
            ENERGETICS$(S(X^{+} \setminus \{c\}, X^{-}, Y'), R, \delta, \sigma)$
         else
            $c \leftarrow \text{pop}(X^{-})$
            $Y' \leftarrow \{v \in Y \mid valid((S^{+}, S^{-} \cup \{c\}))\}$
            if $Y' \neq \emptyset$ then
               ENERGETICS$((S^{+}, S^{-} \cup \{c\}), (X^{+}, X^{-} \setminus \{c\}), Y', R, \delta, \sigma)$
               ENERGETICS$(S(X^{+}, X^{-} \setminus \{c\}), Y, R, \delta, \sigma)$

2) Sub-graph exploration: Once the characteristic $S$ is fixed, its associated set of vertices $V(S)$ has to be processed to identify subsets $K'$ that (1) have a size greater or equal to $\sigma$, (2) satisfy $WRAcc(S, K') \geq \delta$ and (3) $K'$ are connected. To compute them efficiently, we propose an approach that combines reverse search \cite{20} and upper-bounds. It is based on the loose anti-monotonicity of the constraint $A(S, K) \geq \tau$.

Property 2 (Loose anti-monotone). $A(S, K) \geq \tau$ is a loose anti-monotone constraint. This implies that if $A(S, K) \geq \tau$, then it exists $v \in K$ such that $A(S, K \setminus \{v\}) \geq \tau$.

Proof. Let $v^{\circ} = \text{arg min}_{u \in K} A(S, u)$ and $K' = K' \cup \{v^{\circ}\}$. By construction, $\forall u \in K, A(S, u) \geq A(S, v^{\circ})$. Thus, the weighted average $\sum_{u \in K'} \frac{\text{sum}(u)}{\text{sum}(K)} A(S, u)$ is to say $A(S, K') \geq A(S, v^{\circ})$. As $A(S, K') = \sum_{v \in K'} \frac{\text{sum}(v)}{\text{sum}(K')} A(S, v)$, we can derive that $A(S, K') = \frac{\text{sum}(K')}{\text{sum}(K)} A(S, K') + \frac{\text{sum}(v^{\circ})}{\text{sum}(K')} A(S, v^{\circ}) \leq A(S, K') = A(S, K \setminus \{v^{\circ}\})$.

To exploit this property, we enumerate the vertex sets $K$ in a way such that $A(S, K)$ decreases during the enumeration. Starting from a sub-graph $K = \emptyset$ and a set of candidates vertices $Y$, $K$ is expanded by adding vertices from $Y$ according to the following order.

Definition 2 (Vertex order $\succeq_{K}$). Let $K \subseteq V$ and $v_{i}, v_{j} \in V \setminus K$. We say that $v_{i} \succeq_{K} v_{j}$ if $A(S, K \cup \{v_{i}\}) \geq A(S, K \cup \{v_{j}\})$.

\footnotesize{\textbf{ENERGETICS: ENumERating G} ExcepTrIonal Connected Sub-graphs}
From the loose anti-monotonicity of $A(S, K)$, we can derive an upper bound of the WRAcc constraint.

**Definition 3 (UB2).** Let $K$ be the current set of vertices and $Y$ the set of candidates. The function $UB_2$ is defined as:

$$UB_2(S, K, Y) = A(S, K) \times \frac{|K \cup Y|}{|V|}$$

**Property 3.** For all $G(K')$ with $K' = K \cup K''$ and $K'' \subseteq Y$, we have

$$WRAcc(S, K') \leq UB_2(S, K, Y)$$

**Proof.** Let $K' = K \cup K''$. By generalizing the result of proof 2, we have $A(S, K') \leq A(S, K)$. Moreover, as $K' \subseteq K \cup Y$, we have $\frac{|K'|}{|V|} \leq \frac{|K \cup Y|}{|V|}$. By multiplying these results we can conclude the proof.

Therefore, if $UB_2(S, K, Y) < \delta$, we are sure that there is no set $K' \subseteq K \cup Y$ such that $WRAcc(S, K') \geq \delta$ and we can stop the enumeration.

We can derive a tighter upper-bound constructed recursively on the size of the set of vertices $Z \subseteq Y$ added to $K$ in the enumeration process. Considering the vertex order $\geq_K$ used during the enumeration, we have $A(S, K) \geq A(S, v)$, $\forall v \in Y$. Thus, the second term in the following weighted average $A(S, K \cup \{v\}) = \frac{\sum \{v\}}{\sum \{K\}} \cdot A(S, K) + \frac{\sum \{v\}}{\sum \{v\}} \cdot A(S, v)$ tends to decrease $A(S, K \cup \{v\})$. To tightly upper bound $A(S, K \cup \{v\})$, we have to give a high value to $\frac{\sum \{v\}}{\sum \{K\} + \sum \{v\}}$, while minimizing $\frac{\sum \{v\}}{\sum \{K\} + \sum \{v\}}$. This is achieved by $d_1 = \min_{v \in Y} \{A(S, v)\}$.

Then, by choosing $d_1 = \max_{v \in Y} A(S, v)$, the above expression is upper bounded. We just have defined $\overline{A}(S, K, Y, 1)$, the upper bound for $K$ extended by a single vertex from $Y$.

Let us now suppose there is an upper bound $\overline{A}(S, K, Y, i)$ of $A(S, K \cup Z)$ with $|Z| = i$. To define $\overline{A}(S, K, Y, i + 1)$ for $|Z| = i + 1$, we consider $A(S, K \cup Z \cup \{v\}) = \frac{\sum \{K\} + \sum \{Z\} + \sum \{v\}}{\sum \{S\} + \sum \{K\} + \sum \{Z\} + \sum \{v\}} \cdot A(S, K \cup Z) + \frac{\sum \{v\}}{\sum \{v\}} \cdot A(S, v)$.

This tends to decrease the weighted sum, thus $(1) \frac{\sum \{S\} + \sum \{K\} + \sum \{Z\}}{\sum \{S\} + \sum \{K\} + \sum \{Z\} + \sum \{v\}}$ has to be maximized and $(2) \frac{\sum \{S\} + \sum \{K\} + \sum \{Z\}}{\sum \{S\} + \sum \{K\} + \sum \{Z\} + \sum \{v\}}$ has to be minimized. Let $|Y| = t$ and $\{d_1, \ldots, d_t\} = \{\sum \{v\} \mid v \in Y\}$ such that $\forall i, j \in [1, t]$, $d_i \leq d_j$. (1) is obtained by $\frac{\sum \{S\} + \sum \{K\} + \sum \{Z\}}{\sum \{S\} + \sum \{K\} + \sum \{Z\} + \sum \{v\}} \cdot A(S, v)$, and (2) by $\frac{\sum \{S\} + \sum \{K\} + \sum \{Z\}}{\sum \{S\} + \sum \{K\} + \sum \{Z\} + \sum \{v\}} \cdot A(S, v)$. This is synthesized in the following definition.

**Definition 4 (UB3).** Let $(K, Y)$ be the current enumerated sub-graph induced by $K$, and $Y$ the candidate vertices such that $|Y| = t$. Let $\{d_1, \ldots, d_t\} = \{\sum \{v\} \mid v \in Y\}$ such that $\forall i, j \in [1, t]$, $d_i \leq d_j$, and $\{a_1, \ldots, a_t\} = \{A(S, v) \mid v \in Y\}$ such that $\forall i, j \in [1, t]$, $a_i \geq a_j$. The upper bound $UB_3$ is

$$UB_3(S, K, Y) = \max_{i \in [0, t]} \left( A(S, Y, i) \times \frac{|K| + i}{|V|} \right)$$

We tackle this problem in a similar way as [22] that proposes an approach to reduce the output set of maximal cliques. The idea consists in returning a concise subset of $E(G)$, denoted $E'(G)$, that represents the whole set of patterns. More precisely, for each pattern $(S, K_1) \in E(G)$ there is a pattern $(S, K_2) \in E'(G)$ such that $K_1$ is included in $K_2$ or $K_2$ covers the majority of $K_1$. This is formally defined below.

**Definition 5 (Coverage measure).** Let $K_1 \subseteq V$ and $K_2 \subseteq V$. The coverage measure $cov$ is defined by $cov(K_1, K_2) = \frac{|K_1 \cap K_2|}{|K_1|}$. This function measures how much $K_2$ covers $K_1$.**
Algorithm 2: SUB-GRAPH(S, validV, δ, σ, R)

Input: S, validV = \{v ∈ V | valid(S, v)\}
Output: R the set of result patterns under construction
1 for each maximal connected component CC ⊆ validV do
2     SUB-CC(S, CC, δ, σ)
3     SUB-CC(S, (K, Y), δ, σ)
Input: S, K the explored sub-graph, Y the set of candidate vertices
Output: R the set of result patterns under construction
4 if UB2(S, K, Y) < δ then
5     return False
6 if UB3(S, K, Y) < δ then
7     return False
8 existsPattern ← False
9 while |K ∪ Y| ≥ σ and Y ≠ ∅ do
10       ν∗ = argmaxν∈Y A(S, K ∪ \{u\})
11       Y ← Y \ {ν∗}
12       if SUB-CC(S, (K ∪ \{ν∗\}, Y), δ, σ) then
13           existsPattern ← True
14       else
15           Y ← \{v ∈ Y | A(S, v) > A(S, ν∗)\}
16       if WRAcc(S, K) ≥ δ and |K| ≥ σ then
17           existsPattern ← True
18       if S is connected then
19           R ← R ∪ \{(S, K)\}
20 return existsPattern

Hence, the problem we consider in this section is how to compute E′(G) ⊆ E(G) such that ∀(S, K) ∈ E(G), ∃(S, K′) ∈ E′(G) with cov(K, K′) ≥ mincov and E′(G) as small as possible.

The problem of finding the set E′(G) of minimum size is NP hard. Thus, we adopt an efficient heuristic approach that is based on the enumeration process and to use the coverage measure to prune large parts of the sub-graph search space. We use the following lower bound to stop the enumeration process of a candidate (K, Y), as it guarantees that all the patterns generated from it are covered by another already found pattern.

**Definition 6 (LB).** Let K be the current enumerated set of vertices, Y the set of candidate vertices, R a set of solutions (R ⊆ E(G), and (S, H) ∈ R. Let n and \(\overline{\text{cov}}\) be:
\[ n = \frac{|V|}{A(S, K)} - |K| \]
\[ \overline{\text{cov}}(K, Y, H) = \frac{|K \cap H| + \max(0, n - |Y \setminus H|)}{|K| + \max(n, |Y \setminus H|)} \]

The lower bound function LB is thus defined:
\[ LB(K, Y, R) = \max\limits_{(S, H) ∈ R} (\overline{\text{cov}}(K, Y, H)) \]

This lower bound is used to prune the exploration of \(\Gamma(K, Y)\) as explained below.

**Property 5.** For each candidate (K′, Y′) ∈ \(\Gamma(K, Y)\) such that WRAcc(S, K′) ≥ δ, there exists at least one pattern (S, H) ∈ R that satisfies cov(K′, H) ≥ LB(K, Y, R).

**Proof. (Sketch)** Considering that A(S, K′) ≤ A(S, K), we can derive a bound on the minimal number of vertices that have to be added to K to obtain WRAcc(S, K′) ≥ δ. This is the value n. If n ≤ |Y \ H|, we can prove that cov is lower bounded by \(\frac{|K|}{|K| + |Y \setminus H|}\). If n > |Y \ H|, there will be necessarily vertices in the intersection and cov is lower bounded by \(\frac{|K| + n - |Y \setminus H|}{|K| + n}\).

Consequently, the exploration of a sub-tree \(\Gamma(K, Y)\) is stopped if LB(K, Y, R) ≥ mincov, because all the candidates (K′, Y′) of \(\Gamma(K, Y)\) with WRAcc(S, K′) ≥ δ are covered by an already found solution.

**C. The exceptional sub-graph space sampling approach**

In practice, the end-user wants to obtain high quality patterns in a short amount of time, especially in interactive data mining processes. This can be achieved by using the solution presented in the previous section. However, this approach does not scale very well with the cardinality of the category set. To overcome this issue, we propose an approach that computes a sampling of the exceptional sub-graphs that respects the distribution of the WRAcc measure within a user-given time-budget.

We adapt the randomized pattern mining technique of [2] to exceptional sub-graphs discovery. This so-called Controlled Direct Pattern Sampling enables the user to specify a time budget and computes a set of high quality patterns whose size directly depends on the specified amount of time. The idea consists of sampling the patterns based on a probability distribution that rewards high quality patterns. In a first attempt, we proposed to first sample the characteristics and then derive the associated sub-graphs. But this strategy failed in computing patterns with high WRAcc values because the graph structure was neglected. Thus, we adopted the reverse approach that consists in randomly generating connected sub-graphs and then deriving the most relevant characteristic \(S^*(K)\) that fulfills the constraint WRAcc(S*(K), K) ≥ δ. We choose \(S^*(K)\) as the characteristic that contains all the characteristics \(S\) valid for \(K\: VS\) such that valid(S, K), \(S ≤ S^*(K)\). We can easily demonstrate that \(S^*(K)\) is the most relevant characteristic for \(K\), that is \(VS\) such that valid(S, K), \(WRAcc(S^*(K), K) ≥ WRAcc(S, K)\).

The sub-graph generation process is based on a random walk on a graph whose vertices are sub-graphs of G and edges (transitions) are chosen following a probability measure that favors high quality patterns \((S^*(K), K)\). The random walk starts from an empty set \((K = ∅)\) that is expanded by adding vertices in the neighborhood of K drawn randomly: (1) The random walk starts by drawing a first vertex. Considering the weight distribution \(w_1(v) = A(S^*(v), v)\), the probability distribution is computed by \(P(\{v\}) = \frac{w_1(\{v\})}{\sum_{v \in V} w_1(\{v\})}\). Thus, the more contrasted is the vertex, the greater is its probability.

After drawing this first vertex v, the current sub-graph is \(K = \{v\}\). (2) The random walk continues by considering the vertices in the neighborhood of the current sub-graph G(K). Let N(K) be the set of neighbors of K: \(N(K) = \{v ∈ V | ∃v' ∈ K, (v, v') ∈ E\}\). Then, the sub-graph generation is either stopped, or a vertex \(v ∈ N(K)\) is drawn and added to K.
### Algorithm 3: EXPRESS

Output: $K$ the generated sub-graph

```plaintext
1 begin
2 for $v \in V$ do
3     $w_1(v) \leftarrow WRAcc(S^*(v), v)$
4     draw $v \sim w_1(v)$
5     $K \leftarrow \{v\}$
6     continue $\leftarrow$ True
7 while continue do
8     // Calculate the probability of drawing $K \cup \{v\}$ for each neighbor $u$
9     for $v \in N(K)$ do
10        $K' \leftarrow K \cup \{v\}$
11        $w_2(K', K) \leftarrow WRAcc(S^*(K'), K') - WRAcc^s(K)$
12     // Calculate the probability of drawing $K$:
13     $w_2(K, K) \leftarrow WRAcc(S^*(K), K) - WRAcc^s(K)$
14     draw $K' \sim w_2(K', K)$
15     if $K' = K$ then
16        continue $\leftarrow$ False
17     else
18        $K \leftarrow K'$
19     return $K$
```

The choice of the action is taken randomly using a weighted distribution that favors the vertex (and possibly none) whose addition to $K$ leads to the largest increase of the WRAcc measure. Let $K' = K \cup \{v\}$. We define $WRAcc^s(K)$ as the lowest score that can be achieved by the current possible actions:

$$WRAcc^s(K) = \min(\min_{v \in N(K)} WRAcc(S^*(K'), K'), WRAcc(S^*(K), K))$$

From the weight $w_2(K', K)$, defined as $w_2(K', K) = WRAcc(S^*(K'), K') - WRAcc^s(K)$, we derive the probability to reach the connected sub-graph $G(K \cup \{v\})$ from $G(K)$ by $P(K \cup \{v\}) = \frac{\sum_{w_2(K \cup \{v\}, K') \in w_2(K, K)} w_2(K \cup \{v\}, K')}{\sum_{w_2(K \cup \{v\}, K')} w_2(K \cup \{v\}, K')}$. This distribution of probabilities favors transitions towards connected sub-graphs $K'$ with large $WRAcc(S^*(K'), K')$ value. The algorithm EXPRESS$^2$ is repetitively called until the specified execution time is consumed. For each generated sub-graph $K$, if $WRAcc(S^*(K), K) \geq \delta$, then the pattern $(S^*(K), K)$ is added to the output result set.

### IV. Experiments

In this section, we report on experimental results to illustrate the interest of the proposed approach. We start by describing the different real-world datasets we use, as well as the questions we aim to answer. Then, we provide a performance study and give some qualitative results. The implementation of the method is in Java and the experiments run on machines equipped with i7-2600 CPUs @ 3.40GHz, and 16GB main memory, running Ubuntu 12.04, and Java Version 1.6. The code and the data are available$^3$.

### A. Datasets and aims

We considered 10 real-world datasets whose characteristics are given in Table I. Eight of them come from [5] and depict Foursquare venues over 4 US and 4 EU important cities. The venues are described by a hierarchy$^4$. We consider the first level (10 attributes) in the first series of experiments and the second level (around 300 attributes) for the second ones. SF. Crimes data$^5$ is provided by a Kaggle challenge and describe the criminal activity in San Francisco. Finally, San Francisco C\&V is the combination – after normalization – of SF. Crimes and Foursquare data over San Francisco. Each city is divided into rectangular zones in such a way that each rectangle contains a minimal number of venues.

<table>
<thead>
<tr>
<th>dataset</th>
<th></th>
<th></th>
<th></th>
</tr>
</thead>
<tbody>
<tr>
<td>New York</td>
<td>292</td>
<td>647</td>
<td>10 (356)</td>
</tr>
<tr>
<td>Los Angeles</td>
<td>159</td>
<td>255</td>
<td>10 (325)</td>
</tr>
<tr>
<td>San Francisco</td>
<td>124</td>
<td>256</td>
<td>10 (328)</td>
</tr>
<tr>
<td>Washington</td>
<td>106</td>
<td>216</td>
<td>10 (316)</td>
</tr>
<tr>
<td>London</td>
<td>138</td>
<td>244</td>
<td>10 (319)</td>
</tr>
<tr>
<td>Paris</td>
<td>155</td>
<td>231</td>
<td>10 (315)</td>
</tr>
<tr>
<td>Rome</td>
<td>90</td>
<td>177</td>
<td>10 (293)</td>
</tr>
<tr>
<td>Barcelona</td>
<td>109</td>
<td>218</td>
<td>10 (394)</td>
</tr>
<tr>
<td>SF. Crimes</td>
<td>996</td>
<td>2172</td>
<td>39</td>
</tr>
<tr>
<td>S.F. C&amp;V</td>
<td>742</td>
<td>767</td>
<td>49 (328)</td>
</tr>
</tbody>
</table>

**TABLE I:** Description of the real-world datasets

ENERGETICS and EXPRESS$^2$ are evaluated regarding the following questions: What is the efficiency of ENERGETICS with regard to the graph characteristics that may affect its execution time? How effective are ENERGETICS’ pruning properties? Does ENERGETICS scale? Does EXPRESS$^2$ provide a good sample of Exceptional sub-graphs? What about the relevancy of Exceptional sub-graphs?

### B. Quantitative study

First, we devised a synthetic generator to evaluate how effective ENERGETICS’ upper bounds are, while varying the attributed graph properties $|V|$, $|E|$ and $|C|$. A baseline algorithm is obtained by deactivating the upper bounds and the pruning abilities of ENERGETICS. Thus, the baseline only pushes monotonic constraints. We also ran ENERGETICS using each upper-bound in turn. Results are given in Fig. 3. As the baseline only works with very small graphs, we report on ENERGETICS’ behavior with much more larger characteristics in the last column. In most of the cases, ENERGETICS algorithm outperforms the baseline and the versions with a single upper-bound by several orders of magnitude. The use of a single upper-bound takes in some cases more time than the baseline computation. This is due to the upper-bound verification that can be expensive while ineffective if not used in conjunction to other upper-bounds. There is no upper bound that outperforms the others. The efficiency of our algorithm is due to the conjunctive use of all the upper-bounds. We can observe that ENERGETICS scales well on the synthetic data according to the number of vertices and the number of edges while its execution time increases exponentially w.r.t. to the number of attributes.

$^2$https://developer.foursquare.com/categorytree
$^3$https://github.com/AnesBendimerad/Exceptional-Sub-graph-Mining
Fig. 3: Impact of the upper bounds on runtime (1st column), search space size – characteristics (2nd column), and subgraph (3rd column) w.r.t. the number of attributes (1st row), edges (2nd row), and vertices (3rd row). Broader parameter values are explored for ENERGETICS in the 4th column. Experiments carried out on synthetic data with $\delta = 0.015, \sigma = 1, \text{mincov} = 0.8$.

Fig. 4 reports the behavior (runtime, number of patterns, and number of explored subgraphs) of ENERGETICS on real-world datasets when varying the input parameters ($\text{mincov}$, $\delta$ and $\sigma$). Notice that ENERGETICS fails on the real-world data when the constraint on $\text{mincov}$ is desactivated. We can observe that the behavior of ENERGETICS remains unchanged when varying $\text{mincov}$. This is due to the fact that most of redundant subgraphs are strictly included in other ones and they are discarded with $\text{mincov} = 1$. Results for the two other parameters $\delta$ and $\sigma$ are as expected: The execution time increases when either the thresholds $\sigma$ and $\delta$ become less stringent. S.F. C&V is the dataset whose execution times are the most important. It confirms the fact that the number of attributes is the most influential data parameter.

We also studied the behavior of our algorithm w.r.t. the replication factor. For a replication factor equal to $n$, the attributed graphs are duplicated $n$ times, so the initial vertices are repeated $n$ times with the same attributes values and the same connections with the corresponding duplicated vertices. Fig. 5 presents the ratio of the execution time divided by the execution time on the original dataset. Thus, the ratio is equal to 1 for $n = 1$. We can observe that ENERGETICS performances does not degrade too much when the dataset size increases.

These first experiments demonstrate that ENERGETICS is only efficient for graphs whose number of attributes is small. EXPRESS has been designed especially to perform on graphs with hundreds of attributes, using a time budget to control the execution time and the number of computed patterns. To evaluate the ability of EXPRESS to compute exceptional sub-graphs of high WRAcc values, we report on Fig. 6 the distributions of the WRAcc measure on both the complete set of exceptional sub-graphs returned by ENERGETICS and the sample set provided by EXPRESS. Several time budgets are used and are all much lower than the execution time required by ENERGETICS. We can observe that the two distributions are similar and thus the sampling approach succeeds in fostering patterns with high WRAcc measure. Also, the higher the time budget, the better the distribution. Fig. 7 reports similar distributions for the real-world datasets with hundreds number of attributes for which an exhaustive search is not possible. The distributions are similar. Thus, EXPRESS makes it possible to discover high quality patterns within a time-budget.

C. Qualitative study

ENERGETICS was applied on Barcelona graph with 10 attributes. Fig. 8 (left) displays 5 patterns discovered. Pattern $P_1$ depicts neighborhoods with a high concentration of venues of type Outdoors & Recreation. Most of these zones are near the sea with the Olympic harbor and the main beaches of Barcelona. This pattern overlaps with two other patterns, $P_3$ and $P_4$. Notice that the ICDM’16 conference venue is covered by both $P_1$ and $P_3$. The attendees can take benefits of the high concentration of Outdoors & Recreation places as well as the nightlife spots and the food places. However, the concentration of shops in this areas is lower than in the rest of the city. Pattern $P_2$ depicts zones with a high concentration of places...
Fig. 4: Behavior of ENERGETICS (runtime in 1st column, #patterns in 2nd column and, #explored subgraphs in 3rd column) according to mincov (1st row), σ (2nd row) and δ (3rd row) for the 10 real-world datasets (δ = 0.001, σ = 1, mincov = 0.8).

Fig. 5: Runtime ratio w.r.t. the replication factor for real world datasets (δ = 0.001 except SF Crimes and SF C&V (0.007 and 0.03), σ = 1, mincov = 0.8).

of the type Outdoors & Recreation or College & Universities while the concentration of food places is low. This pattern is sensible since it contains the famous camp nou stadium as well as several universities and colleges.

We also applied EXPRESS on Barcelona graph with 304 attributes. Some patterns are reported in Fig. 8 (right). Pattern P₆ identifies an area with a high concentration of metro stations and state and municipalities as well as Home (private) in the North of the city. P₇ depicts an area with a high concentration of bars and typical Mediterranean restaurants while P₈ highlights areas with a larger variety of restaurants (e.g., Salad Place, Asian and Italian Restaurants). P₉ contains a high concentration of hotels and bars. Notice that P₇, P₈ and P₉ are close to ICDM’16 conference venue which attendees could take advantage of these patterns to prepare their venue according to their preferences.

Besides, we mined exceptional sub-graphs on the different cities. In most of them (e.g., Barcelona, Paris, Rome, Los Angeles, London), the nightlife spots are mainly located in the city center. For New York, they are in the south of Manhattan and the west of Brooklyn. The higher concentration of outdoor & recreation places is surrounding for London. For seaside towns, they are concentrated on the coasts.

V. RELATED WORK

Several approaches have been designed to discover new insights in vertex attributed graphs. The pioneering work of Moser et al. [14] presents a method to mine dense homogeneous sub-graphs, i.e., sub-graphs whose vertices share a large set of attributes. Similar to that work, Günnelmann et al. [7] introduce a method based on subspace clustering and dense sub-graph mining to extract non redundant sub-graphs that are homogeneous with respect to the vertex attributes. Silva et al. [18] extract pairs made of a dense sub-graph and a Boolean attribute set such that the Boolean attributes are strongly associated with the dense sub-graphs. In [16], the authors propose to mine the graph topology of a large attributed graph by finding regularities among numerical vertex descriptors. The main objective of all these approaches is to find regularities instead of peculiarities within a large graph,
Fig. 6: Distributions of the patterns from ENERGETICS and EXPRESS with different time budgets ($\delta = 0.001$).

whereas Exceptional Sub-graph Mining computes sub-graphs with their distinguishing characteristics.

Interestingly, a recent work [1] proposes to mine descriptions of communities from vertex attributes, with a Subgroup Discovery approach. In this supervised setting, each community is treated as a target that can be assessed by well-established measures, as the WRAcc measure used in this paper. More generally, Subgroup Discovery [9], [15] aims to find descriptions of subpopulations for which the distribution of a predefined target value is significantly different from the distribution in the whole data. When there are multiple targets, Subgroup Discovery consists in finding descriptions that significantly change the joint distribution of the target attributes – a task that has been introduced as Exceptional Model Mining [10], [3]. A variety of measures of changes have been used: pairwise correlation and entropy measures [10], Bayesian networks [4], Kullback-Leibler Divergence and encoding difference based on Minimum Description Length [21]. The common point to all these approaches is that the combination of large target space and non-monotonic measures leads to the use of heuristic search methods, i.e., beam search. Furthermore, these methods are supervised since the target attributes are given. The algorithms proposed in this paper extend many of these results to the unsupervised setting.

Motivated by both scalability issue and user interaction needs, sampling the output space of patterns has received much attention in the past decade. Many approaches have been proposed in the context of transactional data [2], [13], [12], [6], [11]. Some researchers have tackled the problem of sampling the output space of frequent sub-graphs in a collection of graphs [8], [17]. These methods are based on random walks. In particular, [17] aims at returning the top $k$ frequent graphs of a specified size. The problem we tackle is different on several points: We consider a single graph and a discriminative measure instead of a frequency measure. Beside, these methods aim at sampling frequent sub-graphs while we address the problem of exceptional attributed sub-
in the city areas and highlight some specific phenomena. Other quality measures can be extended to take into account other graph topological properties (e.g., diameter). Our approach is based on a random walk over sub-graphs that fosters patterns with a high WRAcc measure.

VI. CONCLUSION

We introduced the exceptional sub-graph mining problem to discover homogenous areas that differ from the rest of the city. We defined an efficient algorithm that computes the complete set of exceptional sub-graphs by taking advantage of several tight upper bounds and other pruning properties. We also illustrated the relevancy of the discovered patterns thanks to a qualitative analysis. Our proposal can be extended to take into account other graph topological properties (e.g., diameter). Other quality measures can be investigated to assess the quality of the characteristics in the city areas and highlight some specific phenomena.

ACKNOWLEDGMENT

This work was supported in part by the Group Image Mining (GIM) which joins researchers of THALES Group and LIRIS Lab. We thank especially Jérôme Kodjabachian and Bertrand Duqueroie of AS&BSIM Lab. of THALES Group. This work is also partially supported by the EU FP7-PEOPLE-2013-IAPP project GRAISearch and the CNRS project PEPS-2015-Préflute.

REFERENCES