# On the Complexity of Computing Probabilistic Bisimilarity

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**Abstract.** Probabilistic bisimilarity is a fundamental notion of equivalence on labelled Markov chains. It has a natural generalisation to a probabilistic bisimilarity pseudometric, whose definition involves the Kantorovich metric on probability distributions. The probabilistic bisimilarity pseudometric has discounted and undiscounted variants, according to whether one discounts the future in observing discrepancies between states.

This paper is concerned with the complexity of computing probabilistic bisimilarity and the probabilistic bisimilarity pseudometric on labelled Markov chains. We show that the problem of computing probabilistic bisimilarity is **P**-hard by reduction from the monotone circuit value problem. We also show that the discounted probabilistic bisimilarity pseudometric is rational and can be computed exactly in polynomial time using the network simplex algorithm and the continued fraction algorithm. In the undiscounted case we show that the probabilistic bisimilarity pseudometric is again rational and can be computed exactly in polynomial time using the ellipsoid algorithm.

# 1 Introduction

*Probabilistic bisimilarity* is a notion of equivalence for probabilistic labelled transition systems, introduced by Larsen and Skou [19]. It is based on Park and Milner's classical notion of bisimilarity for (non-deterministic) labelled transition systems [22]. A very similar and widely used concept on Markov chains, called *lumpability*, can be found as far back as the classical text of Kemeny and Snell [18]. A system and its probabilistic bisimilarity quotient can be considered indistinguishable, and quotienting by probabilistic bisimilarity is a widely used compression technique in verification and performance analysis [15, 17, 20].

The first part of this paper concerns the complexity of computing probabilistic bisimilarity. It is known that this can be done in polynomial time, e.g., by partition refinement [2, 10, 30]. Our first result shows that probabilistic bisimilarity is **P**-hard, and therefore **P**-complete. As a consequence probabilistic bisimilarity is not in **NC** unless  $\mathbf{P} = \mathbf{NC}$ . (Recall that **NC** is a subclass of **P** comprising problems that can be solved in polylogarithmic time using PRAMs

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of polynomial size [13]. Informally such problems are considered to be efficiently parallelisable.) By contrast, language equivalence of probabilistic automata is in **NC** [29], as are related equivalence problems such as tree isomorphism [13].

For (non-deterministic) labelled transition systems it is known that computing bisimilarity is **P**-complete [3, 24]. However the proof in the probabilistic case requires a different construction than in *op. cit.* 

For probabilistic systems it is natural to generalise from bivalent notions of equivalence, such as probabilistic bisimilarity or language equivalence [28], to quantitative measures of similarity. As well as being more informative, such measures are more meaningful in the presence of rounding errors in computation and modelling.

In the second part of this paper we consider a *probabilistic bisimilarity pseu*dometric on labelled Markov chains. This generalises the notion of probabilistic bisimilarity by assigning a *similarity distance* to pairs of states of a labelled Markov chain. The smaller the distance, the more alike the states, with states at zero distance if and only if they are probabilistic bisimilar. This pseudometric was first introduced in [11] and, together with closely related notions, has subsequently been studied in the context of systems biology [26], games [8], planning [9] and security [7], among others. The definition of the pseudometric is based on the classical Kantorovich metric on probability distributions. The pseudometric has discounted versions, which discount the future in observing discrepancies between states, and an undiscounted version, which does not discount the future.

We show that for labelled Markov chains with rational transition probabilities the discounted probabilistic bisimilarity pseudometric is rational and can be computed exactly by a polynomial-time algorithm. In particular, we show that the distances can be approximated by using the network simplex algorithm repeatedly and the exact distances can be obtained from the approximated ones by means of the continued fraction algorithm. In the undiscounted case we also obtain a polynomial-time algorithm to exactly compute the pseudometric, this time using the heavier machinery of the ellipsoid algorithm. These results go beyond previous work which only showed how to approximate the pseudometric up to some desired level of precision [6]. In the undiscounted case it was only known how to approximate the pseudometric using polynomial space [5].

In combination with our lower bound on computing probabilistic bisimilarity we conclude that computing the probabilistic bisimilarity pseudometric is  $\mathbf{P}$ -complete. By contrast, a generalisation of the probabilistic bisimilarity pseudometric from labelled Markov chains to stochastic games has been shown to be as hard as the sum-of-square-roots problem [8], a problem not known even to be in  $\mathbf{NP}$ .

### 2 Probabilistic Bisimilarity

In this section we introduce labelled Markov chains and probabilistic bisimilarity. The main result of this section is that computing probabilistic bisimilarity is **P**-hard.

A labelled Markov chain is a tuple  $\mathcal{M} = (S, \Sigma, \pi, \ell)$  consisting of a finite set of states S, a finite set of labels  $\Sigma$ , a rational transition matrix  $\pi$  such that  $\sum_{t \in S} \pi_{s,t} = 1$  for all  $s \in S$ , and a labelling function  $\ell : S \to \Sigma$ .

A probabilistic bisimulation on  $\mathcal{M}$  is an equivalence relation  $R \subseteq S \times S$  such that if  $s \ R \ t$  then  $\ell(s) = \ell(t)$  and

$$\sum_{u \in E} \pi_{s,u} = \sum_{u \in E} \pi_{t,u}$$

for each *R*-equivalence class *E*, i.e., related states have the same label and the same probability to transition into any given equivalence class. It is a standard result that there is a largest probabilistic bisimulation on  $\mathcal{M}$  and that this relation is an equivalence relation (see, e.g., [23, Section 7.6]). The maximum probabilistic bisimulation is called *probabilistic bisimilarity* and is denoted  $\sim$ . From now on, we mostly refer to probabilistic bisimilarity as simply bisimilarity.

We are interested in the problem of computing bisimilarity  $\sim$  on  $\mathcal{M}$ . The decision version of the problem asks whether  $s \sim t$  for two designated states  $s, t \in S$ .

The above formulation of the bisimilarity problem is convenient for our hardness proof, however variations, such as replacing state labels with labels on transitions, can easily be accommodated. It is also not difficult to reduce the problem above to the restricted case in which the set of labels has two elements.

For a state s, let  $Succ(s) = \{u : \pi_{s,u} > 0\}$ . We say that a transition matrix  $\pi$  is *uniform* if for all  $s \in S$  and  $u, v \in Succ(s), \pi_{s,u} = \pi_{s,v}$ . That is, the transition probability out of each state is a uniform distribution over its support.

**Lemma 1 (Matching Lemma).** Assume that  $\pi$  is uniform. Suppose that  $s, t \in S$  are such that |Succ(s)| = |Succ(t)|. Then  $s \sim t$  if and only if  $\ell(s) = \ell(t)$  and there exists a bijection  $f: Succ(s) \to Succ(t)$  with  $u \sim f(u)$  for each  $u \in Succ(s)$ .

*Proof.* Suppose that  $s \sim t$ . Since  $\sim$  is a bisimulation,  $\ell(s) = \ell(t)$  and for each  $\sim$ -equivalence class E,

$$\sum_{x \in E} \pi_{s,x} = \sum_{x \in E} \pi_{t,x}.$$
(1)

Since  $\pi$  is uniform,  $|E \cap Succ(s)| = |E \cap Succ(t)|$  for each  $\sim$ -equivalence class *E*. Hence there exists a bijection  $f : Succ(s) \to Succ(t)$  with  $u \sim f(u)$  for all  $u \in Succ(s)$ .

Conversely, assume that  $\ell(s) = \ell(t)$  and suppose that f is a bijection as above. To conclude that  $s \sim t$  we prove that the smallest equivalence relation containing  $\sim \cup \{(s,t)\}$ , which we denote by R, is a bisimulation.

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Since  $\sim$  is a bisimulation and  $\ell(s) = \ell(t)$ , R only relates states with the same label. Moreover, since every R-equivalence class is a union of  $\sim$ -equivalence classes, it suffices to show (1) for  $\sim$ -equivalences classes only.

Assume u R v. We distinguish three cases. First, let u = s and v = t. Because of the existence of the bijection f, we have that  $|E \cap Succ(s)| = |E \cap Succ(t)|$ for each  $\sim$ -equivalence class E. Because  $\pi$  is uniform, (1) holds for each  $\sim$ equivalence class E. Second, let  $u \sim s$  and  $v \sim t$ . Recall that  $\sim$  is a bisimulation. Hence for each  $\sim$ -equivalence class E,

$$\sum_{x \in E} \pi_{u,x} = \sum_{x \in E} \pi_{s,x} \quad [u \sim s]$$
$$= \sum_{x \in E} \pi_{t,x} \quad [\text{by the previous case}]$$
$$= \sum_{x \in E} \pi_{v,x} \quad [v \sim t]$$

The third and final case,  $u \sim v$ , follows immediately from the fact that  $\sim$  is a bisimulation.

#### Theorem 2. Deciding probabilistic bisimilarity is P-hard.

*Proof.* We reduce from the MONOTONE CIRCUIT VALUE problem which is **P**-hard [13, Theorem 6.2.2]. Recall that a monotone circuit is a finite directed acyclic graph in which nodes have in-degree either two or zero. Nodes with indegree two are labelled  $\land$  or  $\lor$ ; nodes with in-degree zero, called input nodes, are labelled either true (1) or false (0). There is a distinguished output node with out-degree zero. The MONOTONE CIRCUIT VALUE problem is to compute the output of a given monotone circuit.

Given a circuit C, we define a Markov chain  $\mathcal{M}(C)$  with a uniform transition matrix. For each node  $n_i$  of C and its incoming edges, we include a gadget consisting of states and their outgoing transitions in  $\mathcal{M}(C)$ . Note that the transitions of the Markov chain go in the opposite direction of the edges of the circuit. Each gadget contains states  $u_i$  and  $v_i$ . We will prove that  $u_i \sim v_i$  if and only if  $n_i$  evaluates to true. We define the labelling function  $\ell$  such that states have the same label if and only if they belong to the same gadget and the gadget does not represent an input node that is labelled false. In the diagrams below, states have the same label if and only if they have the same index and the same colour.

We describe  $\mathcal{M}(C)$  by giving gadgets for each input node, and-gate and or-gate of C.

The gadget for input node labelled true is shown below.



The gadget for input node labelled false is shown below.

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Note that  $u_i$  and  $v_i$  have the same label if and only if  $n_i$  is labelled true and therefore  $u_i \sim v_i$  if and only if  $n_i$  is labelled true.

The gadget for an *and*-gate is shown below.



Note that  $u_j$ ,  $v_j$  and  $u_k$ ,  $v_k$  are states of the gadgets corresponding to the nodes  $n_j$  and  $n_k$ . The correctness of this gadget amounts to showing that  $u_i \sim v_i$  if and only if both  $u_j \sim v_j$  and  $u_k \sim v_k$ . This follows immediately from the Matching Lemma and the fact that the definition of  $\ell$  precludes  $u_j \sim v_k$  and  $v_j \sim u_k$  in case  $n_j$  and  $n_k$  are different nodes. If  $n_j$  and  $n_k$  are one and the same node, the *and*-gate can be removed from the circuit.

The gadget for an *or*-gate is shown below.



The correctness of this gadget amounts to showing that  $u_i \sim v_i$  if and only if  $u_j \sim v_j$  or  $u_k \sim v_k$ .

 $\begin{array}{l} u_i \sim v_i \text{ iff } (w_i \sim y_i \wedge x_i \sim z_i) \lor (w_i \sim z_i \wedge x_i \sim y_i) \quad \text{[Matching Lemma]} \\ \text{ ff } u_i \sim v_i \lor u_k \sim v_k \quad \text{[Matching Lemma]} \end{array}$ 

In the last step we use again the fact that the definition of  $\ell$  precludes  $u_j \sim v_k$ and  $v_j \sim u_k$  in case  $n_j$  and  $n_k$  are different nodes. If  $n_j$  and  $n_k$  are one and the same node, the *or*-gate can be removed from the circuit.

This completes the description of the gadgets. The Markov chain  $\mathcal{M}(C)$  is obtained by composing the gadgets for each node of C. The transduction of a circuit to a Markov chain is done gate by gate. To produce the output gadget corresponding to each circuit gate one only needs to store the indices of the gate and its two inputs, and the states of the output gadget. Thus the reduction can be done in deterministic logarithmic space.  $\hfill \Box$ 

The proofs of **P**-hardness of ordinary bisimilarity for labelled transition systems by Balcázar, Gabarró and Sántha [3] and Sawa and Jančar [24] are also by reduction from MONOTONE CIRCUIT VALUE. However in the probabilistic case disjunction cannot be translated directly as in the non-deterministic case. Interestingly a formally identical gadget to the above disjunction gadget appears in Toran's proof of **DET**-hardness of graph isomorphism [27]. However **DET** is a subclass of **P** and the graph isomorphism problem is not known to be **P**-hard.

### 3 The Bisimilarity Pseudometric

In this section we recall the definition of a bisimilarity pseudometric on labelled Markov chains. We first give a logical characterisation, due to Desharnais, Gupta, Jagadeesan and Panangaden [11], based on a real-valued semantics for Larsen and Skou's probabilistic modal logic [19]. This characterisation illustrates the sense in which states that are close in the pseudometric satisfy similar behavioural properties. In the next section we give a more abstract fixed-point characterisation of the pseudometric, which will be used in our algorithms.

The logic  $\mathcal{L}$  is defined by the grammar

$$\varphi ::= \sigma \mid \varphi \lor \varphi \mid \neg \varphi \mid \varphi \ominus q \mid \diamond \varphi \tag{2}$$

where  $\sigma \in \Sigma$  and  $q \in [0, 1]$  is rational.

We consider a real-valued semantics of  $\mathcal{L}$ , which is parameterised by a discount factor  $c \in (0, 1]$ . The smaller the value of c, the more the future is discounted, with c = 1 being the *undiscounted* case. Given a labelled Markov chain  $\mathcal{M} = (S, \Sigma, \pi, \ell)$ , the interpretation of a formula  $\varphi$  is a function  $\llbracket \varphi \rrbracket : S \to [0, 1]$ defined by the following clauses:

$$\llbracket \sigma \rrbracket(s) = \begin{cases} 1 \text{ if } \ell(s) = \sigma \\ 0 \text{ otherwise} \end{cases}$$
$$\llbracket \varphi \lor \psi \rrbracket(s) = \max(\llbracket \varphi \rrbracket(s), \llbracket \psi \rrbracket(s)) \\ \llbracket \neg \varphi \rrbracket(s) = 1 - \llbracket \varphi \rrbracket(s) \\ \llbracket \varphi \ominus q \rrbracket(s) = \max(\llbracket \varphi \rrbracket(s) - q, 0) \\ \llbracket \diamond \varphi \rrbracket(s) = c \cdot \sum_{t \in S} \pi_{s,t} \cdot \llbracket \varphi \rrbracket(t) \end{cases}$$

A pseudometric is a relaxation of the notion of an ordinary metric in which different states can have distance zero. Formally a (1-bounded) pseudometric on a set S is a map  $d: S \times S \rightarrow [0, 1]$  such that for all  $s, t, u \in S$ :

1. d(s,s) = 02. d(s,t) = d(t,s)3.  $d(s,u) \le d(s,t) + d(t,u)$ . Given a discount factor  $c \in (0, 1]$  the function  $\mathsf{d}_c : S \times S \to [0, 1]$  assigns a distance to every pair of states of a labelled Markov chain according to the following definition:

$$\mathsf{d}_{c}(s,t) = \sup_{\varphi \in \mathcal{L}} \left| \llbracket \varphi \rrbracket(s) - \llbracket \varphi \rrbracket(t) \right|.$$
(3)

It is straightforward that, with this definition,  $d_c$  is a pseudometric. The following theorem justifies our description of  $d_c$  as a bisimilarity pseudometric.

**Theorem 3.** [23, Section 8.2]  $d_c(s,t) = 0$  if and only if  $s \sim t$ .

In [8], Chatterjee, de Alfaro, Majumdar and Raman enriched the logic  $\mathcal{L}$ by the addition of fixed-point operators, yielding a quantitative  $\mu$ -calculus  $\mathcal{L}_{\mu}$ which can express reachability and  $\omega$ -regular specifications. For example, the  $\mathcal{L}_{\mu}$ -formula  $\mu x.(\sigma \lor \diamond x)$  represents the probability to reach a  $\sigma$ -labelled state, while  $\nu y.\mu x.((\sigma \land y) \lor \diamond x)$  represents the probability to infinitely often visit a  $\sigma$ -labelled state. It is shown in [8] that  $\mathsf{d}_c(s,t) = \sup_{\varphi \in \mathcal{L}_{\mu}} |\llbracket \varphi \rrbracket(s) - \llbracket \varphi \rrbracket(t)|$  for any pair of states  $s, t \in S$ ; thus  $\mathsf{d}_c$  can equivalently be defined in terms of the more powerful logic  $\mathcal{L}_{\mu}$ .

In the appendix we show that our bisimilarity pseudometric is an upperbound for the variational distance between trace distributions. This can be seen as a quantitative version of the folklore that bisimilar states satisfy the same lineartime properties. Whereas the variational distance between trace distributions is **NP**-hard to compute as shown by Lyngsø and Pedersen in [21], we will show that our bisimilarity pseudometric can be computed in polynomial time.

### 4 Matchings and the Kantorovich Metric

The characterisation of bisimilarity in Section 2 only admits equivalence relations as bisimulations. The following more general account of bisimilarity is useful both in the technical development below and to motivate the Kantorovich metric.

Say that a probability distribution  $\omega$  on  $S \times S$  is a *matching* of probability distributions  $\mu, \nu$  on S if

$$\begin{split} \sum_{v \in S} \omega(u, v) &= \mu(u) \quad \text{ for all } u \in S \\ \sum_{u \in S} \omega(u, v) &= \nu(v) \quad \text{ for all } v \in S \,. \end{split}$$

In other words,  $\omega$  is a joint probability distribution whose marginals are  $\mu$  and  $\nu$ .

**Proposition 4.** [16, Theorem 4.6] Let  $\mathcal{M} = (S, \Sigma, \pi, \ell)$  be a labelled Markov chain. Then  $s \sim t$  if and only if there exists a relation  $R \subseteq S \times S$  containing (s,t) such that (i)  $\ell(u) = \ell(v)$  for all  $(u,v) \in R$ ; (ii) for each  $(u,v) \in R$  there exists a probability distribution  $\omega$  on  $S \times S$  that is a matching of  $\pi_{u,-}$  and  $\pi_{v,-}$  and whose support is contained in R.

Suppose that (S, d) is a finite metric space. The Kantorovich metric  $d_K$  on the set of probability distributions on S is defined by

$$d_K(\mu,\nu) = \min_{\omega \in \Omega_{\mu,\nu}} \sum_{u,v \in S} d(u,v) \cdot \omega(u,v) \,,$$

where  $\Omega_{\mu,\nu}$  is the set of matchings of  $\mu$  and  $\nu$ .

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Informally we can characterise the bisimilarity pseudometric  $\mathsf{d}_c(s,t)$  as the distance between the distributions  $\pi_{s,-}$  and  $\pi_{t,-}$  in the Kantorovich metric over  $(S,\mathsf{d}_c)$ . This characterisation is recursive, and accordingly we will show that  $\mathsf{d}_c$  is a fixed point of a functional  $\Delta_c$  based on the Kantorovich metric.

Define  $\Delta_c : [0,1]^{S \times S} \to [0,1]^{S \times S}$  as follows. If  $\ell(s) \neq \ell(t)$  then  $\Delta_c(d)(s,t) = 1$  and if  $\ell(s) = \ell(t)$  then

$$\Delta_c(d)(s,t) = c \cdot \min_{\omega \in \Omega_{s,t}} \sum_{u,v \in S} d(u,v) \cdot \omega(u,v), \qquad (4)$$

where  $\Omega_{s,t}$  is the set of matchings of  $\pi_{s,-}$  and  $\pi_{t,-}$ .

The set  $[0,1]^{S\times S}$  is a complete lattice in the pointwise order. It is shown in [4, Proposition 38] that  $\Delta_c$  is a monotone selfmap on  $[0,1]^{S\times S}$  and thus, by Tarski's fixed point theorem, has a least fixed point. This turns out to be the pseudometric  $\mathsf{d}_c$ .

#### **Theorem 5.** [5, Theorem 4.6] $d_c$ is the least fixed point of $\Delta_c$ .

*Remark 6.* In the relational setting it is traditional to view bisimilarity as a greatest fixed point. Intuitively the situation is opposite in the pseudometric setting because the bottom element of [0, 1] represents relatedness.

**Theorem 7.** If c < 1 then  $d_c$  is the unique fixed point of  $\Delta_c$ .

*Proof sketch.* We can show that  $\Delta_c$  is *c*-Lipschitz. From Banach's fixed point theorem we can conclude that the fixed point is unique.

However,  $\Delta_1$  need not have a unique fixed point. For example, consider the labelled Markov chain with a single state. Then  $\Delta_1$  is the identity mapping.

*Example 8.* Consider the Markov chain below, where  $\ell(s) = \ell(t) \neq \ell(u)$ :

$$\bigcap_{s}^{1} \qquad \bigcap_{t \xrightarrow{1-c} u}^{c} \bigcap_{t \xrightarrow{1-c} u}^{1}$$

For c < 1,  $\mathbf{d}_c(s,t) = \frac{c-c^2}{1-c^2}$ . The pseudometric **0** assigns to every pair of states distance zero. For all  $n \in \mathbb{N}$ ,  $\Delta_c^n(\mathbf{0})(s,t) \leq \frac{c-c^{2n+1}}{1+c}$ . This shows that the fixed point may not simply be reached by a finite number of iterations of  $\Delta_c$ .

### 5 Algorithms for Bisimilarity Pseudometrics

#### 5.1 The Discounted Case

Let c < 1 be a fixed rational discount factor. Given a labelled Markov chain  $\mathcal{M}$ , we show that  $\mathsf{d}_c$  is rational and can be computed exactly in time polynomial in  $\operatorname{size}(\mathcal{M})$  and  $\operatorname{size}(c)$ .<sup>3</sup>

<sup>&</sup>lt;sup>3</sup> We denote by size(X) the size of the representation of an object X. We represent rational numbers as quotients of integers written in binary. For example, the size of a rational number is the sum of the bit lengths of its numerator and denominator and the size of a matrix is the sum of the sizes of its entries.

In Theorem 5 we have characterised  $\mathsf{d}_c$  as the least fixed point of  $\Delta_c$ . While the stipulation that  $\mathsf{d}_c$  be the least fixed point is essential in the undiscounted case, it is redundant in the discounted case. In the latter case,  $\Delta_c$  has a unique fixed point (see Theorem 7). As a consequence,  $\mathsf{d}_c$  is also the greatest fixed point of  $\Delta_c$  for c < 1. Thus, by Tarski's fixed point theorem, we have

$$\mathsf{d}_c = \bigsqcup \{ d \in \mathbb{R}^{S \times S} : d \le \Delta_c(d) \land 0 \le d \le 1 \}.$$
(5)

This simple change in perspective is fruitful because the characterisation (5) directly yields a translation of the problem of computing  $d_c$  to the following linear program:

maximise 
$$\sum_{s,t\in S} d_{s,t}$$
such that  $d_{s,t} \leq c \cdot \sum_{u,v\in S} d_{u,v} \cdot \omega_{u,v}$   $\omega \in \Omega_{s,t}, \ \ell(s) \neq \ell(t)$   
 $d_{s,t} = 1$   
 $0 \leq d_{s,t} \leq 1$ 

$$(6)$$

As we will see, the linear program (6) can be solved in polynomial time using the ellipsoid algorithm. We pursue this option in the undiscounted setting below. However, here we do not require such powerful techniques. Instead we just use the linear programming formulation to observe that the fixed point of  $\Delta_c$  is rational and bounded in size by a polynomial in size( $\mathcal{M}$ ) and size(c). We then approximate the fixed point by repeating the network simplex algorithm, obtaining the exact solution by rounding by means of the continued fraction algorithm.

Recall that the set of matchings  $\Omega_{s,t}$  is a polytope in  $\mathbb{R}^{S \times S}$  defined by the following constraints:

$$\sum_{\substack{v \in S \\ \sum_{u \in S} \omega_{u,v} = \pi_{t,v} \\ \omega_{u,v} \ge 0}} \sum_{\substack{u \in S \\ \omega_{u,v} \ge 0}} \omega_{u,v} = \pi_{t,v}$$
(7)

In general,  $\Omega_{s,t}$  is infinite and therefore the set of constraints in (6) is infinite also. However, for each fixed *d* the linear function mapping a matching  $\omega$  to  $c \cdot \sum_{u,v} d(u,v) \cdot \omega(u,v)$  achieves its minimum on  $\Omega_{s,t}$  at some vertex. Thus, writing  $V(\Omega_{s,t})$  for the (finite) set of vertices of  $\Omega_{s,t}$ , we can reformulate (6) as the following linear program, which has the same feasible region:

maximise 
$$\sum_{s,t\in S} d_{s,t}$$
  
such that  $d_{s,t} \leq c \cdot \sum_{u,v\in S} d_{u,v} \cdot \omega_{u,v}$   $\omega \in V(\Omega_{s,t}), \ \ell(s) \neq \ell(t)$   
 $d_{s,t} = 1$   
 $0 \leq d_{s,t} \leq 1$ 

$$(8)$$

We denote the polytope defined by the set of constraints of (8) by D. To prove that the distances are rational, we first observe the following.

**Proposition 9.** Each  $\omega \in V(\Omega_{s,t})$  is rational of size polynomial in size( $\mathcal{M}$ ).

Proof sketch. Since a vertex of  $\Omega_{s,t}$  is by definition an intersection of hyperplanes given by the (in)equalities defining  $\Omega_{s,t}$  and the coefficients of the (in)equalities are rationals bounded in size by size( $\mathcal{M}$ ), we can conclude that each  $\omega \in V(\Omega_{s,t})$ is rational of size polynomial in size( $\mathcal{M}$ ).

### **Proposition 10.** $d_c$ is rational of size polynomial in size( $\mathcal{M}$ ) and size(c).

*Proof sketch.* Along a similar line of reasoning as used in the proof of Proposition 9, we can conclude that the vertices of polytope D are rational of size polynomial in size( $\mathcal{M}$ ) and size(c).

Since the function mapping any d of the polytope D to  $\sum_{s,t\in S} d(s,t)$  is linear, it attains its maximum,  $\mathsf{d}_c$ , at some vertex of D, which, as we have just shown, is rational of size polynomial in size( $\mathcal{M}$ ) and size(c).

Note that the proofs of Proposition 9 and 10 are also valid for c = 1 and, hence,  $d_1$  is rational as well. Having established that  $d_c$  is rational, we now give a simple iterative algorithm to approximate  $d_c$  starting from the pseudometric **0**.

**Proposition 11.** For all  $n \in \mathbb{N}$ ,  $\Delta_c^n(\mathbf{0})$  is rational of size polynomial in size( $\mathcal{M}$ ) and size(c) and can be computed in time polynomial in size( $\mathcal{M}$ ) and size(c).

*Proof sketch.* We prove this property by induction on n. Obviously, the property holds for n = 0. Let n > 0. Obviously, the property holds when  $\ell(s) \neq \ell(t)$ . Otherwise,

$$\Delta_c^n(\mathbf{0})(s,t) = c \cdot \min_{\omega \in \Omega_{s,t}} \sum_{u,v \in S} \Delta_c^{n-1}(\mathbf{0})(u,v) \cdot \omega(u,v).$$

The above minimum is attained at a vertex of  $\Omega_{s,t}$ . As we have seen in Proposition 9, these vertices are rationals of size polynomial in size( $\mathcal{M}$ ). Furthermore, by induction,  $\Delta_c^{n-1}(\mathbf{0})$  is rational of size polynomial in size( $\mathcal{M}$ ) and size(c). Hence,  $\Delta_c^n(\mathbf{0})(s,t)$  is a rational of size polynomial in size( $\mathcal{M}$ ) and size(c). Computing  $\Delta_c(d)(s,t)$  is a minimum-cost flow problem for which there are versions of the network simplex algorithm that are strongly polynomial time [1].

To get  $\epsilon$ -close to  $\mathsf{d}_c$ , we need to iterate  $\lceil \log_c(\epsilon) \rceil$  times.

**Proposition 12.** For all  $\epsilon > 0$ ,  $\|\Delta_c^{\lceil \log_c(\epsilon) \rceil}(\mathbf{0}) - \mathsf{d}_c\| \leq \epsilon$ .

*Proof.* It suffices to show that for all  $n \in \mathbb{N}$ ,  $\|\Delta_c^n(\mathbf{0}) - \mathsf{d}_c\| \leq c^n$  by induction on n. Obviously, the property holds for n = 0. Let n > 0. Then

$$\begin{aligned} \|\Delta_c^n(\mathbf{0}) - \mathsf{d}_c\| &= \|\Delta_c(\Delta_c^{n-1}(\mathbf{0})) - \Delta_c(\mathsf{d}_c)\| \quad \text{[Theorem 7]} \\ &\leq c \cdot \|\Delta_c^{n-1}(\mathbf{0}) - \mathsf{d}_c\| \quad [\Delta_c \text{ is } c\text{-Lipschitz}] \\ &\leq c \cdot c^{n-1} \quad \text{[induction hypothesis]} \end{aligned}$$

¿From Proposition 11 and 12 we can conclude that we can approximate  $\mathsf{d}_c$  in time polynomial in  $\operatorname{size}(\mathcal{M})$ ,  $\operatorname{size}(c)$  and  $\log_2(\frac{1}{\epsilon})$ . Once we have iterated close enough to  $\mathsf{d}_c$ , we can use the continued fraction algorithm (see, e.g. [14, Section 5.1]) to obtain  $\mathsf{d}_c$ .

**Theorem 13.** The pseudometric  $d_c$  can be computed in time polynomial in  $size(\mathcal{M})$  and size(c).

*Proof sketch.* This follows now immediately from the observation made by Etessami and Yannakakis [12, page 2540] that for problems whose solutions are rational, of size polynomial in the input size, if we can solve the approximation problem in polynomial time, then we can also solve the exact computation problem in polynomial time by using the continued fraction algorithm.

### 5.2 The Undiscounted Case

Throughout this section we refer to the undiscounted bisimilarity pseudometric as d, rather than  $d_1$  and likewise use  $\Delta$  instead of  $\Delta_1$ .

In the previous section we gave a reduction of the problem of computing  $d_c$  to linear programming by characterising  $d_c$  as the greatest fixed point of  $\Delta_c$  for c < 1. However, recall from Section 4 that d is not in general the greatest fixed point of  $\Delta$ . Nevertheless we can recover a greatest-fixed-point characterisation of d by separately handling the set of bisimilar states, i.e. the states at distance zero. This will allow us to use linear programming to compute d.

As a first step we define  $\Delta' : [0,1]^{S \times S} \to [0,1]^{S \times S}$  by

$$\Delta'(d)(s,t) = \begin{cases} \Delta(d)(s,t) & \text{if } s \not\sim t \\ 0 & \text{if } s \sim t. \end{cases}$$

**Proposition 14.**  $\Delta'$  has a unique fixed point.

*Proof.* Since  $\Delta$  is monotone, we can easily deduce that  $\Delta'$  is monotone as well. According to Tarski's fixed point theorem,  $\Delta'$  has a least and a greatest fixed point. Hence, it is sufficient to prove that if  $d \leq d'$  are both fixed points of  $\Delta'$  then d = d'.

To this end, let

$$m = \max\{ d'(s,t) - d(s,t) : s, t \in S \},\$$

and let

$$R = \{ (s,t) \in S \times S : d'(s,t) - d(s,t) = m \}$$

be the set of pairs which maximise the discrepancy between d' and d. We will show that m = 0, which implies that d = d'. We distinguish two cases.

Assume that  $(s,t) \in R$  such that  $\ell(s) \neq \ell(t)$ . Then

$$d'(s,t) - d(s,t) = \Delta'(d')(s,t) - \Delta'(d)(s,t) = 1 - 1 = 0$$

and, hence, m = 0.

Otherwise, for all  $(s,t) \in R$  we have that  $\ell(s) = \ell(t)$ . In this case, we claim that  $R \subseteq \sim$ . From this claim it follows that for all  $(s,t) \in R$ ,

$$d'(s,t) - d(s,t) = \Delta'(d')(s,t) - \Delta'(d)(s,t) = 0 - 0 = 0$$

and, hence, m = 0. It just remains to prove the claim.

According to Proposition 4, it suffices to show that if  $(s,t) \in R$  then there exists  $\omega \in \Omega_{s,t}$  such that  $\omega(u,v) > 0$  implies  $(u,v) \in R$ , that is, d'(u,v) - d(u,v) = m.

Suppose  $\varDelta'(d)(s,t) = \sum_{u,v \in S} d(u,v) \cdot \omega(u,v),$  where  $\omega \in \varOmega_{s,t}.$  Then

$$\begin{split} m &= d'(s,t) - d(s,t) \\ &= \Delta'(d')(s,t) - \Delta'(d)(s,t) \\ &= \left( \min_{\omega' \in \Omega_{s,t}} \sum_{u,v \in S} d'(u,v) \cdot \omega'(u,v) \right) - \sum_{u,v \in S} d(u,v) \cdot \omega(u,v) \\ &\leq \sum_{u,v \in S} d'(u,v) \cdot \omega(u,v) - \sum_{u,v \in S} d(u,v) \cdot \omega(u,v) \\ &= \sum_{u,v \in S} (d'(u,v) - d(u,v)) \cdot \omega(u,v). \end{split}$$

Since  $d'(u,v) - d(u,v) \le m$  and  $\sum_{u,v\in S} \omega(u,v) = 1$ , we can conclude from  $\sum_{u,v\in S} (d'(u,v)-d(u,v)) \cdot \omega(u,v) \ge m$  that  $\omega(u,v) > 0$  implies d'(u,v)-d(u,v) = m.

**Corollary 15.** d is the unique fixed point of  $\Delta'$ .

*Proof.* It is enough to prove that d is a fixed point of  $\Delta'$ . On the one hand, suppose that  $s \sim t$ . Then  $d(s,t) = 0 = \Delta'(d)(s,t)$  by Theorem 3. On the other hand, suppose that  $s \not\sim t$ . Then  $d(s,t) = \Delta(d)(s,t) = \Delta'(d)(s,t)$ .

Corollary 15 implies that d is the greatest fixed point of  $\Delta'$ . Thus, following the development in Section 5.1, we can compute d as the solution to the following linear program:

maximise 
$$\sum_{s,t\in S} d_{s,t}$$
such that  $d_{s,t} = 0$ 
 $d_{s,t} = 1$ 
 $d_{s,t} \leq \sum_{u,v\in S} d_{u,v} \cdot \omega_{u,v}$ 
 $\omega \in V(\Omega_{s,t}), s \not\sim t, \ell(s) = \ell(t)$ 

$$(9)$$

Unfortunately we cannot solve (9) using the iterative method adopted in the discounted case. The reason is that it may require exponentially many iterations of  $\Delta'$  to achieve a sufficiently close approximation to d.

*Example 16.* Consider the Markov chain below, where  $\ell(s) = \ell(t) \neq \ell(u)$ :



Then d(s,t) = 1 and  $(\Delta')^n(\mathbf{0})(s,t) \leq n \cdot 2^{-m}$  for all  $n \in \mathbb{N}$ . This shows that it may require exponentially many iterations in size( $\mathcal{M}$ ) to approximate the fixed point of  $\Delta'$ .

Instead we use the ellipsoid algorithm (see, e.g. [25, Chapter 14]) to solve the linear program (9). According to Proposition 9 (which also holds for c = 1), the coefficients of the constraints of the linear program (9) are rational of size polynomial in size( $\mathcal{M}$ ). By, e.g. [25, Corollary 14.1a], to conclude that the linear program (9) can be solved in time polynomial in size( $\mathcal{M}$ ), it suffices to show that there exists a polynomial time separation algorithm. In our setting, given a  $d \in \mathbb{R}^{S \times S}$  rational of size polynomial in size( $\mathcal{M}$ ), the separation algorithm has to decide whether d satisfies the constraints of (9) or not, and, in the latter case, find in time polynomial in size( $\mathcal{M}$ ) a separating hyperplane, i.e., an  $\alpha \in \mathbb{Q}^{S \times S}$ such that

$$\sum_{u,v\in S} d(u,v) \cdot \alpha(u,v) < \sum_{u,v\in S} d'(u,v) \cdot \alpha(u,v)$$
(10)

for all  $d' \in \mathbb{R}^{S \times S}$  that satisfy the constraints of (9).

Let  $d \in \mathbb{R}^{S \times S}$  be rational of size polynomial in size( $\mathcal{M}$ ). For each pair of states  $s, t \in S$ , we consider the following linear program:

minimise 
$$\sum_{\substack{u,v \in S \\ v \in S \\ u,v \in S \\ u,v = \pi_{s,u} \\ \sum_{\substack{u \in S \\ u \in S \\ \omega_{u,v} = \pi_{t,v} \\ \omega_{u,v} \ge 0}} (11)$$

This linear program is a minimum-cost flow problem for which there are versions of the network simplex algorithm that can compute an  $(\omega_{u,v})_{u,v\in S}$ , which satisfies the constraints of (11) and minimizes the objective function, and that are strongly polynomial time [1].

Note that d satisfies the constraints of (11) if and only if d(s,t) is smaller than or equal to the optimal value of (11) for each pair of states  $s, t \in S$ . Otherwise, there exists a pair of states  $s, t \in S$  such that d(s,t) is greater than the optimal value of (11). Let  $\omega \in V(\Omega_{s,t})$  be a vertex that realizes the optimal value of (11). As we have seen in Proposition 9,  $\omega$  is rational of size polynomial in size( $\mathcal{M}$ ).

It remains to define an  $\alpha$  that satisfies (10). We define  $\alpha$  in terms of  $\omega$  as follows:

$$\alpha(u, v) = \begin{cases} \omega(u, v) - 1 \text{ if } (u, v) = (s, t) \\ \omega(u, v) & \text{otherwise.} \end{cases}$$

**Proposition 17.** Assume that d does not satisfy the constraints of (11). Then for all  $d' \in \mathbb{R}^{S \times S}$  that satisfy the constraints of (11), we have (10).

*Proof.* Since d does not satisfy the constraints of (11), there exists a pair of states  $s, t \in S$  such that  $d(s,t) > \sum_{u,v \in S} d(u,v) \cdot \omega(u,v)$ . Hence,

$$\sum_{u,v\in S} d(u,v) \cdot \alpha(u,v) < 0.$$
(12)

Let  $d' \in \mathbb{R}^{S \times S}$  satisfying the constraints of (11). Then  $d'(s,t) \leq \sum_{u,v \in S} d'(u,v) \cdot \omega(u,v)$ . Hence,

$$\sum_{u,v\in S} d'(u,v) \cdot \alpha(u,v) \ge 0.$$
(13)

¿From (12) and (13) we can immediately conclude (10).

# 6 Conclusions

It is interesting to compare the problem of computing the bisimilarity pseudometric to that of computing the value of a Shapley stochastic game [12]. Both problems involve computing a fixed point of a function  $F : \mathbb{R}^n \to \mathbb{R}^n$  such that computing F requires solving a linear program. However the function F is simpler in the case of the pseudometric, being in particular polynomial piecewise linear in the sense of [12]. On the other hand, for Shapley games the problem of deciding whether the value of the game exceeds a given threshold is as hard as the sum-of-square-roots problem—a problem not known even to be in **NP**. The problem of approximating the value of a Shapely game is known to be in **PPAD**, but is not known to be in **P** either.

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### A Some proofs

Due to the lack of space, for some results we only provide proof sketches. Below, we provide complete proofs for some of those results.

#### A.1 The Distances are Rational

Recall that for each  $s, t \in S$ , the set  $\Omega_{s,t}$  is defined by

$$\Omega_{s,t} = \{ \omega \in \mathbb{R}^{S \times S} : \forall u, v \in S \ \omega(u,v) \ge 0 \land \\ \forall u \in S \ \sum_{v \in S} \omega(u,v) = \pi_{s,u} \land \\ \forall v \in S \ \sum_{u \in S} \omega(u,v) = \pi_{t,v} \}$$

Note that

$$\sum_{u,v\in S} \omega(u,v) = \sum_{u\in S} \pi_{s,u} = 1$$

Hence, for all  $u, v \in S$ ,  $\omega(u, v) \leq 1$ . Therefore, the set  $\Omega_{s,t}$  is bounded. We leave it to the reader to verify that the set  $\Omega_{s,t}$  is convex. Hence,  $\Omega_{s,t}$  is a polytope. Recall that  $V(\Omega_{s,t})$  denotes the set of vertices of  $\Omega_{s,t}$ .

Next, we give a bound for the denominators of the vertices. We will use this to bound the denominators of the distances.

**Proposition 18.** Let  $\delta$  be the least common divisor of the denominators of  $\pi$ . For each  $\omega \in V(\Omega_{s,t})$  and  $u, v \in S$ ,  $\omega_{u,v} = \frac{p}{\delta}$  for some  $0 \le p \le \delta$ .

*Proof.* By definition, a vertex is an intersection of hyperplanes given by the (in)equalities defining  $\Omega_{s,t}$ . Obviously, the (in)equalities  $\omega(u, v) \geq 0$  give rise to zeroes. Let us focus on the other equalities. We multiply all  $\pi_{s,u}$  and  $\pi_{t,v}$  by  $\delta$ , turning them into integers. Obviously, it suffices to show that this modified system of equations has an integral solution. By viewing  $(\omega(u, v))_{u,v\in S}$  and  $(\delta \cdot \pi_{s,u})_{u\in S}(\delta \cdot \pi_{t,v})_{v\in S}$  both as a vector, the system of equations can be captured as the equation  $A\omega = \delta\pi$ . Let n be the number of states. Then A is an  $2n \times n^2$ -matrix. The first n rows are defined by

$$A_{(s,w),(u,v)} = \begin{cases} 1 \text{ if } w = u\\ 0 \text{ otherwise} \end{cases}$$

and the second n rows are defined by

$$A_{(t,w),(u,v)} = \begin{cases} 1 \text{ if } w = v \\ 0 \text{ otherwise.} \end{cases}$$

Note that all entries of A are zero or one, each column of A contains two ones, one in the first n rows and one in the second n rows. Hence, we can conclude from e.g. [25, page 276] that A is totally unimodular. From the proof of Theorem 15 of [2, page 111] we can conclude that the vector  $\omega$  is integral.

Now consider the set D defined by

$$\begin{aligned} D &= \{ d \in \mathbb{R}^{S \times S} : \forall s, t \in S \ \ell(s) \neq \ell(t) \Rightarrow d(s, t) = 1 \land \\ \ell(s) &= \ell(t) \Rightarrow \forall \omega \in V(\Omega_{s,t}) \ d(s, t) \leq c \cdot \sum_{u,v \in S} d(u,v) \cdot \omega(u,v) \land \\ d(s,t) \geq 0 \land \\ d(s,t) \leq 1 \}. \end{aligned}$$

Obviously, the set D is bounded. Again, we leave it to the reader to verify that the set D is convex. Let V(D) be the set of vertices of the polytope D. Next, we give a bound for the denominators of these vertices as well.

**Proposition 19.** Let  $\gamma$  be the denominator of c. Let  $\delta$  be the least common divisor of the denominators of  $\pi$ . For each  $d \in V(D)$ , there exists a  $q \leq (\gamma \delta)^{n^2+1} (n_2)^{\frac{n^2}{2}}$  such that for all  $u, v \in S$ ,  $d(u, v) = \frac{p}{q}$  for some  $0 \leq p \leq q$ .

*Proof.* By definition, a vertex is an intersection of hyperplanes given by the (in) equalities defining D. We can rewrite the inequality

$$d(s,t) \leq c \cdot \sum_{u,v \in S} d(u,v) \cdot \omega(u,v)$$

 $\operatorname{to}$ 

$$\sum_{u,v\in S} \omega'(u,v) \cdot d(u,v) \ge 0$$

where

$$\omega'(u,v) = \begin{cases} c \cdot \omega(u,v) - 1 \text{ if } (u,v) = (s,t) \\ c \cdot \omega(u,v) & \text{otherwise} \end{cases}$$

We multiply all  $\omega'(u, v)$  by  $\gamma\delta$ , turning them into integers. Note that  $\gamma\delta\cdot\omega'(u, v) \leq \gamma\delta$ , since  $\omega(u, v) \leq 1$ . We view  $(d(u, v))_{u,v\in S}$  as a vector. The system of equations can be formulated as Ad = 0 where the non-zero entries of the  $n^2 \times n^2$ -matrix A are  $\omega'_{u,v}$ . Cramer's rule tells us that the denominators of d(u, v) divide the determinant of A (multiplied by  $\gamma\delta$ ). From the Hadamard's inequality, we can conclude that  $(\gamma\delta)^{n^2}(n_2)^{\frac{n^2}{2}}$  is a bound for the the determinant of A. Hence, the denominators of d(u, v) are bounded by  $(\gamma\delta)^{n^2+1}(n_2)^{\frac{n^2}{2}}$ .

**Corollary 20.** Let  $\gamma$  be the denominator of c. Let  $\delta$  be the least common divisor of the denominators of  $\pi$ . There exists a  $q \leq (\gamma \delta)^{n^2+1} (n_2)^{\frac{n^2}{2}}$  such that for all  $s, t \in S$ ,  $\mathsf{d}_c(s,t) = \frac{p}{q}$  for some  $0 \leq p \leq q$ .

*Proof.* Since the linear function mapping  $d \in D$  to  $\sum_{s,t\in S} d(s,t)$  achieves its maximum at a vertex of D and  $\mathsf{d}_c$  is that maximum, the result follows from Proposition 19.

Corollary 21.  $d_c$  is rational.

### A.2 Couplings

We have already remarked that bisimilarity is closely related to the notion of lumpability in Markov chains. The characterisation in Proposition 4 also illustrates a connection with the notion of coupling of Markov chains. For each  $s, t \in S$ , let  $\omega_{(s,t),(-,-)}$  be a matching of  $\pi_{s,-}$  and  $\pi_{t,-}$ . Then the Markov chain  $\mathcal{C} = (S \times S, \omega)$  is a coupling (see, e.g., [3, Chapter 11] for a discussion of couplings).

Such a coupling can be seen as two copies of  $\mathcal{M}$  running sychronously, although not necessarily independently. Couplings are used to give upper bounds on convergence to stationary distributions. Here we use them to a slightly different end. Given a coupling  $\mathcal{C}$ , as above, define the *discrepancy* of a state  $(s,t) \in S \times S$ , denoted  $d_{\mathcal{C}}(s,t)$ , to be the probability that a trajectory of  $\mathcal{C}$ starting in state (s,t) reaches a state (u,v) with  $\ell(u) \neq \ell(v)$ .

Formally, given a coupling  $\mathcal{C}$ , we define  $\Gamma_{\mathcal{C}} : [0,1]^{S \times S} \to [0,1]^{S \times S}$  as follows. If  $\ell(s) \neq \ell(t)$  then  $\Gamma_{\mathcal{C}}(d)(s,t) = 1$  and if  $\ell(s) = \ell(t)$  then

$$\Gamma_{\mathcal{C}}(d)(s,t) = \sum_{u,v \in S} d(u,v) \cdot \omega_{(s,t),(u,v)}.$$

We leave it to the reader to check that  $\Gamma_{\mathcal{C}}$  is a monotone selfmap on  $[0, 1]^{S \times S}$ . By Tarski's fixed point theorem,  $\Gamma_{\mathcal{C}}$  has a least fixed point, which we denote by  $d_{\mathcal{C}}$ . As we will show next,  $d_{\mathcal{C}}$  is closely related to our bisimilarity pseudometric  $d_1$ .

**Proposition 22.** For every coupling C,  $d_1 \leq d_C$ .

*Proof.* According to Tarski's fixed point theorem,  $d_1$  is the least pre-fixed point of  $\Delta_1$ . Hence, it suffices to show that  $d_{\mathcal{C}}$  is a pre-fixed point of  $\Delta_1$  as well. We distinguish two cases. If  $\ell(s) \neq \ell(t)$ , then  $\Delta_1(d_{\mathcal{C}})(s,t) = 1 = d_{\mathcal{C}}(s,t)$ . If  $\ell(s) = \ell(t)$  then

$$\Delta_1(d_{\mathcal{C}})(s,t) = \min_{\omega \in \Omega_{s,t}} \sum_{u,v \in S} d_{\mathcal{C}}(u,v) \cdot \omega(u,v)$$
$$\leq \sum_{u,v \in S} d_{\mathcal{C}}(u,v) \cdot \omega_{(s,t),(u,v)}$$
$$= d_{\mathcal{C}}(s,t).$$

#### **Proposition 23.** For some coupling C, $d_1 \ge d_C$ .

*Proof.* We start by constructing a coupling, i.e., for each  $s, t \in S$  we give a  $\omega_{(s,t),(-,-)}$  which is a matching of  $\pi_{s,-}$  and  $\pi_{t,-}$ . Let  $s, t \in S$ . If  $\ell(s) \neq \ell(t)$ , then we define  $\omega_{(s,t),(u,v)} = \pi_{s,u} \cdot \pi_{t,v}$ . We leave it to the reader to verify that  $\omega_{(s,t),(-,-)}$  which is a matching of  $\pi_{s,-}$  and  $\pi_{t,-}$ . If  $\ell(s) = \ell(t)$ , then

$$\mathsf{d}_1(s,t) = \min_{\omega \in \Omega_{s,t}} \sum_{u,v \in S} d_{\mathcal{C}}(u,v) \cdot \omega(u,v)$$

since  $\mathsf{d}_1$  is a fixed point of  $\Delta_1$ . Let  $\omega_{(s,t),(-,-)}$  be a vertex of  $\Omega_{s,t}$  which minimizes the linear function mapping a matching  $\omega$  to  $\sum_{u,v\in S} d_{\mathcal{C}}(u,v) \cdot \omega(u,v)$ , i.e.

$$\mathsf{d}_1(s,t) = \sum_{u,v \in S} d_{\mathcal{C}}(u,v) \cdot \omega_{(s,t),(u,v)}.$$
(14)

By definition,  $\omega_{(s,t),(-,-)}$  is a matching of  $\pi_{s,-}$  and  $\pi_{t,-}$ .

Since  $d_{\mathcal{C}}$  is the least fixed point of  $\Gamma_{\mathcal{C}}$ , it suffices to show that  $\mathsf{d}_1$  is a fixed point of  $\Gamma_{\mathcal{C}}$ . If  $\ell(s) \neq \ell(t)$  then  $\Gamma_{\mathcal{C}}(\mathsf{d}_1)(s,t) = 1 = \mathsf{d}_1(s,t)$ . If  $\ell(s) = \ell(t)$  then

$$\Gamma_{\mathcal{C}}(\mathsf{d}_1)(s,t) = \sum_{u,v \in S} d_{\mathcal{C}}(u,v) \cdot \omega_{(s,t),(u,v)}$$
$$= \mathsf{d}_1(s,t) \quad [(14)]$$

As a consequence of the above two propositions, the bisimilarity pseudometric  $d_1$  corresponds to the minimal coupling. Next, we will show that two states have discrepancy zero in some coupling if and only if they are bisimilar.

## **Proposition 24.** $d_{\mathcal{C}}(s,t) = 0$ for some coupling $\mathcal{C}$ if and only if $s \sim t$ .

*Proof.* From Proposition 22 and 23 we can conclude that  $d_{\mathcal{C}}(s,t) = 0$  for some coupling  $\mathcal{C}$  if and only if  $\mathsf{d}_1(s,t) = 0$ . In combination with Theorem 3, this gives us the desired result.

The following simple lemma shows that the discrepancy can be used to bound the variational distance between trace distributions. This can be seen as a quantitative version of the folklore that bisimilar states satisfy the same linear-time properties.

**Lemma 25 (Coupling Lemma).** Let C be a coupling of the labelled Markov chain  $\mathcal{M} = (S, \Sigma, \pi, \ell)$ . Then for any measurable set  $A \subseteq \Sigma^{\omega}$  and  $s, t \in S$ ,

$$|\Pr_{\mathcal{M},s}(A) - \Pr_{\mathcal{M},t}(A)| \le d_{\mathcal{C}}(s,t).$$

*Proof.* Let the random variable  $(X, Y) \in \Sigma^{\omega} \times \Sigma^{\omega}$  denote the infinite sequence of labels along a run of  $\mathcal{C}$  that starts in state (s, t). Then

$$Pr_{\mathcal{M},s}(A) = Pr(X \in A)$$
  

$$\geq Pr((X = Y) \cap (Y \in A))$$
  

$$= 1 - Pr((X \neq Y) \cup (Y \notin A))$$
  

$$\geq 1 - Pr(X \neq Y) - Pr(Y \notin A)$$
  

$$= Pr(Y \in A) - Pr(X \neq Y)$$
  

$$= Pr(Y \in A) - d_{\mathcal{C}}(s, t)$$
  

$$= Pr_{\mathcal{M},t}(A) - d_{\mathcal{C}}(s, t) .$$

As a consequence of the Coupling Lemma and Proposition 23, we get that our bisimilarity pseudometric is a bound for the variational distance between trace distributions. In the lemma, we use  $\Pr_{\mathcal{M},s}(A)$  to denote the probability that a run of the labelled Markov chain  $\mathcal{M}$  started in state *s* is in the set *A*. For a formal definition of  $\Pr_{\mathcal{M},s}(A)$  and a definition of measurable subset of the set  $\Sigma^{\omega}$  of infinite sequences over  $\Sigma$ , we refer the reader to, e.g., [1, Chapter 10].

**Corollary 26.** For any measurable set  $A \subseteq \Sigma^{\omega}$  and  $s, t \in S$ ,

$$|\Pr_{\mathcal{M},s}(A) - \Pr_{\mathcal{M},t}(A)| \le \mathsf{d}_1(s,t)(s,t)$$

Whereas our bisimilarity pseudometric can be computed in polynomial time, the variational distance between trace distributions is **NP**-hard to compute as shown by Lyngsø and Pedersen in [21].

# References

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