Entity-Linking via Graph-Distance Minimization

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Entity-linking is a natural-language–processing task that consists in identifying the entities mentioned in a piece of text, linking each to an appropriate item in some knowledge base; when the knowledge base is Wikipedia, the problem comes to be known as *wikification* (in this case, items are wikipedia articles). One instance of entity-linking can be formalized as an optimization problem on the underlying concept graph, where the quantity to be optimized is the average distance between chosen items. Inspired by this application, we define a new graph problem which is a natural variant of the Maximum Capacity Representative Set. We prove that our problem is NP-hard for general graphs; nonetheless, under some restrictive assumptions, it turns out to be solvable in linear time. For the general case, we propose two heuristics: one tries to enforce the above assumptions and another one is based on the notion of hitting distance; we show experimentally how these approaches perform with respect to some baselines on a real-world dataset.

1 Introduction

Wikipedia [1] is a free, collaborative, hypertextual encyclopedia that aims at collecting articles on different (virtually, all) branches of knowledge. The usage of wikipedia for automatically tagging documents is a well-known methodology, that includes in particular a task called *wikification* [13]. Wikification is a special instance of *entity-linking*: a textual document is given and within the document various fragments are identified (either manually or automatically) as being *(named) entities* (e.g., names of people, brands, places...); the purpose of entity-linking is assigning a specific reference (a wikipedia article, in the case of wikification) as a tag to each entity in the document.

Entity-linking happens typically in two stages: in a first phase, every entity is assigned to a set of items, e.g., wikipedia articles (the *candidate nodes* for that entity); then a second phase consists in selecting a single node for each entity, from within the set of candidates. The latter task, called *candidate selection*, is the topic on which this paper focuses.

To provide a concrete example, suppose that the target document contains the entity "jaguar" and the entity "jungle". Entity "jaguar" is assigned to a set of candidates that contains (among others) both the wikipedia article about the feline living in America and the one about the Jaguar car producer. On the other hand, "jungle" is assigned to the article about tropical forests and to the one about the electronic music genre. Actually, there are more than 30 candidates for "jaguar", and more about 20 for "jungle".

In this paper, we study an instance of the candidate selection problem in which the selection takes place based on some cost function that depends on the average distance between the selected candidates, where the distance is measured on the wikipedia graph[2]: the rationale should be clear enough—concepts

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[1] http://en.wikipedia.org/

[2] The undirected graph whose vertices are the wikipedia articles and whose edges represent hyperlinks between them.

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appearing in the same text are related, and so we should choose, among the possible candidates for each entity, those that are more closely related to one another.

Getting back to the example above, there is an edge connecting "jaguar" the feline with "jungle" the tropical forest, whereas the distance between, say, the feline and the music genre is much larger.

The approach we assume here highlights the *collective* nature of the entity-linking problem, as mentioned already in [10]: accuracy of the selection can be improved by a global (rather than local) optimization of the choices. As [10] observes, however, trying to optimize all-pair compatibility is a computationally difficult problem.

In this paper, we prove that the problem itself, even in the simple instance we take into consideration, is NP-hard; however, it becomes efficiently solvable under some special assumptions. We prove that, although these assumptions fail to hold in real-world scenarios, we can still provide heuristics to solve real instances.

We test our proposals on a real-world dataset showing that one of our heuristics is very effective, actually more effective than other methods previously proposed in the literature, and more than a simple greedy approach using the same cost function adopted here.

2 Related Work

Named-entity linking (NEL)- also referred to as *named entity disambiguation* grounds mentions of entities in text (*surface forms*) into some knowledge base (e.g. Wikipedia, Freebase). Early approaches to NEL [13] make use of measures derived from the frequency of the keywords to be linked in the text and in different Wikipedia pages. These include *tf-idf*, χ^2 and *keyphraseness*, which stands for a measure of how much a certain word is used in Wikipedia links in relation to its frequency in general text. Cucerzan [7] employed the context in which words appears and Wikipedia page categories in order to create a richer representation of the input text and candidate entities. These approaches were extended by Milne and Witten [14] who combined commonness (i.e., prior probability) of an entity with its relatedness to the surrounding context using machine learning. Further, Bunescu [4] employed a *disambiguation* kernel which uses the hierarchy of classes in Wikipedia along with its word contents to derive a finer-grained similarity measure between the candidate text and its context with the potential named entities to link to. In this paper we will make use of Kulkarni et al.'s dataset [11]. They propose a general collective disambiguation approach, under the premise that coherent documents refer to entities from one or a few related topics. They introduce formulations that account for the trade-off between local spot-to-entity compatibility and measures of global coherence between entities. More recently, Han et al. [10] propose a graph-based representation which exploits the global interdependence of different linking decisions. The algorithm infers jointly the disambiguated named mentions by exploiting the graph.

It is worth to remark that NEL is a task somehow similar to Word Sense Disambiguation (determining the right sense of a word given its context) in which the role of the knowledge base is played by Wordnet [8]. WSD is a problem that has been extensively studied and its explicitly connection with NEL was made by Hachey et al [9]. WSD has been an area of intense research in the past, so we will review here the approaches that are directly relevant to our work. Graph-based approaches to word sense disambiguation are pervasive and yield state of the art performance [15]; however, its use for NEL has been restricted to ranking candidate named entities with different flavors of centrality measures, such as in-degree or PageRank [9].

Mihalcea [12] introduced an unsupervised method for disambiguating the senses of words using random walks on graphs that encode the dependencies between word senses.

Navigli and Lapata [18, 16, 17] present subsequent approaches to WSD using graph connectivity metrics, in which nodes are ranked with respect to their local *importance*, which is regarded using centrality measures like in-degree, centrality, PageRank or HITS, among others.

Importantly, even if the experimental section of this paper deals with a NEL dataset exclusively, the theoretical findings could be equally applied to WSD-style problems. Our *greedy* algorithm is an adaptation of Navigli and Velardi's Structural Semantic Interconnections algorithms for WSD [18, 6]. The original algorithm receives an ordered list of words to disambiguate. The procedure first selects the *unambiguous* words from the set (the ones with only one synset), and then for every ambiguous word, it iteratively selects the sense that is *closer* to the sense of disambiguated words, and adds the word to the unambiguous set. This works in the case that a sufficiently connected amount of words is unambiguous; this is not the case in NEL and in our experimental set-up, where there could potentially exists hundreds of candidates for a particular piece of text.

3 Problem statement and NP-completeness

In this section we will introduce the general formal definition of the problem, in the formulation we decided to take into consideration. We will make use of the classical graph notation: in particular, given an undirected graph $G = (V,E)$, we will denote with $G[W]$ the graph induced by the vertices in W, and with $d(u,v)$ the distance between the nodes u and v, that is, the number of edges in the shortest path from u to v (or the sum of the weights of the lightest path, if G is weighted).

If G is a graph and e is an edge of G, $G-e$ is the graph obtained by removing e from G; we say that e is a *bridge* if the number of connected components of $G-e$ is larger than that of G. A connected bridgeless graph is called *biconnected*; a maximal set of vertices of G inducing a biconnected subgraph is called a *biconnected component* of G.

We call our main problem the *Minimum Distance Representative*, in short MINDR, and we define it as follows. Given an undirected graph $G = (V,E)$ (possibly weighted) and k subsets of its set of vertices, $X_1,\ldots,X_k \subseteq V$, a feasible solution for MINDR is a sequence of vertices of G, $x_1,\ldots,x_k$, such that for any i, with $1 \leq i \leq k$, $x_i \in X_i$ (i.e., the solution contains exactly one element from every set, possibly with repetitions).

Given the instance $G,\{X_1,\ldots,X_k\}$, the measure (the *distance cost*) of a solution S, $x_1,\ldots,x_k$, is $f(S) = \sum_{i=1}^{k}\sum_{j=1}^{k} d(x_i,x_j)$. The goal is finding the solution of minimum distance cost, i.e., a feasible solution S such that $f(S)$ is minimum.

We call the restriction of this problem, in which the sets of vertices in input $\{X_1,\ldots,X_k\}$ are disjoint, MINDIR(Minimum Independent Distance Representative). In this case, for the sake of simplicity, we will refer to a solution as the multiset composed by its elements.[3]

3.1 NP-completeness of MINDR

The MINDIR problem seems to be similar and related to the so-called Maximum Capacity Representatives [5], in short MAXCRS. The Maximum Capacity Representatives problem is defined as follows: given some disjoint sets $X_1,\ldots,X_m$ and for any $i \neq j$, $x \in X_i$, and $y \in X_j$, a nonnegative capacity $c(x,y)$, a

[3] We shall make free use of multiset membership, intersection and union with their standard meaning: in particular, if A and B are multisets with multiplicity function a and b, respectively, the multiplicity functions of $A \cup B$ and $A \cap B$ are $x \mapsto \max(a(x),b(x))$ and $x \mapsto \min(a(x),b(x))$, respectively.

solution is a set $S = \{x_1, \ldots x_m\}$, such that, for any i, $x_i \in X_i$; such a solution is called *system of representatives*. The measure of a solution is the capacity of the system of representatives, that is $\sum_{x \in S} \sum_{y \in S} c(x,y)$, and the MAXCRS problem aims at *maximizing* it. The MAXCRS problem was introduced by [1], who showed that it is NP-complete and gave some non-approximability results. Successively, in [19], tight inapproximability results for the problem were presented.

The MINDIR problem differs from MAXCRS just for in the sense that we are dealing with distances instead of capacities, and therefore we ask for a minimum instead of a maximum. Nonetheless the following Lemma, whose proof is given in Appendix A, shows that also MINDIR problem is NP-complete.

Lemma 1. *The* MINDIR *(hence,* MINDR*) problem is NP-complete.*

4 The decomposable case

In this section we study the MINDR problem under some restrictive hypothesis and we will show that in this case a linear exact algorithm exists.

Even if it may seem that these hypothesis are too strong to make the algorithm useful in practice, in the next section we will use our algorithm to design an effective heuristic for the general problem. In particular, we assume that the graph G (possibly weighted) is such that:

- any set X_i induces a connected subgraph on G, i.e., $G[X_i]$ is connected,
- for any $i \neq j$, for any $x \in X_i$ and $y \in X_j$, x and y do not belong to the same biconnected component.

The problem, under these further restrictions, will be called *decomposable* MINDR. Note that the second condition implies that a decomposable MINDR is in fact an instance of MINDIR, because it implies that no two sets can have nonempty intersection.

Let us consider an instance $(G, \{X_1, \ldots X_k\})$ of decomposable MINDR problem on a graph $G = (V,E)$.

An edge $e = (x,y) \in E$ is called *useful* if it is a bridge, x and y do not belong to the same set X_i, and there are at least two indices i and j such that X_i and X_j are in different components of $G - e$ (since e is a bridge, the graph obtained removing the edge e from G is no more connected).

4.1 Decomposing the problem

The main trick that allows to obtain a linear-time solution for the decomposable case is that we can actually decompose the problem (hence the name) through useful edges. First observe that, trivially:

Remark 1. *Let $e = (x,y)$ be a useful edge and let Z_x and Z_y be the two connected components of $G - e$ containing x and y, respectively. In G, all paths from any $x' \in Z_x$ to any $y' \in Z_y$ must contain e.*

Moreover:

Remark 2. *Let $e = (x,y)$ be a useful edge. There cannot be an index i such that X_i has a nonempty intersection with both components of $G - e$.*

In fact, assume by contradiction that one such X_i exists, and let $u, w \in X_i$ be two vertices living in the two different components of $G - e$: since $G[X_i]$ is connected, there must be a path connecting u and w and made only of elements of X_i; because of Remark 1, this path passes through e, but this would imply that $x, y \in X_i$, in contrast with the definition of useful edge.

Armed with the previous observations, we can give the following further definitions. Let Y_x (respectively, Y_y) be the set of sets X_i such that $X_i \subseteq Z_x$ (respectively, $X_i \subseteq Z_y$); we denote the sets of nodes in Y_x and Y_y by $V(Y_x) \subseteq Z_x$ and $V(Y_y) \subseteq Z_y$, respectively.

By virtue of Remark 1, all the paths in G from any $x' \in V(Y_x)$ to any $y' \in V(Y_y)$ pass through e. This implies also that there is no simple cycle in the graph including both $x' \in V(Y_x)$ and $y' \in V(Y_y)$.

Given a solution S for MINDIR$(G, \{X_1, \ldots, X_k\})$, and a useful edge (x,y), we have:

$$\begin{aligned}\sum_{x_i,x_j \in S} d(x_i,x_j) \quad = \quad & \sum_{x_i,x_j \in S \cap V(Y_x)} d(x_i,x_j) + \sum_{x_i,x_j \in S \cap V(Y_y)} d(x_i,x_j) + \\ & 2 \sum_{x_i \in S \cap V(Y_x), x_j \in S \cap V(Y_y)} (d(x_i,x) + d(x,y) + d(y,x_j)) .\end{aligned}$$

Indeed all the shortest paths from any $x_i \in S \cap V(Y_x)$ to any $x_j \in S \cap V(Y_y)$ pass through the useful edge (x,y) by Remark 1. Moreover, since the sets $X_1, \ldots, X_k$ are disjoint, we have that $|S \cap V(Y_x)| = |Y_x|$ and $|S \cap V(Y_y)| = |Y_y|$, that is, a solution has exactly one element for each set in Y_x (respectively, Y_y). Hence we can rewrite the last summand of the above equation as follows:

$$\begin{aligned}\sum_{x_i \in S \cap V(Y_x), x_j \in S \cap V(Y_y)} (d(x_i,x) + d(y,x_j) + d(x,y)) \quad = \quad & |Y_y| \cdot \sum_{x_i \in S \cap V(Y_x)} d(x_i,x) + \\ & |Y_x| \cdot \sum_{x_j \in S \cap V(Y_y)} d(y,x_j) + \\ & |Y_x| \cdot |Y_y| \cdot d(x,y).\end{aligned}$$

By combining the two equations, we can conclude that finding a solution for MINDIR$(G, \{X_1, \ldots, X_k\})$ can be decomposed into the following two subproblems:

1. finding S_x minimizing $\sum_{x_i,x_j \in S \cap V(Y_x)} d(x_i,x_j) + 2 \sum_{x_i \in S \cap V(Y_x)} |Y_y| d(x_i,x)$ in the instance $(G[Z_x], Y_x)$;
2. finding S_y minimizing $\sum_{x_i,x_j \in S \cap V(Y_y)} d(x_i,x_j) + 2 \sum_{x_j \in S \cap V(Y_y)} |Y_x| d(y,x_j)$ in the instance $(G[Z_y], Y_y)$.

Note that both instances are smaller than the original one because of the definition of a useful edge. The idea of our algorithm generalizes this principle; note that the new objective function we must take into consideration is slightly more complex than the original one: in fact, besides the usual all-pair–distance cost there is a further summand that is a weighted sum of distances from some fixed nodes (such as x for the instance $G[Z_x], Y_x$ and y for the instance $G[Z_y], Y_y$).

We hence define an extension of the MINDR problem, that we call EXTMINDR (for *Extended Minimum Distance Representatives*). In this problem, we are given:

- an undirected graph $G = (V,E)$ (possibly weighted)
- k subsets of its set of vertices, $X_1, \ldots, X_k \subseteq V$
- a multiset B of vertices, each $x \in B$ endowed with a weight $b(x)$.

A feasible solution for the EXTMINDR is a multiset $S = \{x_1, \ldots, x_k\}$ of vertices of G, such that for any i, with $1 \leq i \leq k$, $S \cap X_i \neq \emptyset$ (i.e., the set contains at least one element from every set). Its cost is

$$f(S) = \sum_{i=1}^{h} \sum_{j=1}^{k} d(x_i,x_j) + \sum_{i=1}^{k} \sum_{z \in B} b(z) d(x_i,z).$$

The goal is finding the solution of minimum cost, i.e., a feasible solution S such that $f(S)$ is minimum. The original version of the problem is obtained by letting $B = \emptyset$.

We are now ready to formalize our decomposition through the following Theorem, whose proof is given in Appendix B.

Theorem 1. *Let us be given a decomposable* EXTMINDR *instance* $(G, \{X_1, \ldots, X_k\}, B, b)$ *and a useful edge* $e = (t_0, t_1)$. *For every* $s \in \{0,1\}$, *let* Z_s *be the connected component of* $G - e$ *containing* t_s, Y_s *be the set of sets* X_i *such that* $X_i \subseteq Z_s$ *and* $V(Y_s)$ *be the union of those* X_i*'s. Let also* B_s *be the intersection of* B *with* Z_s. *Define a new instance* $I_s = (T[Z_s], \{X_i, i \in Y_s\}, B_s \cup \{t_s\}, b_s)$ *where*

$$b_s(t_s) = 2|Y_{1-s}| + \sum_{z \in B_{1-s}} b(z) \text{ and } b_s(z) = b(z), \text{ for any } z \in B.$$

Then the cost $f(S)$ *of an optimal solution* S *of the original problem is equal to*

$$f(S_0) + f(S_1) + 2|Y_0||Y_1|d(t_0,t_1) + \sum_{s \in \{0,1\}} \left(|S \cap V(Y_s)| \cdot \sum_{z \in B \cap Z_{1-s}} b(z) d(t_s, z) \right)$$

where S_s *is an optimal solution for the instance* I_s.

For completeness, we need to consider the base case of an instance with just one set $G, \{X_1\}, B, b$: the solution in this case is just one node $x \in X_1$ and the objective function to be minimized is simply $\sum_{z \in B} d(x,z)b(z)$. The optimal solution can be found by performing a BFS from every $z_j \in B$ (in increasing order of j), maintaining for each node $y \in X_1$, $g(y) = \sum_{z_t \in B, t<j} d(x, z_t)b(z_t)$, and picking the node having maximum final $g(y)$. This process takes $O(|B| \cdot |E(G[X_1])|)$. It is worth observing that in our case the size of the multiset B is always bounded by k. Moreover since $\sum_{i=1}^{k} |E(G[X_i])| \leq |E(G)| = m$, the overall complexity for all these base cases is bounded by $O(k \cdot m)$.

4.2 Finding useful edges

For every instance with more than one set, given an useful edge e the creation of the subproblems as described above is linear, so we are left with the issue of finding useful edges. This task can be seen as a variant of the standard depth-first search of bridges, as shown in Algorithm 2 and 3, in Appendix C. Recall that bridges can be found by performing a standard DFS that numbers the nodes as they are found (using the global counter visited, and keeping the DFS numbers in the array dfs); every visit returns the index of the least ancestor reachable through a back edge while visiting the DFS-subtree rooted at the node where the visit starts from. Every time a DFS returns a value that is larger than the number of the node currently being visited, we have found a bridge.

The variant consists in returning not just the index of the least ancestor reachable, but also the set of indices i that are found while visiting the subtree. If the set of indices and its complement are both different from $\emptyset$ then the bridge is useful: at this point, a "rapid ascent" is performed to get out of the recursive procedure.

4.3 The final algorithm

Combining the observations above, we can conclude that the overall complexity of the algorithm is $O(k \cdot m)$. The algorithm is presented in Algorithm 1.

5 The general case

As we observed at the beginning, the MINDR problem is NP-complete in general, although the decomposable version turns out to be linear. We want to discuss how we can deal with a general instance of the problem. To start with, let us consider a general connected MINDR instance, that is:

Algorithm 1: DECOMPOSABLEMINDR

Input: A graph $G = (V,E)$, $X_1 \ldots, X_k \subseteq V$, a weighted multiset B of nodes in V, where each element in B has a weight b. $G[X_i]$ is connected for every i and moreover for all $i \neq j$ and $x \in X_i$, $y \in X_j$, the two vertices x and y do not belong to the same biconnected component of G.

Output: A solution $S = \{x_1, \ldots, x_k\}$ such that for any i, with $1 \leq i \leq k$, $x_i \in X_i$, minimizing $\sum_{i=1}^{h} \sum_{j=1}^{k} d(x_i, x_j) + \sum_{i=1}^{k} \sum_{z \in B} b(z) d(x_i, z)$

Find a useful edge $e = (x,y)$, if it exists, using Algorithm 2

if *the useful edge does not exist* **then**

  if $k \neq 1$ **then**

    Fail!

  end

  Output the element $x_1 \in X_1$ minimizing $\sum_{z \in B} b(z) d(x_1, z)$

else

  Let Z_x (respectively Z_y) be the connected component of $T - e$ containing x (respectively y) .

  Let Y_x (respectively Y_y) be the indices i such that $X_i \subseteq Y_x$ ($X_i \subseteq Y_y$, respectively)

  $B' \leftarrow B \cup \{x\}$ (multiset union) with $b(x) = 2|Y_y| + \sum_{z \in B \cap Z_y} b(z)$

  $B' \leftarrow B' \cap Z_x$ (multiset intersection)

  $S' \leftarrow$ DECOMPOSABLEMINDR$(T[Z_x], Y_x, B')$

  $B'' \leftarrow B \cup \{x\}$ (multiset union) with $b(y) = 2|Y_x| + \sum_{z \in B \cap Z_x} b(z)$

  $B'' \leftarrow B'' \cap Z_y$ (multiset intersection)

  $S'' \leftarrow$ DECOMPOSABLEMINDR$(T[Z_y], Y_y, B'')$

  return $S' \cup S''$

end

- a connected undirected (possibly weighted) graph $G = (V,E)$,
- k subsets of its set of vertices, $X_1, \ldots, X_k \subseteq V$,

with the additional assumption that $G[X_i]$ is connected for every i. Recall that a feasible solution is a sequence S of vertices of G, $x_1, \ldots, x_k$, such that for any i, with $1 \leq i \leq k$, we have $x_i \in X_i$; its (distance) cost is $f(S) = \sum_{i=1}^{k} \sum_{j=1}^{k} d(x_i, x_j)$.

We shall discuss two heuristics to approach this problem: the first is related to Algorithm 1 in that it tries to modify the problem to make it into a decomposable one, whereas the second is based on the notion of hitting distance.

Before describing the two heuristics, let us briefly explain the rationale behind the additional assumption (i.e., that every $G[X_i]$ be connected). In our main application (entity-linking) the structure of the graph within each X_i is not very important, and can actually be misleading: a very central node in a large candidate set may seem very promising (and may actually minimize the distance to the other sets) but can be blatantly wrong. It is pretty much like the distinction between nepotistic and non-nepotistic links in PageRank computation: the links *within* each host are not very useful in determining the importance of a page—on the contrary, they may be confusing, and are thus often disregarded.

Based on this observation, we can (and probably want to) modify the structure of the graph within each set X_i to avoid this kind of trap. This is done by preserving the *external* links (those that connect vertices of X_i to the outside), but at the same time adding or deleting edges within each X_i in a suitable way. In our experiments, we considered two possible approaches:

- one consists in making $G[X_i]$ *maximally connected*, i.e., transforming it into a clique;
- the opposite approach makes $G[X_i]$ *minimally connected* by adding the minimum number of edges needed to that purpose; this can be done by computing the connected components of $G[X_i]$ and then adding enough edges to join them in a single connected component.

Both approaches guarantee that $G[X_i]$ is connected, so that the two heuristics described below can be applied.

5.1 The spanning-tree heuristic

The first heuristic aims at modifying the graph G in such a way that the resulting instance becomes decomposable. For the moment, let us assume that the sets X_i are pairwise disjoint. To guarantee that the problem be decomposable, we proceed as follows. Define an equivalence relation $\sim$ on V by letting $x \sim y$ whenever x and y belong to the same X_i.[4] The quotient graph $G/\sim = (V/\sim, E/\sim)$ has vertices $V/\sim$ and an edge between $[x]$ and $[y]$ whenever there is some edge $(x',y') \in E$ with $x' \sim x$ and $y' \sim y$ (here, and in the following, $[x]$ denotes the $\sim$-equivalence class including x). Thus, there is a surjective (but not injective) map $\iota : E \to E/\sim$.

Since G is connected, so is $G/\sim$, and we perform a breadth-first traversal of G building a spanning tree T. Every tree edge is an edge of $G/\sim$, so its pre-image with respect to ι is a nonempty set of edges in G. Let us arbitrarily choose one edge of G from $\iota^{-1}(t)$ for every tree edge t, and let T' be the resulting set of edges of G.

Define the new graph $G' = (V,E')$ where $E' = T' \cup \bigcup_{i=1}^{k} E(G[X_i])$: this graph cointains all the edges within each set X_i, plus the set T' of external edges.

It is easy to see that $G'[X_i]$ is connected (it is in fact equal to $G[X_i]$), and moreover all the elements of T' are bridges dividing all the X_i's in distinct biconnected components. In other words, we have turned the instance into a *decomposable* one, where Algorithm 1 can be run.

The non-disjoint case If the sets X_i are not pairwise disjoint, we can proceed as follows. Let us define maximal mutually disjoint sets of indices $I_1, \ldots, I_h \subseteq \{1, \ldots, k\}$ such that for all $t \neq s$, $\cup_{i \in I_t} X_i \cap \cup_{i \in I_s} X_i = \emptyset$.

Now, take the new problem instance with the same graph and sets $Y_1, \ldots, Y_h$ where $Y_t = \cup_{i \in I_t} X_i$: this instance is disjoint, so the previous construction applies. The only difference is that, at the very last step of Algorithm 1, when we are left with a graph and a *single* Y_t, we will not select a single $y \in Y_t$ optimizing the cost function

$$\sum_{z \in B} b(z) d(y,z).$$

Rather, we will choose one element x_i for every $i \in I_t$ optimizing

$$\sum_{i \in I_t} \sum_{z \in B} b(z) d(x_i,z).$$

Discussion Both steps presented above introduce some level of imprecision, that make the algorithm only a heuristic in the general case. The first approximation is due to the fact that building a tree on G will produce distances (between vertices living in different X_i) much larger than they are in G; the second approximation is that when we have non-disjoint sets, we only optimize with respect to bridges, disregarding the sum of distances of the nodes of different sets. Actually, we should optimize

$$\sum_{i \in I_t} \sum j \in I_t d(x_i,x_j) + \sum_{i \in I_t} \sum_{z \in B} b(z) d(x_i,z).$$

but this would make the final optimization step NP-complete.

[4] Note that, since the sets X_i are pairwise disjoint, $\sim$ is transitive.

5.2 The hitting-distance heuristic

The second heuristic we propose is based on the notion of *hitting distance*: given a vertex x and a set of vertices Y, define the hitting distance of x to Y as $d(x,Y) = \min_{y \in Y} d(x,y)$. The hitting distance can be easily found by a breadth-first traversal starting at x and stopping as soon as an element of Y is hit. Given a general connected instance of MINDR, as described above, we can consider, for every i and every $x \in X_i$, the average hitting distance of x to the other sets:

$$\frac{\sum_{j=1}^{k} d(x,X_j)}{k}.$$

The element $x_i^* \in X_i$ minimizing the average hitting distance (or any such an element, if there are many) is the candidate chosen for the set X_i in that solution.

The main problem with this heuristic is related to its locality (optimization is performed separately for each X_i); moreover the worst-case complexity is $O(m\sum_i |X_i|)$, that reduces to $O(k \cdot m)$ only under the restriction that the sets X_i have $O(1)$ size.

6 Experiments

All our experiments were performed on a snapshot of the English portion of Wikipedia as of late February 2013; the graph (represented in the BVGraph format [3]) was symmetrized and only the largest component was kept. The undirected graph has 3 685 351 vertices (87.2% of the vertices of the original graph) and 36 066 162 edges (99.9% of the edges of the original graph). Such a graph will be called the "Wikipedia graph" and referred to as G throughout this experimental section.

Our experiments use actual real-world entity-linking problems for which we have a human judgment, and tries the two heuristics proposed in Section 5, as well as a greedy baseline and other heuristics.

The greedy baseline works as follows: it first chooses an index i at random, and draws an element $x_i \in X_i$ also at random. Then, it selects a vertex of $x_{i+1} \in X_{i+1}, x_{i+2} \in X_{i+2}, \ldots, x_k \in X_k, x_1 \in X_1, \ldots, x_{i-1} \in X_{i-1}$ (in this order) minimizing each time the sum of the distances to the previously selected vertices; the greedy algorithm continues doing the same also for $x_i \in X_i$ to get rid of the only element (the first one) that was selected completely at random. Moreover we have considered also two other heuristics, that have been observed to be effective in practice [9]: these are *degree* and *PageRank based*. They respectively select the highest degree and the highest PageRank vertex for each set.

The real-world entity-linking dataset has been taken from [11] which contains a larger number of human-labelled annotations. For retrieving the candidates, we created an index over all Wikipedia pages with different fields (title, body, anchor text) and used a variant of BM25F [2] for ranking, returning the top 100 scoring candidate entities. Since the candidate selection method was the same for every graph-based method employed, there should be no bias in the experimental outcomes.

The problem instances contained in the dataset have 11.73 entities on average (with a maximum of 53), and the average number of candidates per entity is 95.90 (with a maximum of 200). Each of the 100 problem instances in the NEL dataset is annotated, and in particular, for every i there is a subset $X_i^* \subseteq X_i$ of *fair* vertices (that is, vertices that are good candidates for that set): typically $|X_i^*| = 1$. Note that, for every instance in the NEL dataset, we deleted the sets X_i such that X_i^* were not included in the largest connected component of the Wikipedia graph. The number of sets X_i deleted was at maximum 2 (for two instances). We have not considered instances in which, after these modifications, we have just one set X_i: this situation happened in 5 cases. So the problem set on which we actually ran our algorithm contains 95 instances.

	Distance-cost ratio		Value	
	Maximal Connection	Minimal Connection	Maximal Connection	Minimal Connection
Heuristic	Average (± Std Error)	Average (± Std Error)	Average (± Std Error)	Average (± Std Error)
Spanning-tree	122.747(±2.812)	130.998 (±2.917)	0.369 (±0.023)	0.360 (±0.023)
Hitting-distance	103.945 (±1.320)	105.797 (±2.322)	**0.454 (±0.027)**	**0.459 (±0.027)**
Greedy	**101.969 (±0.429)**	**102.785 (± 0.426)**	0.428 (±0.025)	0.426 (±0.026)
Degree based	114.182 (±2.386)	113.285 (±2.305)	0.411 (±0.024)	0.394 (±0.023)
PageRank based	114.894 (±2.452)	112.392 (±2.266)	0.407 (±0.025)	0.398 (±0.023)
Ground truth	115.117 (±1.782)	119.243 (±1.873)		

Table 1: Distance-cost ratio and value.

For every instance, we considered the maximal and minimal connection [5] approach, and then ran both heuristics described in Section 5, comparing them with the greedy baseline, and also with the degree and PageRank heuristics.

For any instance, when comparing the distance cost f of the solutions S_j returned by some algorithm A_j, we have computed the *distance-cost ratio* of each algorithm A_j, defined as

$$\frac{f(S_j)}{\min_j f(S_j)} \cdot 100.$$

Intuitively this corresponds to the approximation ratio of each solution with respect to the best solution found by all the considered algorithms: hence the best algorithm has minimum distance-cost ratio and it equals 100.

Besides evaluating the distance cost of the solutions found by the various heuristics, we can compute how many of the elements found are fair: we normalize this quantity by k, so that 1.0 means that all the k candidates selected are fair. We call such a quantity the *value* of a solution.

In the last two columns of Table 1 we report, for each heuristic, the average value (across all the instances) along with the standard error. For both the connection approaches, we have that the hitting-distance heuristic outperforms all the other heuristics, and it selects more than 45% of fair candidates. The variability of the results seems not to differ too much for all the methods. The second best heuristic is the greedy baseline, that selects almost 42.8% and 42.6% fair candidates respectively in a maximal and minimal connected scenario.

It is worth observing that the greedy approach comes second (as far as the value is concerned), and outperforms the baseline techniques (degree and PageRank). The spanning tree heuristic, instead, perform worse than any other method.

The latter outcome is easily explained by the fact that it transforms completely the topology of the graph in order to make the instance decomposable, and the distances between vertices are mostly scrambled. This interpretation of the bad result obtained can also be seen looking at the distance cost (central columns of Table 1): the spanning-tree heuristic is the one that is less respectful of distances, selecting candidates that are far apart from one another.

In the central columns of Table 1, we report also the distance-cost ratio for all the other heuristics. For both the maximal and the minimal connection approaches, the greedy baseline seems to obtain more

[5] To obtain the minimal connection of each $G[X_i]$, we chose to connect the vertex of maximum degree of its largest component with an (arbitrary) vertex of each of its remaining components.

often a minimum distance cost solution. The second best option is the hitting distance heuristic, while the other methods seems to be more far away from an optimal result.

In the last row of Table 1, we report the distance-cost ratio for the ground-truth solution given by the fair candidates. It seems that for any instance, the ground truth has distance cost averagely 15%-20% higher than the best solution we achieve by using the heuristics. This observation suggests that probably our objective function (that simply aims at minimizing the graph distances) is too simplistic: the distance cost is an important factor to be taken into account but certainly not the unique one.

It is interesting to remark, though, that the average Jaccard coefficient between the solution found by the degree based and the hitting-distance heuristic is 0.3 (for both maximal and minimal connection approaches): this fact means that the degree and distance can be probably used as complementary features that hint at different good candidates, although we currently do not know how to combine these pieces of information.

Finally, we remark that we also tried to apply the degree and PageRank based heuristics by using the same problem set but *in the original directed graph*; in this case, we did not enforce any connectivity of the subgraphs $G[X_i]$: the resulting average values ($\pm$ standard error) are respectively 0.327 ($\pm$0.020) and 0.336 ($\pm$0.022), and they are both worse than the values achieved by degree and PageRank heuristics in Table 1. This fact suggests that our experimental approach (of considering the undirected version and of enforcing some connectivity on the subgraphs) not only guarantees the applicability of our heuristics in a more suitable scenario, but also improves the effectiveness of the other existing techniques.

7 Conclusions and future work

Inspired by the entity-linking task in NLP, we defined and studied a new graph problem related to Maximum Capacity Representative Set and we proved that this problem is NP-hard in general (although it remains an open problem to determine its exact approximability). Morevoer, we showed that the problem can be solved efficiently in some special case, and that we can anyway provide reasonable heuristics for the general scenario. We tested our proposals on a real-world dataset showing that one of our heuristics is very effective, actually more effective than other methods previously proposed in the literature, and more than a simple greedy approach using the same cost function adopted here.

The other heuristic proposed in this paper seem to work poorly (albeit it reduces to a case where we know how to produce the optimal solution), but we believe that this is just because of the very rough preprocessing phase it adopts; we plan to devise a more refined way to induce the conditions needed for Algorithm 1 to work, without having to resort to the usage of a spanning tree—the latter scrambles the distances too much, resulting in a bad selection of candidates.

Finally, we observed that a distance-based approach is complementary to other methods (e.g., the local techniques based solely on the vertex degree), hinting at the possibility of obtaining a new, better cost function that exploits both features at the same time.

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A Proof of Lemma 1

Proof. We reduce MAXCRS to MINDIR. Given an instance of MAXCRS, $\{X_1,\dots X_k\}$ and for any $i \neq j$, $x \in X_i$, and $y \in X_j$, a nonnegative capacity $c(x,y)$, we construct the instance of MINDIR $G, \{X_1,\dots,X_k\}$; the vertices of G are $X_1 \cup \dots \cup X_k$, and for any pair $x \in X_i$, $y \in X_j$, with $i \neq j$, we add a weighted edge between x and y, i.e., for each pair for which MAXCRS defines a capacity we create a corresponding edge in G. In particular the weight of the edge between x and y is set to $\alpha - c(x,y)$, where $\alpha = 2\max_{z\in X_i, t\in X_j, i\neq j} c(z,t)$.

Observe that for any pair of nodes $u \in X_i$, $v \in X_j$, with $i \neq j$, $d(u,v)$ in G is equal to the weight of (u,v), i.e., it is not convenient to pass through other nodes when going from u to v: in fact, for any path $z_1,\dots,z_p$ from u to v in G, with $p \geq 1$, we always have $\alpha - c(u,v) \leq \alpha - c(u,z_1) + \dots + \alpha - c(z_p,v)$, since $\alpha - c(u,v) \leq \alpha$ and the weight of such a path is at least $\frac{p+1}{2}\alpha \geq \alpha$. Moreover, observe that any optimal solution in G has exactly one element for each set X_i: thus, we have $k(k-1)$ pairs of elements (x,y), whose distance is always given by the weight of the single edge (x,y), that is $\alpha - c(x,y)$.

Hence it is easy to see that MAXCRS admits a system of representatives whose capacity is greater than h, if and only if MINDIR admits a solution S such that $f(S)$ is less than $k(k-1)\alpha - h$.

Since MINDIR is a restriction of MINDR we can conclude that also MINDR is NP-complete. □

B Proof of Theorem 1

Proof. We can rewrite the objective function as follows.

$$\begin{aligned}\sum_{x_i,x_j\in S} d(x_i,x_j) + \sum_{x_i\in S}\sum_{z\in B} d(x_i,z)b(z) \;=\; & 2|Y_0||Y_1|d(t_0,t_1) + \sum_{x_i,x_j\in S\cap V(Y_0)} d(x_i,x_j) + \sum_{x_i,x_j\in S\cap V(Y_1)} d(x_i,x_j) + \\ & 2|Y_1|\sum_{x_i\in S\cap V(Y_0)} d(x_i,t_0) + \sum_{x_i\in S\cap V(Y_0)}\sum_{z\in B} d(x_i,z)b(z) + \\ & 2|Y_0|\sum_{x_j\in S\cap V(Y_1)} d(t_1,x_j) + \sum_{x_i\in S\cap V(Y_1)}\sum_{z\in B} d(x_i,z)b(z).\end{aligned}$$

This is because if $z \in B\cap Z_1$, for any node $x_i \in S\cap V(Y_0)$, we have $d(x_i,z) = d(x_i,t_0) + d(t_0,z)$ (and analogously, if $z \in B\cap Z_0$, for any node $x_i \in S\cap V(Y_1)$, we have $d(x_i,z) = d(x_i,t_1) + d(t_1,z)$). Hence:

$$\sum_{x_i\in S\cap V(Y_0)}\sum_{z\in B} d(x_i,z)b(z) \;=\; \sum_{x_i\in S\cap V(Y_0)}\sum_{z\in B\cap Z_0} d(x_i,z)b(z) + \sum_{x_i\in S\cap V(Y_0)}\sum_{z\in B\cap Z_1} d(x_i,t_0)b(z) + d(t_0,z)b(z)$$

and

$$\sum_{x_i\in S\cap V(Y_1)}\sum_{z\in B} d(x_i,z)b(z) \;=\; \sum_{x_i\in S\cap V(Y_1)}\sum_{z\in B\cap Z_1} d(x_i,z)b(z) + \sum_{x_i\in S\cap V(Y_1)}\sum_{z\in B\cap Z_0} d(x_i,t_1)b(z) + d(t_1,z)b(z).$$

Observe that t_0 or t_1 might already belong to B: this is why we assumed that B is a multiset.

Then, we have that:

$$f(S_0) = \sum_{x_i,x_j\in S\cap V(Y_0)} d(x_i,x_j) + \sum_{x_i\in S\cap V(Y_0)}\sum_{z\in B\cap Z_0} d(x_i,z)b(z) + \sum_{x_i\in S\cap V(Y_0)} d(x_i,t_0)\cdot\left(2|Y_1| + \sum_{z\in B\cap Z_1} b(z)\right)$$

$$f(S_1) = \sum_{x_i,x_j \in S \cap V(Y_1)} d(x_i,x_j) + \sum_{x_i \in S \cap V(Y_1)} \sum_{z \in B \cap Z_1} d(x_i,z)b(z) + \sum_{x_i \in S \cap V(Y_1)} d(x_i,t_1) \cdot \left(2|Y_0| + \sum_{z \in B \cap Z_0} b(z) \right)$$

Hence, by adding t_s to $B \cap Z_s = B_s$, with weight equal to $b_s = 2|Y_{1-s}| + \sum_{z \in B \cap Z_{1-s}} b(z)$, $f(S)$ can be reduced to $f(S_0)$ and $f(S_1)$. □

C The algorithm for finding useful edges

Algorithm 2: USEFULEDGE

Input: An instance $G, \{X_1, \ldots, X_k\}, B, b$
Output: A useful edge, or null
Pick a node u of the set X_i of the instance $G, \{X_1, \ldots, X_k\}, B, b$
Mark all the nodes as unseen
dfs[] $\leftarrow -1$, visited $\leftarrow 0$, usefulEdgeFound $\leftarrow$ *false*, usefulEdge $\leftarrow$ *null*
DFS(u,-1)
if usefulEdgeFound **then**
  return usefulEdge
else
  return *null*
end

Algorithm 3: DFS

Input: A node u, its parent p
Output: A pair (t,Y), where t is an integer and Y is a set of indices
if usefulEdgeFound **then return** *null* Mark u as seen
dfs[u] $\leftarrow$ visited
visited $\leftarrow$ visited $+ 1$
furthestAncestor $\leftarrow$ visited
$Y \leftarrow \emptyset$
if $t \in X_i$ **then** $Y \leftarrow Y \cup \{i\}$ **for** $v \in N(u)$ *s.t.* $w \neq p$ **do**
  if v *is unseen* **then**
    $(t', Y') \leftarrow$ DFS(v,u)
    if $t' >$ dfs[u] *and* $\emptyset \neq Y' \neq \{1, \ldots, k\}$ **then**
      usefulEdgeFound $\leftarrow$ *true*
      usefulEdge $\leftarrow (u, v)$
      return *null*
    end
    furthestAncestor $\leftarrow$ min(furthestAncestor,t')
    $Y \leftarrow Y \cup Y'$
  else
    furthestAncestor $\leftarrow$ min(furthestAncestor, dfs[v])
  end
end
return (furthestAncestor,Y)