A NEW METHOD OF MEASURING SIMILARITY FOR A SPECIAL CLASS OF DIRECTED GRAPHS

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ABSTRACT. The problem of graph similarity is challenging and important in many areas of science, e.g., mathematics [Sobik, F.: Graphmetriken und Klassifikation strukturierter Objekte, ZKI-Informationen, Akad. Wiss. DDR, 2, (1982), 63-122]; [Zelinka, B.: On a certain distance between isomorphism classes of graphs, Cas. Pest. Mat., 100, (1975), 371-373], biology [Koch, I., Lengauer, T.; Wanke, E.: An algorithm for finding maximal common subtopologies in a set of protein structures, J. Comput. Biology, 3, (1996), 289-306], and chemistry [Skvortsova, M. I., Baskin, I. I., Stankevich, I. V., Palyulin, V. A., Zerzow, N. S.: Molecular similarity in structure-property relationship studies. Analytical description of the complete set of graph similarity measures, International Symposium CACR’96, (1996) pp. 542-646]. In this paper, we design a new method, which uses sequence alignment techniques [Altschul, S. F., Gish, W., Miller, V., Myers, E. W., Lipman, D. J.: Basic local alignment search tool, J. Molecular Biology 125, (1991), 403-410]; [Altschul, S. F., Madden, T. L., Miller, W., Schaffer, A. A., Zhang, J., Zhang, Z., Lipman, D. J.: Gapped BLAST and PSI-BLAST: a new generation of protein database search programs, Nucleic Acids Res. 25, (1997), 3389-3402]. [Kilian, J., Hoos, H. H.: MusicBLAST—gapped sequence alignment for MIR, in: Proceedings of the 5th International Conference on Music Information Retrieval (ISMIR), (2004)], to measure the structural similarity of unlabeled, hierarchical, and directed graphs. More precisely, if $h$ denotes the maximal length of a path from the root to a leaf of a given hierarchical and directed graph $\mathcal{H}$, we align out-degree and in-degree sequences induced by the vertex sequences on a level $i$, $0 \leq i \leq h$. On the basis of the level alignments, we construct measured values and prove that they are similarity measures. In our algorithm, which uses the well-known technique of dynamic programming, the alignments of out-degree and in-degree sequences are decoupled. Therefore, we obtain a family $(d_i(\mathcal{H}_1, \mathcal{H}_2))_{1 \leq i \leq n}$ of graph similarity measures. As an application, we examine the measures on a graph corpus of 464 graphs, where the graphs represent web-based hypertext structures (websites).

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1. Introduction

In this paper, we introduce novel graph similarity measures for a special class of directed graphs. First, we have to define the term graph similarity. In general, the concept of similarity is not unique and thus cannot be completely formalized. If we consider two structured objects, we can refer to several similarity aspects, e.g., semantic similarity, structural similarity, and functional similarity. In the following, we introduce similarity measures on unlabeled, hierarchical and directed graphs, which are based on structural similarity. In order to construct such a measure, we have to take into account the structural properties and parameters of the graphs under consideration.

Measures of distances between graphs have been frequently investigated. In the literature, the problem of computing the distance between graphs is often called *inexact graph matching* [7]. Many similarity measures on graphs are based on isomorphic relations and subgraph isomorphism [25], [31], respectively. An example of such a similarity measure is the well-known *Zelinka*-distance [32]. The *Zelinka*-distance is based on the principle that the more similar two graphs are, the bigger the common induced subgraph is. In other words, graphs which have a large common induced subgraph, have a small distance and vice versa. *Zelinka* was the first to introduce this measure for unlabeled graphs. *Sobick* [28] [29] and *Kaden* [16], [17] generalized this measure for arbitrary (also labeled) graphs of different order and proved that it is a metric. But, for large graphs, the complexity is considered to be unacceptable for practical use.

*Kaden* has obtained further similarity measures on graphs by transforming the graphs by injective mapping. For example, *Kaden* [16] considered *line graphs* [4] and used the *Zelinka*-distance to compute the similarity of the transformed graphs.

*Sapiro* [26] introduced a known similarity metric for graphs based on the corresponding adjacency matrices. Let the graphs be $G_1, G_2$ and the corresponding adjacency matrices $A_1, A_2$. Permute the rows and the columns in the matrix $A_2$ in such a way that the matrix elements conform with the matrix elements of $A_1$ as much as possible. Now, *Sapiro* defines the graph distance between $G_1, G_2$ by the minimal number of dissenting matrix elements and proves that it is a metric.

In addition, an important class of similarity measures based on the *edit distance* of graphs has been investigated by [33] [34]. The edit distance is based on basic weighted transformation steps like deletions, substitutions, and insertions of vertices and edges. Since there is an infinite number of different possibilities for transforming $G_2$ in $G_1$, the similarity of the graphs is defined as the minimum cost of transformations.
Other approaches to inexact graph matching consider distances between graphs on the basis of graph grammars [10], [24]. These methods are primarily interesting for theoretical aspects but not for practical use, since the specific grammar is difficult to obtain.

This paper is organized as follows: In the next Section (2) we state some fundamental definitions and explain topological properties of our graphs in terms of out-degree and in-degree sequences. In Section 3 we present our new approach for measuring the structural similarity of unlabeled, hierarchical, and directed graphs. In Section (4) we apply our new method to a graph corpus $C_G$ and state the experimental results. The paper finishes in Section (5) with a summary and conclusions.

2. Fundamental definitions and topological aspects of graphs

**Definition 2.1.** Let $\mathcal{H} = (V, E)$, $E \subseteq V \times V$, $|V| < \infty$ be a directed graph.

$\mathcal{N}^+(v) := \{ \hat{v} \in V \setminus \{v\} | (v, \hat{v}) \in E \}$ denotes the set of out-neighbours of $v$,

$\mathcal{N}^-(v) := \{ \hat{u} \in V \setminus \{v\} | (\hat{u}, v) \in E \}$ denotes the set of in-neighbours of $v$,

$s^{\text{out}}_j(\mathcal{H}) \in \mathbb{N}$, $0 \leq j \leq k^{\text{out}} := \max \{ \delta^{\text{out}}_j(v) \}$ (or $s^{\text{in}}_i(\mathcal{H}) \in \mathbb{N}$, $0 \leq i \leq k^{\text{in}} := \max \{ \delta^{\text{in}}_i(v) \}$) denotes the number of vertices of $\mathcal{H}$ with out-degree $j$ (or in-degree $i$).

The vector $s^{\text{out}}(\mathcal{H}) := (s^{\text{out}}_0(\mathcal{H}), s^{\text{out}}_1(\mathcal{H}), \ldots, s^{\text{out}}_{k^{\text{out}}}(\mathcal{H}))$, or

$s^{\text{in}}(\mathcal{H}) := (s^{\text{in}}_0(\mathcal{H}), s^{\text{in}}_1(\mathcal{H}), \ldots, s^{\text{in}}_{k^{\text{in}}}(\mathcal{H}))$

is called the out-degree (or in-degree) sequence of $\mathcal{H}$. 

![Figure 1.1](image-url) Two directed graphs with the same degree sequences but $G_1 \neq G_2$. 

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*Figures and equations are not provided in this text.*
Degree sequences can be useful for describing the structure of a graph. As a preliminary, we note an important property of degree sequences. If $G_1, G_2$ are arbitrary directed graphs, then the following assertion holds: $G_1 \neq G_2 \Rightarrow s^{out}(G_1) = s^{out}(G_2) \land s^{in}(G_1) = s^{in}(G_2)$. The reverse assertion is not always true. For example, the degree sequences of the graphs $G_1, G_2$ shown in Figure (1.1) are $s^{out}(G_1) = (0, 0, 8) = s^{out}(G_2) = (0, 0, 8), s^{in}(G_1) = (0, 0, 8) = s^{in}(G_2) = (0, 0, 8)$, but $G_1 \neq G_2$. Therefore, we have to examine how far degree sequences are applicable for measuring the structural similarity of our graphs.

Another important question is: Given the sequences of natural numbers $(\delta_0, \delta_1, \ldots, \delta_p)$ and $(\delta_0, \delta_1, \ldots, \delta_p)$ under what conditions are the two sequences (out-degree and in-degree) sequences of a certain graph $H$? This problem has been solved by [12], [13], [14]. Now, we will construct the class of graphs that we want to examine in this paper.

**Definition 2.2.** Let be $\mathcal{H} = (V, E), E \subseteq V \times V, |V| < \infty$ and

$\hat{V} := \{v_{0,1}, v_{1,1}, v_{1,2}, \ldots, v_{1,\sigma_1}, v_{2,1}, v_{2,2}, \ldots, v_{2,\sigma_2}, \ldots, v_{k,1}, v_{k,2}, \ldots, v_{k,\sigma_k}\}$. Define $h$ as the maximal length of a path from the root $v_{0,1}$ to a leaf. $v_{i,j}$ denotes the $j$th vertex on the $i$th level, $0 \leq i \leq h, 1 \leq j \leq \sigma_i$. $\sigma_i$ is maximal in the sense that there is no other vertex sequence such that $v_{i,1}, v_{i,2}, \ldots, v_{i,\sigma_i}$ with $\hat{\sigma}_i > \sigma_i$. $L: \hat{V} \rightarrow \mathbb{N}, L(v_{i,j}) := i$ is a function which determines the level of a vertex $v_{i,j}$. The edge types are now defined by:

\[
\hat{E}_1 := \{(v_{i,\nu}, v_{i+1,\nu_1}) | v_{i,\nu}, v_{i+1,\nu_1} \in \hat{V}, 1 \leq j \leq k, k := \delta_{out}(v_{i,\nu}) ,
\]

\[
L(v_{i+1,\nu_1}) = L(v_{i,\nu}) + 1 \land ((\hat{\delta}(v_{i,\nu}, v_{i+1,\nu_1}), \nu > \nu)
\]

\[
\lor (\hat{\delta}(v_{i,\nu}, v_{i+1,\nu_1}), \nu < \nu) , \quad v_1 < v_2 < \cdots < v_k \}
\]
\[ \hat{E}_2 := \{ (v_{i+s}, v_{i+p}, v_{i+p}, v_{i+p}) \mid v_{i+s}, v_{i+p} \in \hat{V}, \mathcal{L}(v_{i+p}) = \mathcal{L}(v_{i+p}) - s, s \leq h \]
\[ \land \exists! \left( (v_{i+p}, v_{i+p+1}, v_{i+p+1}), \ldots, (v_{i+s-1}, v_{i+s-1}, v_{i+s}), 1 \leq \bar{p} \leq \sigma_j, 1 \leq \nu \leq \sigma_{i+1}, \ldots, \right) \]
\[ 1 \leq v_j \leq \sigma_{i-s-1}, 1 \leq \nu \leq \sigma_{i+s} \}, \tag{2.2} \]

\[ \hat{E}_3 := \{ (v_{i+p}, v_{i+s}, v_{i+s}, v_{i+s}) \mid v_{i+p}, v_{i+s} \in \hat{V}, \mathcal{L}(v_{i+s}) = \mathcal{L}(v_{i+p}) + s, 1 < s \leq h \]
\[ \land \exists! \left( (v_{i+p}, v_{i+p+1}, v_{i+p+1}), \ldots, (v_{i+s-1}, v_{i+s-1}, v_{i+s}), 1 \leq \bar{p} \leq \sigma_j, 1 \leq \nu \leq \sigma_{i+1}, \ldots, \right) \]
\[ 1 \leq v_j \leq \sigma_{i-s-1}, 1 \leq \nu \leq \sigma_{i+s} \}, \tag{2.3} \]

\[ \hat{E}_4 := \{ (v_{i+u}, v_{i+p}, v_{i+p}) \mid v_{i+p}, v_{i+p} \in \hat{V}, \mathcal{L}(v_{i+p}) = \mathcal{L}(v_{i+p}) - s, s \leq h \}
\[ \cup \{ (v_{i+s}, v_{i+s}, v_{i+s}, v_{i+s}) \mid v_{i+s}, v_{i+s}, v_{i+s} \in \hat{V}, (v_{i+s}, v_{i+s}, v_{i+s}) \notin \hat{E}_2, \mathcal{L}(v_{i+s}) \]
\[ = \mathcal{L}(v_{i+s}) - s, s \leq h \}
\[ \cup \{ (v_{i+s}, v_{i+s}, v_{i+s}, v_{i+s}) \mid v_{i+s}, v_{i+s}, v_{i+s} \in \hat{V}, (v_{i+s}, v_{i+s}, v_{i+s}) \notin \hat{E}_1, \hat{E}_3, \mathcal{L}(v_{i+s}) \}
\[ = \mathcal{L}(v_{i+s}) + s, s \leq h \}. \tag{2.4} \]

Then \( \mathcal{H} := (\hat{V}, \hat{E}) \), \( \hat{E} := \hat{E}_1 \cup \hat{E}_2 \cup \hat{E}_3 \cup \hat{E}_4 \) denotes the hierarchical and directed graph of \( \mathcal{H} \).

We call the elements of the edge set of \( \mathcal{H} \) Kernel-edges (2.1), Up-edges (2.2), Down-edges (2.3) and Across-edges (2.4). The properties of each edge type are, briefly [23]:

- **Kernel-edges**: The Kernel hierarchy is induced by the Kernel-edges. Kernel-edges associate dominating nodes with their immediately dominated successor nodes.
- **Up-edges** associate analogously nodes of the Kernel hierarchy with one of their (dominating) predecessor nodes.
- **Down-edges** associate nodes of the Kernel hierarchy with one of their (dominated) successor nodes in terms of that Kernel hierarchy.
- **Across-edges** associate nodes of the Kernel hierarchy, none of which is an (immediate) predecessor of the other in terms of the Kernel hierarchy.

The Definition (2.2) provides a structural property of \( \mathcal{H} \) which is evident.

**Proposition 2.1.** Let \( \mathcal{H} := (\hat{V}, \hat{E}) \), \( \mathcal{H}_T := (\hat{V}, E_T) \), \( E_T := \hat{E} \setminus \{ \hat{E}_2, \hat{E}_3, \hat{E}_4 \} \) is a directed rooted tree.
Corollary 2.2. $\hat{\mathcal{H}}_{\text{rec}}^{\text{top}} := (w, \hat{\mathcal{H}}_{\mathcal{H}_1}, \hat{\mathcal{H}}_{\mathcal{H}_2}, \ldots, \hat{\mathcal{H}}_{\mathcal{H}_{\text{out}(w)}})$ is the recursive description of $\hat{\mathcal{H}}_{\mathcal{H}}$.

Now we will provide examples of graphs which fulfill Definition (2.2), but whose topology cannot be adequately described by degree sequences. With reference to the next chapter, we note that the vertices on each level $i$ induce an out-degree and in-degree sequence in a left to right order. In particular, the out-degrees from each vertex $v_{i,j}$ on a level $i$, $0 \leq i \leq h$ are significant for the embeddings of the substructures associated with $v_{i,j}$.

By Definition (2.1), the graphs in Figure (2.2) have the same out-degree and in-degree sequences, $s^{\text{out}}(\hat{\mathcal{H}}_1) = (5, 0, 2, 1) = s^{\text{out}}(\hat{\mathcal{H}}_2) = (5, 0, 2, 1)$ and $s^{\text{in}}(\hat{\mathcal{H}}_1) = (1, 7) = s^{\text{in}}(\hat{\mathcal{H}}_2) = (1, 7)$. But they are not symmetrically located on the symmetry axis, indicated by the dashed lines. Therefore, we see that degree sequences of the given graphs have no influence on the embeddings of substructures in the graph.

An other example is Figure (2.3). It holds that $s^{\text{out}}(\hat{\mathcal{H}}_1) = (6, 1, 3, 0, 1) = s^{\text{out}}(\hat{\mathcal{H}}_2) = (6, 1, 3, 0, 1)$ and $s^{\text{in}}(\hat{\mathcal{H}}_1) = (0, 11) = s^{\text{in}}(\hat{\mathcal{H}}_2) = (0, 11)$. Nonetheless, $\hat{\mathcal{H}}_1$ and $\hat{\mathcal{H}}_2$ possess different topologies. Altogether, it follows that simple comparisons of out-degree and in-degree vectors are not suitable for determining the structural similarity of our graphs.

3. New approach for measuring the structural similarity of graphs

In Section (2) we saw that degree sequences cannot describe the topology of a graph completely. Since we are examining unlabeled, hierarchical and directed graphs, in the following, we will look at the out-degree and in-degree sequences (on a level $i$), induced by the vertex sequences $v_{i,1}, v_{i,2}, \ldots, v_{i,n}$, and their edge relations in terms of Definition (2.2). Now, with respect to a cost function $\alpha$ the
more similar the out-degree and in-degree alignments on the levels $i$, $0 \leq i \leq h$ are, the more similar the common structure of the graphs is and vice versa. The cost function $\alpha$, which we will define later, weighs alignments on the sequences $S_0^\hat{H}$, $S_1^\hat{H}$, $\ldots$, $S_h^\hat{H}$ under certain conditions. With $w_i^\hat{H} := v_{0,i}^\hat{H}$, the structural embedding of a graph $\hat{H}$ is essentially described by the sequences

$$S_0^\hat{H} := v_1^\hat{H},$$

$$S_1^\hat{H} := v_{1,1}^\hat{H} \circ v_{1,2}^\hat{H} \circ \cdots \circ v_{1,\delta_{out}(w_1^\hat{H})}^\hat{H},$$

$$\vdots$$

$$S_h^\hat{H} := v_{h,1}^\hat{H} \circ v_{h,2}^\hat{H} \circ \cdots \circ v_{h,\sigma_h}^\hat{H},$$

together with their out-degree and in-degree sequences on a level $i$.

Now, the problem of determining the structural similarity between $\hat{H}_1$ and $\hat{H}_2$ is equivalent to determining the optimal alignment of

$$S_0^{\hat{H}_1} := w_1^{\hat{H}_1},$$

$$S_1^{\hat{H}_1} := v_{1,1}^{\hat{H}_1} \circ v_{1,2}^{\hat{H}_1} \circ \cdots \circ v_{1,\delta_{out}(w_1^{\hat{H}_1})}^{\hat{H}_1},$$

$$\vdots$$

$$S_h^{\hat{H}_1} := v_{h,1}^{\hat{H}_1} \circ v_{h,2}^{\hat{H}_1} \circ \cdots \circ v_{h,\sigma_h}^{\hat{H}_1},$$

and
with respect to a cost function $\alpha$. Based on these vertex sequences, we will later implement alignments of the corresponding out-degree and in-degree sequences. We regard this task as an optimization problem, in the sense of finding a minimum score path in an alignment graph (see Definition (3.1)). For this purpose, we use the method of dynamic programming, which is based on the optimality principle of Bellman [6].

Generally, the method of dynamic programming can be viewed as an optimization of a process $\mathcal{P}_{fin} = \mathcal{P}_1 \circ \mathcal{P}_2 \circ \cdots \circ \mathcal{P}_n$. Starting from the initial state $\mathcal{P}_0$ and a control $c_1$, the new state $\mathcal{P}_1$ will be calculated with a transition function $\mathcal{P}_1 = f_1(\mathcal{P}_0, c_1)$, analogous for $\mathcal{P}_2 = f_2(\mathcal{P}_1, c_2)$, ..., $\mathcal{P}_n = f_n(\mathcal{P}_{n-1}, c_n)$. Now, we have to optimize a target function $F(\mathcal{P}_0, \mathcal{P}_1, \ldots, \mathcal{P}_n, c_1, \ldots, c_n)$. $F$ is additive such that there exists function $\text{add}_i$ with $F = \sum_{i=1}^{n} \text{add}(\mathcal{P}_{i-1}, c_i)$.

Now, let $\mathcal{H}_1, \mathcal{H}_2, v_{i,j}^{\mathcal{H}_1}, 0 \leq i \leq h_1, 1 \leq j \leq \sigma_i$ denote the $j$th vertex on the $i$th level of $\mathcal{H}_1$, analogous to $v_{i,j}^{\mathcal{H}_2}$ for $\mathcal{H}_2$. We define the sequences representing graphs

$$S_1 := v_{1,1}^{\mathcal{H}_1} \circ v_{1,2}^{\mathcal{H}_1} \circ \cdots \circ v_{h_1,\sigma_{h_1}}^{\mathcal{H}_1},$$

$$S_2 := v_{2,1}^{\mathcal{H}_2} \circ v_{2,2}^{\mathcal{H}_2} \circ \cdots \circ v_{h_2,\sigma_{h_2}}^{\mathcal{H}_2},$$

(3.1)

$S_k[i]$ denotes the $i$th position of the sequences $S_k$ and it holds $S_1[n] = v_{h_1,\sigma_{h_1}}, S_2[m] = v_{h_2,\sigma_{h_2}}, \mathbb{N} \ni n, m \geq 1$, $S_k[1] = w_k^{\mathcal{H}_k}, k \in \{1,2\}$. In order to find the optimal alignment between $S_1, S_2$, we construct the alignment graph $G_{S_1, S_2} := (V_{S_1, S_2}, E_{S_1, S_2}, f_{E_{S_1, S_2}})$ with an edge labeling function $f_{E_{S_1, S_2}} : E_{S_1, S_2} \rightarrow \mathbb{R}_+^+$.

**Definition 3.1.** Let $V_{S_1, S_2} := \{(i, j) | 0 \leq i \leq n, 0 \leq j \leq m\}$, $e_{Del} := (i-1, j) \rightarrow (i, j)$, $e_{Ins} := (i, j-1) \rightarrow (i, j)$, $e_{Subst} := (i-1, j-1) \rightarrow (i, j)$. The edge set $E_{S_1, S_2}$ is now defined by

$$E_{S_1, S_2} := \{e_{Del} | f_{E_{S_1, S_2}}(e_{Del}) = [S_1[i], -], i \in [1, n]\} \cup \{e_{Ins} | f_{E_{S_1, S_2}}(e_{Ins}) = [\-, S_2[j]], j \in [1, m]\} \cup \{e_{Subst} | f_{E_{S_1, S_2}}(e_{Subst}) = [S_1[i], S_2[j]], i \in [1, n], j \in [1, m]\}.$$
$G_{S_1,S_2} := (V_{S_1,S_2}, E_{S_1,S_2}, f_{E_{S_1,S_2}})$ denotes the alignment graph of the sequences $S_1$ and $S_2$.

$(i-1, j) \rightarrow (i, j)$ equals the deletion of $S_1[i]$ in $S_1$, $(i, j-1) \rightarrow (i, j)$ equals the insertion of $S_2[j]$ in $S_1$ at the $i$th position, and $(i-1, j-1) \rightarrow (i, j)$ equals the substitution $S_1[i]$ to $S_2[j]$.

![Alignment graph $G_{S_1,S_2}$ of the sequences $S_1$, $S_2$.](image)

As an application, we consider the two graphs $\hat{\mathcal{H}}_1, \hat{\mathcal{H}}_2$ represented as the sequences $S_1 := v_0 \circ v_1$ and $S_2 := v_1 \circ v_0 \circ v_1$ (simplified notation). The corresponding alignment graph is shown in Figure (3.1). From this figure we can read off (bold edges) the alignment:

```
- v_0 v_1
  v_1 v_0 v_1
```

According to the edge labeling function $f_{E_{S_1,S_2}} : E_{S_1,S_2} \rightarrow \mathbb{R}_+$ and for each possible aligned pair $[a, b]$ a cost function $a([a, b]) \in \mathbb{R}_+$ is assigned, where $a, b$ are sequence entries of $S_1$ and $S_2$ or the gap symbol $\cdot$. Our algorithm with the complexity $O(|\hat{V}_1| \cdot |\hat{V}_2|)$ for finding the optimal alignment of the Sequences (3.1), (3.2) generates a matrix $(M(i,j))_{ij}$, $0 \leq i \leq n$, $0 \leq j \leq m$, where $M(i,j)$ is equivalent to the minimal edit distance between the sequences $\hat{S}_1, \hat{S}_2$. Thereby $\hat{S}_1, \hat{S}_2$ consist of the first $i$th, $j$th characters of $S_1$ and $S_2$. 

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Hence, we are searching for \( \mathcal{M}(n, m) \) because \( \mathcal{M}(n, m) \) is the minimal edit distance
\[
d(S_1, S_2) := \min_{S_1 \to S_2} \sum \alpha([a, b]). \tag{3.3}
\]
We find the optimal alignment by tracing back along the minimal values from \( \mathcal{M}(n, m) \) to \( \mathcal{M}(0, 0) \), since we notionally assign pointers to the edit operations. It is well known that the Distance (3.3) is a metric developed by Levenshtein [21]. The algorithm is recursive and may be stated as
\[
\begin{align*}
\mathcal{M}(0, 0) &:= 0, \\
\mathcal{M}(i, 0) &:= \mathcal{M}(i - 1, 0) + \alpha(S_i[i], -): 1 \leq i \leq n, \\
\mathcal{M}(0, j) &:= \mathcal{M}(0, j - 1) + \alpha(-, S_j[j]): 1 \leq j \leq m, \\
\mathcal{M}(i, j) &:= \min \left\{ \mathcal{M}(i - 1, j) + \alpha(S_i[i], -), \mathcal{M}(i, j - 1) + \alpha(-, S_j[j]) : i \in [1, n], j \in [1, m] \right\}.
\end{align*}
\tag{3.7}
\]
The Equations (3.4), (3.5), (3.6) indicate the initial conditions. They also define the penalty of unpaired elements at the outset of the sequences \( S_1, S_2 \). Definition (3.7) states that all possible transformations (edit operations), that is to say deletions, insertions, and substitutions, have influence on the algorithm.
We have now described how the recursive algorithm, by means of the Equations (3.4), (3.5), (3.6) and (3.7), finds the optimal alignment of the sequences \( S_1, S_2 \). In order to construct our graph similarity measures, in the following we define the functions \( a^{\text{out}}, a^{\text{in}} \).
\[
a^{\text{out}}(v_{i_1, j_1}, v_{i_2, j_2}) := \begin{cases} 
\beta^{\text{out}}(\frac{\delta^{\text{out}}(v_{i_1, j_1}, v_{i_2, j_2}), \sigma^{1}_{\text{out}}}{1} + \infty & : i_1 = i_2, \\
\beta^{\text{out}}(\frac{\delta^{\text{out}}(v_{i_1, j_1}, v_{i_2, j_2}), \sigma^{2}_{\text{out}}}{1} & : \text{else},
\end{cases}
\tag{3.8}
\]
with \( 0 \leq i_k \leq h_k, 1 \leq j_k \leq \sigma_{ih} \), whereas \( \beta^{\text{out}}(x, y, \sigma^{k}_{\text{out}}) := 1 - e^{-\frac{(x - y)^2}{2(\sigma^{k}_{\text{out}})^2}}} \), \( x, y, \sigma^{k}_{\text{out}} \in \mathbb{R}, k \in \{1, 2\} \), and
\[
a^{\text{out}}(v_{i, j_1}, -) := \beta^{\text{out}}(\frac{\delta^{\text{out}}(v_{i, j_1}, -), \xi, \sigma^{2}_{\text{out}}}{1}), \\
a^{\text{out}}(-, v_{i, j_2}) := \beta^{\text{out}}(\frac{\xi, \delta^{\text{out}}(v_{i, j_2}), \sigma^{2}_{\text{out}}}{1}).
\tag{3.9}
\]
With \( \beta^{\text{in}}(x, y, \sigma^{k}_{\text{in}}) := 1 - e^{-\frac{(x - y)^2}{2(\sigma^{k}_{\text{in}})^2}}} \) we define straightforward
\[
a^{\text{in}}(v_{i_1, j_1}, v_{i_2, j_2}) := \begin{cases} 
\beta^{\text{in}}(\frac{\delta^{\text{in}}(v_{i_1, j_1}, v_{i_2, j_2}), \sigma^{1}_{\text{in}}}{1} + \infty & : i_1 = i_2, \\
\beta^{\text{in}}(\frac{\delta^{\text{in}}(v_{i_1, j_1}, v_{i_2, j_2}), \sigma^{2}_{\text{in}}}{1} & : \text{else},
\end{cases}
\tag{3.11}
\]
Given graphs \( \mathcal{G} \), we do not align vertices on different levels. To prevent this, we set the gap penalty to \(+\infty\), whereby our dynamic programming algorithm will never choose this cost-intensive path. In order to validate the alignments on each level of the given graphs \( \mathcal{H}_1, \mathcal{H}_2 \), we define the functions

\[
\gamma_{\text{out}}^k(\mathcal{H}_k, i) := \frac{\sum_{j=1}^{\sigma^k_i} \hat{\alpha}_{\text{out}}(v^k_{i,j}, \text{align}(v^k_{i,j}))}{\sigma^k_i}, \quad (3.14)
\]

\[
\gamma_{\text{in}}^k(\mathcal{H}_k, i) := \frac{\sum_{j=1}^{\sigma^k_i} \hat{\alpha}_{\text{in}}(v^k_{i,j}, \text{align}(v^k_{i,j}))}{\sigma^k_i}, \quad (3.15)
\]

\( k \in \{1, 2\} \). Once again, \( \sigma^k_i \) denotes the upper index of a vertex on level \( i \) related to \( \mathcal{H}_k \). To complete the Definitions (3.14), (3.15), we define the mapping align and \( \hat{\alpha}_{\text{out}}, \hat{\alpha}_{\text{in}} \) as follows:

\[
\text{align}(v^k_{i,j_1}) := \begin{cases} 
  v^k_{i,j_2} & : \text{align}^{-1}(v^k_{i,j_2}) = v^k_{i,j_1}, \\
  - & : \text{else},
\end{cases} \quad (3.16)
\]

\[
\hat{\alpha}_{\text{out}}(v^k_{i,j_1}, -) := \beta_{\text{out}}(\delta_{\text{out}}(v^k_{i,j_1}), \xi, \hat{\sigma}_{\text{out}}^1), \quad (3.17)
\]

\[
\hat{\alpha}_{\text{out}}(-, v^k_{i,j_2}) := \beta_{\text{out}}(\xi, \delta_{\text{out}}(v^k_{i,j_2}), \hat{\sigma}_{\text{out}}^1), \quad (3.18)
\]

\[
\hat{\alpha}_{\text{out}}(v^k_{i,j_1}, v^k_{i,j_2}) := \beta_{\text{out}}(\delta_{\text{out}}(v^k_{i,j_1}), \delta_{\text{out}}(v^k_{i,j_2}), \hat{\sigma}_{\text{out}}^2), \quad (3.19)
\]

\[
\hat{\alpha}_{\text{in}}(v^k_{i,j_1}, -) := \beta_{\text{in}}(\delta_{\text{in}}(v^k_{i,j_1}), \xi, \hat{\sigma}_{\text{in}}^1), \quad (3.20)
\]

\[
\hat{\alpha}_{\text{in}}(-, v^k_{i,j_2}) := \beta_{\text{in}}(\xi, \delta_{\text{in}}(v^k_{i,j_2}), \hat{\sigma}_{\text{in}}^1), \quad (3.21)
\]

\[
\hat{\alpha}_{\text{in}}(v^k_{i,j_1}, v^k_{i,j_2}) := \beta_{\text{in}}(\delta_{\text{in}}(v^k_{i,j_1}), \delta_{\text{out}}(v^k_{i,j_2}), \hat{\sigma}_{\text{in}}^2), \quad (3.22)
\]
Finally, if we set
\[
\gamma^{\text{out}}(i) := 1 - \\
\frac{1}{\sigma^1_i + \sigma^2_i} \cdot \left\{ \sum_{j=1}^{\sigma^1_i} \hat{\alpha}^{\text{out}}(v_{i,j}, \text{align} (v_{i,j})) \right\} \\
+ \frac{1}{\sigma^1_i + \sigma^2_i} \cdot \left\{ \sum_{j=1}^{\sigma^2_i} \hat{\alpha}^{\text{out}}(v_{i,j}, \text{align} (v_{i,j})) \right\},
\]
(3.23)

and
\[
\gamma^{\text{in}}(i) := 1 - \\
\frac{1}{\sigma^1_i + \sigma^2_i} \cdot \left\{ \sum_{j=1}^{\sigma^1_i} \hat{\alpha}^{\text{in}}(v_{i,j}, \text{align} (v_{i,j})) \right\} \\
+ \frac{1}{\sigma^1_i + \sigma^2_i} \cdot \left\{ \sum_{j=1}^{\sigma^2_i} \hat{\alpha}^{\text{in}}(v_{i,j}, \text{align} (v_{i,j})) \right\},
\]
(3.24)
we obtain measures which indicate how similar the out-degree and in-degree alignments of two sequences on a level \(i\) are. It holds by construction \(\gamma^{\text{out}}(i), \gamma^{\text{in}}(i) \in [0,1]\). In the following we construct a family of graph similarity measures.

Generally, similarity measures have wide applications in several areas of science, for example in cluster analysis [3], [9], sociology [22] and psychology [11], [30]. Now, our similarity measures will be based on the structural description of our objects. We define similarity measures which are similar to the definition of Batagelj [5].

**Definition 3.2.** Let \(U\) be a set of units and a mapping \(\phi: U \times U \rightarrow [0,1]\). \(\phi\) is called a similarity measure if
\[
\phi(u, v) = \phi(v, u), \forall u, v \in U \quad \text{(Symmetry)},
\]
(3.25)
and either
\[
\phi(u, u) \leq \phi(u, v), \forall u, v \in U \quad \text{(Forward)},
\]
(3.26)
or
\[
\phi(u, u) \geq \phi(u, v), \forall u, v \in U \quad \text{(Backward)}.
\]
(3.27)

Finally, we prove the key result for measuring the structural similarity of unlabeled, hierarchical, and directed graphs.
**Theorem 3.1.** Let $\hat{H}_1, \hat{H}_2$, $0 \leq i \leq \rho$, $\rho := \max(h_1, h_2)$.

\[
d_1(\hat{H}_1, \hat{H}_2) := \frac{\sum_{i=0}^{\rho} \lambda_i \cdot \gamma^{\text{fin}}(i)}{\sum_{i=0}^{\rho} \lambda_i},
\]

\[
d_2(\hat{H}_1, \hat{H}_2) := \frac{\sum_{i=0}^{\rho} \gamma^{\text{fin}}(i)}{\rho + 1},
\]

\[
d_3(\hat{H}_1, \hat{H}_2) := \frac{\prod_{i=0}^{\rho} \gamma^{\text{fin}}(i)}{d_2(\hat{H}_1, \hat{H}_2)},
\]

are a family $(d_i(\hat{H}_1, \hat{H}_2))_{1 \leq i \leq 3}$ of Backward similarity measures, where $\gamma^{\text{fin}}(i)$ is defined as

\[
\gamma^{\text{fin}}(i) := \zeta \cdot \gamma^{\text{out}}(i) + (1 - \zeta) \cdot \gamma^{\text{in}}(i).
\]

It holds $(d_i(\hat{H}_1, \hat{H}_2))_{1 \leq i \leq 3} \in [0, 1]$.

**Proof.** First, we consider the function

\[
\gamma^{\text{out}}(i) := 1 - \frac{1}{\sigma_1^2 + \sigma_i^2} \cdot \left\{ \sum_{j=1}^{\sigma_1} \hat{\alpha}^{\text{out}}(v_{i,j}^{\hat{H}_1}, \text{align} \left( v_{i,j}^{\hat{H}_2} \right)) \right\} + \frac{1}{\sigma_1^2 + \sigma_i^2} \cdot \left\{ \sum_{j=1}^{\sigma_2} \hat{\alpha}^{\text{out}}(v_{i,j}^{\hat{H}_2}, \text{align} \left( v_{i,j}^{\hat{H}_2} \right)) \right\}.
\]

With Equation (3.16), it follows that we have to distinguish the three cases for the function $\hat{\alpha}^{\text{out}}$: $\hat{\alpha}^{\text{out}}(v_{i,j_1}^{\hat{H}_1}, -)$, $\hat{\alpha}^{\text{out}}(-, v_{i,j_2}^{\hat{H}_2})$, $\hat{\alpha}^{\text{out}}(v_{i,j_1}^{\hat{H}_1}, v_{i,j_2}^{\hat{H}_2})$. We infer by the Equations (3.17), (3.18), (3.19) that

\[
\hat{\alpha}^{\text{out}}(v_{i,j_1}^{\hat{H}_1}, \text{align} \left( v_{i,j_1}^{\hat{H}_1} \right)) \leq 1 \quad \text{and} \quad \hat{\alpha}^{\text{out}}(v_{i,j_2}^{\hat{H}_2}, \text{align} \left( v_{i,j_2}^{\hat{H}_2} \right)) \leq 1.
\]

Hence, by Definition (3.23) we obtain $\gamma^{\text{out}}(i) \leq 1$. The Proof $\gamma^{\text{in}}(i) \leq 1$ is identical. Since

\[
\gamma^{\text{fin}}(i) \leq \zeta + (1 - \zeta) = 1,
\]

we obtain

\[
d_1(\hat{H}_1, \hat{H}_2) \leq \frac{\sum_{i=0}^{\rho} \lambda_i}{\sum_{i=0}^{\rho} \lambda_i} = 1.
\]
To prove that $d_1(\hat{H}_1, \hat{H}_2)$ is symmetric, we see that
\[
\gamma^{\text{out}}(i) := 1 - \frac{1}{\sigma_i^1 + \sigma_i^2} \cdot \left\{ \sigma_i^1 \sum_{j=1}^{\alpha} \hat{\gamma}_1^{\text{out}}(v_{i,j}, \text{align}(v_{i,j})) \right\} + \frac{1}{\sigma_i^1 + \sigma_i^2} \cdot \left\{ \sigma_i^2 \sum_{j=1}^{\alpha} \hat{\gamma}_2^{\text{out}}(v_{i,j}, \text{align}(v_{i,j})) \right\} = 1 - \frac{1}{\sigma_i^2 + \sigma_i^1} \cdot \left\{ \sigma_i^2 \sum_{j=1}^{\alpha} \hat{\gamma}_2^{\text{out}}(v_{i,j}, \text{align}(v_{i,j})) \right\} + \frac{1}{\sigma_i^2 + \sigma_i^1} \cdot \left\{ \sigma_i^1 \sum_{j=1}^{\alpha} \hat{\gamma}_1^{\text{out}}(v_{i,j}, \text{align}(v_{i,j})) \right\},
\]
and
\[
\gamma^{\text{in}}(i) := 1 - \frac{1}{\sigma_i^1 + \sigma_i^2} \cdot \left\{ \sigma_i^1 \sum_{j=1}^{\alpha} \hat{\gamma}_1^{\text{in}}(v_{i,j}, \text{align}(v_{i,j})) \right\} + \frac{1}{\sigma_i^1 + \sigma_i^2} \cdot \left\{ \sigma_i^2 \sum_{j=1}^{\alpha} \hat{\gamma}_2^{\text{in}}(v_{i,j}, \text{align}(v_{i,j})) \right\} = 1 - \frac{1}{\sigma_i^2 + \sigma_i^1} \cdot \left\{ \sigma_i^2 \sum_{j=1}^{\alpha} \hat{\gamma}_2^{\text{in}}(v_{i,j}, \text{align}(v_{i,j})) \right\} + \frac{1}{\sigma_i^2 + \sigma_i^1} \cdot \left\{ \sigma_i^1 \sum_{j=1}^{\alpha} \hat{\gamma}_1^{\text{in}}(v_{i,j}, \text{align}(v_{i,j})) \right\}.
\]
Therefore, we conclude with Equation (3.28) that
\[
d_1(\hat{H}_1, \hat{H}_2) = d_1(\hat{H}_2, \hat{H}_1).
\]
To finalize the proof for the Similarity Measure (3.28), we have to show that
\[
d_1(\hat{H}_1, \hat{H}_1) \geq d_1(\hat{H}_1, \hat{H}_2).
\]
If $\mathcal{H}_1 = \mathcal{H}_2$, then $\gamma^{\text{out}}(i) = 1$, $\gamma^{\text{in}}(i) = 1$ and $\gamma^{\text{fin}}(i) = 1$. Therefore we infer from Equation (3.28) that $d_1(\mathcal{H}_1, \mathcal{H}_1) = 1$ and see

$$d_1(\mathcal{H}_1, \mathcal{H}_1) = 1 \geq \frac{\sum_{i=0}^{\rho} \lambda_i \cdot \gamma^{\text{fin}}(i)}{\sum_{i=0}^{\rho} \lambda_i} = d_1(\mathcal{H}_1, \mathcal{H}_2),$$

the Backward property. In the case of Similarity Measure (3.29) we have nothing to prove, because if we set $1 = \lambda_0 = \lambda_1 = \cdots = \lambda_{\rho}$ in Equation (3.28), we obtain Equation (3.29). To prove the assertion of the theorem for Equation (3.30), we consider the well-known inequality [15]

$$(p_1 \cdot p_2 \cdots p_n)^{\frac{1}{n}} \leq \frac{p_1 + p_2 + \cdots + p_n}{n}, \quad p_i > 0, \quad 1 \leq i \leq n. \quad (3.33)$$

Since $\gamma^{\text{fin}}(i) \leq 1$, we can apply Inequality (3.33). We obtain

$$\gamma^{\text{fin}}(0) \cdot \gamma^{\text{fin}}(1) \cdots \gamma^{\text{fin}}(\rho) \leq \left[ \gamma^{\text{fin}}(0) \cdot \gamma^{\text{fin}}(1) \cdots \gamma^{\text{fin}}(\rho) \right]^{\frac{1}{\rho+1}} \leq \frac{\gamma^{\text{fin}}(0) + \gamma^{\text{fin}}(1) + \cdots + \gamma^{\text{fin}}(\rho)}{\rho + 1}$$

and especially

$$1 \geq \frac{\gamma^{\text{fin}}(0) \cdot \gamma^{\text{fin}}(1) \cdots \gamma^{\text{fin}}(\rho)}{\gamma^{\text{fin}}(0) + \gamma^{\text{fin}}(1) + \cdots + \gamma^{\text{fin}}(\rho)}. \quad (3.34)$$

The symmetry condition is clear, because the expression in the denominator of Equation (3.30) is a special case of $d_1$. The Backward condition follows immediately from Inequality (3.34),

$$1 = d_3(\mathcal{H}_1, \mathcal{H}_1) \geq \prod_{i=0}^{\rho} \frac{\gamma^{\text{fin}}(i)}{d_2(\mathcal{H}_1, \mathcal{H}_2)}.$$

\[\square\]

4. Experimental results

In this chapter, we report on the results obtained by testing the algorithm on a graph corpus $C_G$ representing 464 conference/workshop websites from computer science and mathematics. The graph corpus has already been used for a study [8], [23] in hypertext categorization.
In Figure 4.1 we depict the distributions of the ranked similarity values of Theorem (3.1) on the basis of different parameter spectra. Therefore, we express

**Definition 4.1.** In terms of $C_G$, we define data classes $D_1 - D_5$ which are manifested by the following parameter spectra:

1. $D_1$: $\zeta = 1.0$ (Solely alignments of Kernel-edges); If $\gamma^{fin}(i) < 0.5$, set $\lambda_i = 100$, else $\lambda_i = 1$; parameter settings:

   \[\sigma_{out}^1 = 1.0, \sigma_{out}^2 = 2.0, \sigma_{in}^1 = 1.0, \sigma_{in}^2 = 2.0, \hat{\sigma}_{out}^1 = 3.0, \hat{\sigma}_{out}^2 = 5.0, \]
   \[\hat{\sigma}_{in}^1 = 3.0, \hat{\sigma}_{in}^2 = 5.0.\]

2. $D_2$: $\zeta = 0.3$; If $\gamma^{fin}(i) < 0.5$, set $\lambda_i = 100$, else $\lambda_i = 1$; parameter settings:

   \[\sigma_{out}^1 = 1.0, \sigma_{out}^2 = 1.0, \sigma_{in}^1 = 1.0, \sigma_{in}^2 = 1.0, \hat{\sigma}_{out}^1 = 1.0, \hat{\sigma}_{out}^2 = 5.0, \]
   \[\hat{\sigma}_{in}^1 = 1.0, \hat{\sigma}_{in}^2 = 5.0.\]

3. $D_3$: $\zeta = 0.5$; If $\gamma^{fin}(i) < 0.5$, set $\lambda_i = 100$, else $\lambda_i = 1$; parameter settings:

   \[\sigma_{out}^1 = 1.0, \sigma_{out}^2 = 1.0, \sigma_{in}^1 = 1.0, \sigma_{in}^2 = 1.0, \hat{\sigma}_{out}^1 = 1.0, \hat{\sigma}_{out}^2 = 5.0, \]
   \[\hat{\sigma}_{in}^1 = 1.0, \hat{\sigma}_{in}^2 = 5.0.\]
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(4) \( D_4; \zeta = 0.5; \) If \( \gamma^{fin}(i) < 0.5 \), set \( \lambda_i = 64 \), else \( \lambda_i = 16 \); parameter settings:
\[
\sigma^1_{out} = 1.0, \sigma^2_{out} = 2.0, \sigma^1_{in} = 1.0, \sigma^2_{in} = 2.0, \sigma^1_{out} = 3.0, \sigma^2_{out} = 5.0, \\
\sigma^1_{in} = 3.0, \sigma^2_{in} = 5.0.
\]

(5) \( D_5; \zeta = 0.5; \) If \( \gamma^{fin}(i) < 0.5 \), set \( \lambda_i = 100 \), else \( \lambda_i = 1 \); parameter settings:
\[
\sigma^1_{out} = 1.0, \sigma^2_{out} = 1.0, \sigma^1_{in} = 1.0, \sigma^2_{in} = 1.0, \sigma^1_{out} = 3.0, \sigma^2_{out} = 5.0, \\
\sigma^1_{in} = 3.0, \sigma^2_{in} = 5.0.
\]

We regard the graph corpus \( C_G \) (see Table (4.3)) as homogeneous in the sense that it covers, structurally, the whole interval. Now, Figure (4.1) shows the ranked similarity values in terms of the data classes \( D_1 - D_5 \) representing the parameter spectra. Then, we observe from the plots of Figure (4.1) that the similarity measures \( d_1 \) and \( d_2 \) are not suitable. They do not cover the whole interval of the possible similarity values. If we set, in \( d_1 \), different values for \( \lambda_i \), \( 0 \leq i \leq \rho = \max(h_1, h_2) \), the measure \( d_1 \) is very sensitive to structural variances of the graphs. The “visible steps” and especially the horizontal segments in the plots of \( d_1 \) induce clusters, which contain the same number of graph pairs with a certain similarity value on the \( X \)-axis. In contrast to this, \( d_3 \) covers the interval of similarity values very well. It exploits the whole interval. Therefore, in the following we use for further experiments only \( d_3 \).

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{figure4.2.png}
\caption{The \( X \)-axis corresponds to the values of \( d_3 \in [0, 1] \) and the \( Y \)-axis represents the cumulative similarity distributions for \( D_1 - D_5 \).
\footnote{The first upper plot belongs to Item (1) of Definition (4.1), the second upper plot belongs to Item (2) of Definition (4.1) etc.}}
\end{figure}
Furthermore, we compute the cumulative similarity distribution (see Figure (4.2)), of $C_G$ based on Definition (4.1). In general, the computation of the cumulative similarity distribution of a graph corpus opens new perspectives: As a preprocessing step for the structural analysis of graph corpora we can decide, on the basis of the cumulative similarity distribution, how structurally different the graphs are.

The application of Figure (4.2) leads us directly, e.g., to the examination of the navigation strategies in terms of our chosen web-genre, that is conference/workshop websites from computer science and mathematics. Thereby, we assume that an unlabeled, hierarchical, and directed graph reflects all possible navigation paths of a graph-based conference website. The computation and interpretation of the cumulative similarity distribution of $C_G$ leads us to the question how different the navigation strategies within the specific web-genre are. In order to discuss the cumulative similarity distribution of $C_G$ (see Figure (4.2)); we note that the data classes $D_1$-$D_5$ are manifested by the same corpus $C_G$. We obtain a certain data class only by varying the parameters mentioned in Definition (4.1). Now, by varying the parameters we find the parameter set

$$\left( \zeta, \sigma^1_{out}, \sigma^2_{out}, \sigma^1_{in}, \sigma^2_{in}, \hat{\sigma}^1_{out}, \hat{\sigma}^2_{out}, \hat{\sigma}^1_{in}, \hat{\sigma}^2_{in} \right),$$

which captures enough structural information during the similarity measuring process. In the following we note that the plot of class $D_1$ differs in principle from the plots of data classes $D_2$-$D_5$. We recognize that, e.g., 20% of the graph-based conference websites have already the similarity value $d \leq 0.5$. Unlike 90% of the conference websites in $D_2$ have the similarity value $d \leq 0.5$. In summary, we conclude from Figure (4.2) that the similarity values of the conference websites in $D_1$ were significantly higher compared to the conference websites of data classes $D_2$-$D_5$. This is plausible, because the conference websites in $D_1$ are treated solely as rooted trees without Across-edges, Up-edges and Down-edges. Hence, the main part of the conference websites of $D_1$ is significantly less structurally different than the websites of the remaining data classes. In terms of $D_2$-$D_5$ the situation is inverted: In consideration of all types of conference websites the main part of the graph-based hypertext structures is structurally dissimilar on the basis of $d_1$. The plot of $D_4$ equals the plot of $D_5$. Finally, we note that for the data classes $D_2$-$D_5$ the main part of all possible navigation strategies is very different within our web-genre. This is reflected by psychological features of hypertext navigation.

5. Summary and conclusions

In this paper, we introduced a new method to measure the structural simi-
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Figure 4.3. Key data of the graph corpus.

<table>
<thead>
<tr>
<th>key data</th>
<th>value</th>
</tr>
</thead>
<tbody>
<tr>
<td>min(</td>
<td>V</td>
</tr>
<tr>
<td>max(</td>
<td>V</td>
</tr>
<tr>
<td>min(diam(H))</td>
<td>1</td>
</tr>
<tr>
<td>max(diam(H))</td>
<td>27</td>
</tr>
<tr>
<td>avg(</td>
<td>V</td>
</tr>
<tr>
<td>avg(diam(H))</td>
<td>3</td>
</tr>
</tbody>
</table>

larity of a special class of directed graphs. Thereby, the graphs are unlabeled, hierarchical, and directed and we applied our algorithm on a graph corpus $C_G$ [23] of graph-based hypertext structures. The main contributions of the paper are:

(1) Starting from observations on degree sequences of directed graphs, we developed a new method for measuring the structural similarity of graphs. The main idea of our new similarity measures are based on the derivation of property strings (out-degree and in-degree sequences on each graph level) for each hierarchical and directed graph and then to align the property strings representing our graphs by a dynamic programming technique. From the resulting alignment we obtain a value of the scoring function, which is minimized during the alignment process. The similarity of two hierarchical and directed graphs will be expressed by a cumulation of local similarity functions $\gamma^{out}(i), \gamma^{in}(i)$ which weighs two types of alignments: out-degree and in-degree alignments on a graph level. These alignments have both global and local significance. On the one hand, the sequence alignments will be implemented in a global sense, to compute the optimal alignment between the sequences $S_1$ and $S_2$. On the other hand, the alignments will be evaluated on the levels of the graphs by the function $\gamma^{fin}(i)$. Since the functions $\gamma^{out}(i), \gamma^{in}(i)$ and $\gamma^{fin}(i)$ are basically decoupled from the similarity measures of Theorem (3.1), we now can define new measured values $d_i$, where they are adapted to a new graph similarity problem. Therefore, we obtain a family of graph similarity measures. From the overall results mentioned above we see that our family of similarity measures $(d_i(\mathcal{H}_1, \mathcal{H}_2))_{1 \leq i \leq 3}$ is also different from graph similarity measures, which are based on isomorphic relations, e.g., [20], [28], [29], [32].

(2) In Section (4) we evaluated the measures $(d_i(\mathcal{H}_1, \mathcal{H}_2))_{1 \leq i \leq 3}$ on a graph corpus $C_G$, consisting of 464 graphs representing web-based hypertext
structures. The results of the evaluation are reflected in Figure (4.1), (4.2). Especially, we showed that the cumulative similarity distribution provides useful information about our graph corpus $C_G$. With Definition (4.1) we answered the important and interesting question how structurally different the graphs of $C_G$ are. Because our measures are parametric similarity measures depending on

$$(\zeta, \sigma^1_{\text{out}}, \sigma^2_{\text{out}}, \sigma^1_{\text{in}}, \sigma^2_{\text{in}}, \hat{r}^1_{\text{out}}, \hat{r}^2_{\text{out}}, \hat{r}^1_{\text{in}}, \hat{r}^2_{\text{in}}),$$

we were able to emphasize different structure types of our graph trees during the alignment process. For example, by setting $\zeta = 1$ we consider an unlabeled hypertext structure as a directed rooted tree. That means, in more detail, that we align only the out-degree property strings induced by edges from the underlying directed rooted tree. If we set $\zeta = 0$, we align the property strings induced by in-degree sequences only. In most of the cases we used $\zeta = \frac{1}{2}$, which weighs in- and out-degree sequences equally, $\gamma^{\text{fin}} = \frac{\gamma_{\text{out}}}{2} + \frac{\gamma_{\text{in}}}{2}$.

In the future we concentrate on a generalization of our new method for measuring the structural similarity of arbitrary graphs.

REFERENCES


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