On Sufficient and Necessary Conditions in transition system with Finite States: a preliminary report

Abstract

This article describes a method to compute the SNC (WSC) of a given property (a CTL formula) under a given transition system (expressed as a Kripke structure) by proposing a semantic forgetting for CTL. We show that the CTL system is close under our definition of forgetting, and this definition satisfies those four postulates of forgetting at first. By changing a transition system $\mathcal{M}$ into its characteristic formula, we can compute the WSC (SNC) of a property under the transition system. We also investigate some properties and algorithm for the forgetting.

1 Introduction

Weakest precondition, we also call *weakest sufficient condition* (WSC), is introduced by Dijkstra in [Dijkstra, 1978]. *Strongest postcondition* (we also call *strongest necessary condition* (SNC)), a dual concept, was introduced subsequently. The SNC is the most general consequent and the WSC is the most general precondition. In program correctness methods, SNC and WSC meet their toughest challenge when they deal with iterative constructs [Mraihi *et al.*, 2011]. It is known that the computing of WSC for code fragment S with respect to assertion Q requires S must terminate [Tremblay, 1996].

However, in model checking of concurrent system with there is at least one successor state for each state s in this system, the termination of this system is impossible. Informally, given a transition system $\mathcal{M}$ with s_0 as an initial state and a specification φ (in this article we suppose it is a *Computation Tree Logic* (CTL in short) [Clarke and Emerson, 1981] formula), we should decide whether $(\mathcal{M}, s_0) \models \varphi$. It is a good thing if $(\mathcal{M}, s_0) \models \varphi$ indeed. However, if $(\mathcal{M}, s_0) \not\models \varphi$, how can we find the WSC ψ under a given set of atoms such that $(\mathcal{M}, s_0) \models \psi \supset \varphi$. It can be shown by the following example.

Example 1 A Beverage Vending Machine, which has been established as standard in the field of process calculi, can be described as a Kripke structure $\mathcal{M} = (S, R, L, s_0)$ on $V_a = \{select, pay, beer, soda\}$ with $S = \{s_0, s_1, s_2\}$, $R = \{(s_0, s_1), (s_0, s_2), (s_1, s_0), (s_2, s_0)\}$, $L(s_0) = \{select\}$, $L(s_1) = \{pay, soda\}$, $L(s_2) = \{pay, beer\}$ and s_0 is an initial state. Which means that when we in s_0 if we select soda and pay for it then we change to the s_1, else if we select beer and pay for it then we change to the s_2, after takeing out the drink we transform to s_0. This is somewhat different from that in [Baier and Katoen, 2008] for simply. For convenience, we use s for $select$, p for pay, b for $beer$ and so for $soda$. Let $\varphi = \text{AGAF}(p \wedge r)$, which means $p \wedge r$ will be satisfied infinite times in the structure, be a CTL formula with r expressing orange juice.

We can decide $(\mathcal{M}, s_0) \not\models \varphi$ easily due to this structure do not contain the atom r. In order for $(\mathcal{M}, s_0)$ satisfy φ, we should find a condition ψ such that $(\mathcal{M}, s_0) \models \psi \supset \varphi$. As we know that if this condition exists, there are many conditions that satisfy the need. In this case, if we are clever enough to judge in advance the set of possible atomic propositions that make up the condition, then we can find this condition in the set only, and the smaller the set, the easier it is to work out the condition. In this paper, we always assume that the condition is a property defined on the specified atomic proposition set V, for our example $V = \{p, r\}$, and find the weakest property (that is, the weakest sufficient condition) satisfying the condition on the set. Finding this property is called discovering theorem by Lin in [Lin, 2018]. Inspired by the forgetting-based method to compute SNC (WSC) [Lin, 2001], in this paper, we tackle this problem by proposing a semantic forgetting for CTL.

However, as we have said that $\mathcal{M}$ is a Kripke structure, which needs to be converted into a logical formula (theory), that is the characteristic formula, which is a CTL formula proposed in [Browne *et al.*, 1988]. Thanks to we find the WSC in a set V of atoms, hence a set-based bisimulation between two K-structures (a Kripke structure with a state in it), V-bisimulation, and characteristic formula on V will be proposed in this paper. Our V-bisimulation is a more general bisimulation relation than others. On the one hand, the above set-based bisimulation is an extension of the bisimulation-equivalence of Definition 7.1 in [Baier and Katoen, 2008] in the sense that if $V = \mathcal{A}$ then our bisimulation is almost same to the latter. On the other hand, the above set-based bisimulation notion is similar to the state equivalence in [Browne *et al.*, 1988]. But it is different in the sense that ours is defined on K-structures, while it is defined on states in [Browne *et al.*, 1988]. What's more, the set-based bisimulation notion is also different from the state-based bisimulation notion of

Definition 7.7 in [Baier and Katoen, 2008], which is defined for states of a given K-structure.

As a logical notion, forgetting was first formally defined in propostional and first order logics by Lin and Reiter [Lin and Reiter, 1994]. Over the last twenty years, researchers have developed forgetting notions and theories not only in classical logic but also in other non-classical logic systems [Eiter and Kern-Isberner, 2019], such as forgetting in logic programs under answer set/stable model semantics [Zhang and Foo, 2006; Eiter and Wang, 2008; Wong, 2009; Wang *et al.*, 2012; Wang *et al.*, 2013], forgetting in description logic [Wang *et al.*, 2010; Lutz and Wolter, 2011; Zhao and Schmidt, 2017] and knowledge forgetting in modal logic [Zhang and Zhou, 2009; Su *et al.*, 2009; Liu and Wen, 2011; Fang *et al.*, 2019]. In application, forgetting has been used in planning [Lin, 2003], conflict solving [Lang and Marquis, 2010; Zhang *et al.*, 2005], createing restricted views of ontologies [Zhao and Schmidt, 2018], strongest and weakest definitions [Lang and Marquis, 2008], SNC (WSC) [Lin, 2001] and so on.

Though forgetting has been extensively investigated from various aspects of different logical systems. However, the existing forgetting method in propositional logic, answer set programming, description logic and modal logic are not directly applicable in CTL. Similar with that in [Zhang and Zhou, 2009], we research forgetting in CTL from the semantic forgetting point of view. And it is shown that our definition of forgetting satisfies those four postulates of forgetting.

The rest of the paper is organised as follows. Section 2 introduces the related notions for forgetting in CTL, including the syntax and semantics of CTL, the language we aimed for. A formal definition of concept forgetting and its properties for CTL follows in Section 3. Section 4 explores the relation between forgetting and SNC (WSC). From the point of view of model, we propose an algorithm for computing forgetting on CTL in Section 5. Finally, we conclude this paper.

2 Preliminaries

We start with some technical and notational preliminaries. Throughout this paper, we fix a finite set $\mathcal{A}$ of propositional variables (or atoms), and use V, V' for subsets of $\mathcal{A}$. In this part, we will introduce the structure we will use for CTL and syntactic and semantic of CTL.

2.1 Model structure in CTL

In general, a transition system [1] is described as a *model structure* (or *Kripke structure*), and a model structure is a triple $\mathcal{M} = (S, R, L)$, where

- S is a finite nonempty set of states,
- $R \subseteq S \times S$ and, for each $s \in S$, there is $s' \in S$ such that $(s, s') \in R$,
- L is a labeling function $S \to 2^{\mathcal{A}}$.

[1] According to [Baier and Katoen, 2008], a *transition system* TS is a tuple $(S, Act, \to, I, AP, L)$ where (1) S is a set of states, (2) Act is a set of actions, (3) $\to \subseteq S \times \text{Act} \times S$ is a transition relation, (4) $I \subseteq S$ is a set of initial states, (5) AP is a set of atomic propositions, and (6) $L : S \to 2^{\text{AP}}$ is a labeling function.

We call a model structure $\mathcal{M}$ on a set V of atoms if $L : S \to 2^V$, *i.e.* , the labeling function L map every state to V (not the $\mathcal{A}$). A *path* π_{s_i} start from s_i of $\mathcal{M}$ is a infinite sequence of states $\pi_{s_i} = (s_i, s_{i+1}s_{i+2}, \dots)$, where for each j $(0 \leq i \leq j)$, $(s_j, s_{j+1}) \in R$. By $s' \in \pi_{s_i}$ we mean that s' is a state in the path π_{s_i}. A sate $s \in S$ is *initial* if for any state $s' \in S$, there is a path π_s s.t $s' \in \pi_s$. We denote this model structure as (S, R, L, s_0), where s_0 is initial.

For a given model structure (S, R, L, s_0) and $s \in S$, the *computation tree* $\text{Tr}_n^{\mathcal{M}}(s)$ of $\mathcal{M}$(or simply $\text{Tr}_n(s)$), that has depth n and is rooted at s, is recursively defined as [Browne *et al.*, 1988], for $n \geq 0$,

- $\text{Tr}_0(s)$ consists of a single node s with label s.
- $\text{Tr}_{n+1}(s)$ has as its root a node m with label s, and if $(s, s') \in R$ then the node m has a subtree $\text{Tr}_n(s')$.

By s_n we mean a nth level node of tree $\text{Tr}_m(s)$ $(m \geq n)$.

A K-*structure* (or K-*interpretation*) is a model structure $\mathcal{M} = (S, R, L, s_0)$ associating with a state $s \in S$, which is written as $(\mathcal{M}, s)$ for convenience in the following. In the case s is an initial state of $\mathcal{M}$, the K-structure is *initial*.

2.2 Syntax and semantics of CTL

In the following we briefly review the basic syntax and semantics of the CTL [Clarke *et al.*, 1986]. The *signature* of $\mathcal{L}$ includes:

- a finite set of Boolean variables, called *atoms* of $\mathcal{L}$: $\mathcal{A}$;
- the classical connectives: $\bot, \vee$ and $\neg$;
- the path quantifiers: A and E;
- the temporal operators: X, F, G U and W, that means 'neXt state', 'some Future state', 'all future states (Globally)', 'Until' and 'Unless', respectively;
- parentheses: (and).

The *(existential normal form or ENF in short) formulas* of $\mathcal{L}$ are inductively defined via a Backus Naur form:

$$\phi ::= \bot \mid p \mid \neg\phi \mid \phi \vee \phi \mid \text{EX}\phi \mid \text{EG}\phi \mid \text{E}[\phi \text{ U } \phi] \quad (1)$$

where $p \in \mathcal{A}$. The formulas $\phi \wedge \psi$ and $\phi \to \psi$ are defined in a standard manner of propositional logic. The other form formulas of $\mathcal{L}$ are abbreviated using the forms of (1). The priorities for the CTL connectives are assumed to be (from the highest to the lowest):

$$\neg, \text{EX}, \text{EF}, \text{EG}, \text{AX}, \text{AF}, \text{AG} \prec \wedge \prec \vee \prec \text{EU}, \text{AU}, \text{EW}, \text{AW}, \to .$$

We are now in the position to define the semantics of $\mathcal{L}$. Let $\mathcal{M} = (S, R, L, s_0)$ be an model structure, $s \in S$ and ϕ a formula of $\mathcal{L}$. The *satisfiability* relationship between $\mathcal{M}, s$ and ϕ, written $(\mathcal{M}, s) \models \phi$, is inductively defined on the structure of ϕ as follows:

- $(\mathcal{M}, s) \not\models \bot$;
- $(\mathcal{M}, s) \models p$ iff $p \in L(s)$;
- $(\mathcal{M}, s) \models \phi_1 \vee \phi_2$ iff $(\mathcal{M}, s) \models \phi_1$ or $(\mathcal{M}, s) \models \phi_2$;
- $(\mathcal{M}, s) \models \neg\phi$ iff $(\mathcal{M}, s) \not\models \phi$;
- $(\mathcal{M}, s) \models \text{EX}\phi$ iff $(\mathcal{M}, s_1) \models \phi$ for some $s_1 \in S$ and $(s, s_1) \in R$;

- $(\mathcal{M}, s) \models \text{EG}\phi$ iff $\mathcal{M}$ has a path $(s_1 = s, s_2, \ldots)$ such that $(\mathcal{M}, s_i) \models \phi$ for each $i \geq 1$;
- $(\mathcal{M}, s) \models \text{E}[\phi_1 \text{U} \phi_2]$ iff $\mathcal{M}$ has a path $(s_1 = s, s_2, \ldots)$ such that, for some $i \geq 1$, $(\mathcal{M}, s_i) \models \phi_2$ and $(\mathcal{M}, s_j) \models \phi_1$ for each $j < i$.

Similar to the work in [Browne *et al.*, 1988; Bolotov, 1999], only initial K-structures are considered to be candidate models in the following, unless explicitly stated. Formally, an initial K-structure $\mathcal{K}$ is a *model* of a formula ϕ whenever $\mathcal{K} \models \phi$. We denote *Mod*(ϕ) the set of models of ϕ. The formula ϕ is *satisfiable* if *Mod*(ϕ) $\neq \emptyset$. Since the states in model structure is finite, *Mod*(ϕ) is finite for any formula ϕ.

Let ϕ_1 and ϕ_2 be two formulas or set of formulas. By $\phi_1 \models \phi_2$ we denote *Mod*(ϕ_1) $\subseteq$ *Mod*(ϕ_2). By $\phi_1 \equiv \phi_2$ we mean $\phi_1 \models \phi_2$ and $\phi_2 \models \phi_1$. In this case ϕ_1 is *equivalent* to ϕ_2. By *Var*(ϕ_1) we mean the set of atoms occurring in ϕ_1. ϕ_1 is V*-irrelevant*, written IR(ϕ_1, V), if there is a formula ψ with *Var*(ψ) $\cap V = \emptyset$ such that $\phi_1 \equiv \psi$.

3 Forgetting in CTL

In this section, we will define the forgetting in CTL by V-bisimulation, set-based bisimulations. Besides, some properties of forgetting are also explored. For convenience, let $\mathcal{M} = (S, R, L, s_0)$, $\mathcal{M}' = (S', R', L', s'_0)$ and $\mathcal{K}_i = (\mathcal{M}_i, s_i)$ with $\mathcal{M}_i = (S_i, R_i, L_i, s^i_0)$, $s_i \in S_i$ and i is an integer.

3.1 Set-based bisimulation

To present a formal definition of forgetting, we need the concepts of V-bisimulation. Inspired by the notion of bisimulation in [Browne *et al.*, 1988], we define the relations $\mathcal{B}_0, \mathcal{B}_1, \ldots$ between K-structures as follows: let $\mathcal{K}_i = (\mathcal{M}_i, s_i)$ with $i \in \{1, 2\}$,

- $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}_0$ if $L_1(s_1) - V = L_2(s_2) - V$;
- for $n \geq 0$, $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}_{n+1}$ if
  - $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}_0$,
  - for every $(s_1, s'_1) \in R_1$, there is $(s_2, s'_2) \in R_2$ such that $(\mathcal{K}'_1, \mathcal{K}'_2) \in \mathcal{B}_n$, and
  - for every $(s_2, s'_2) \in R_2$, there is $(s_1, s'_1) \in R_1$ such that $(\mathcal{K}'_1, \mathcal{K}'_2) \in \mathcal{B}_n$,

  where $\mathcal{K}'_i = (\mathcal{M}_i, s'_i)$ with $i \in \{1, 2\}$.

Now, we define the notion of V-bisimulation between K-structures:

Definition 1 (V-bisimulation) *Let $V \subseteq \mathcal{A}$. The V-*bisimular relation $\mathcal{B}$ *between* K*-structures is defined as:*

$(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}$ *if and only if* $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}_i$ *for all* $i \geq 0$.

In this case, $\mathcal{K}_1$ and $\mathcal{K}_2$ are called V-bisimular.

Proposition 1 *Let $V \subseteq \mathcal{A}$ and $\mathcal{K}_i = (\mathcal{M}_i, s_i)$ $(i = 1, 2)$ be* K*-structures. Then $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}$ if and only if*

(i) $L_1(s_1) - V = L_2(s_2) - V$,

(ii) for every $(s_1, s'_1) \in R_1$, *there is* $(s_2, s'_2) \in R_2$ *such that* $(\mathcal{K}'_1, \mathcal{K}'_2) \in \mathcal{B}$, *and*

(iii) for every $(s_2, s'_2) \in R_2$, *there is* $(s_1, s'_1) \in R_1$ *such that* $(\mathcal{K}'_1, \mathcal{K}'_2) \in \mathcal{B}$,

where $\mathcal{K}'_i = (\mathcal{M}_i, s'_i)$ *with* $i \in \{1, 2\}$.

Two pathes $\pi_i = (s_{i,1}, s_{i,2}, \ldots)$ of $\mathcal{M}_i$ with $i \in \{1, 2\}$ are V*-bisimular* if

$$(\mathcal{K}_{1,j}, \mathcal{K}_{2,j}) \in \mathcal{B} \text{ for every } j \geq 0$$

where $\mathcal{K}_{i,j} = (\mathcal{M}_i, s_{i,j})$.

In the following we abbreviated $(\mathcal{K}_1, \mathcal{K}_2) \in \mathcal{B}$ by $(s_1, s_2) \in \mathcal{B}$ when the underlying model structures of states s_1 and s_2 are clear from their contexts or there is no confusion. The V-bisimular relation is uniformly abbreviated as $\leftrightarrow_V$ for convenience.

Lemma 1 *The relation $\leftrightarrow_V$ is an equivalence relation.*

Besides, we have the following properties:

Proposition 2 *Let $i \in \{1, 2\}$, $V_1, V_2 \subseteq \mathcal{A}$, s'_is be two states and π'_is be two pathes, and $\mathcal{K}_i = (\mathcal{M}_i, s_i)$ $(i = 1, 2, 3)$ be* K*-structures such that $\mathcal{K}_1 \leftrightarrow_{V_1} \mathcal{K}_2$ and $\mathcal{K}_2 \leftrightarrow_{V_2} \mathcal{K}_3$. Then:*

(i) $s'_1 \leftrightarrow_{V_i} s'_2$ $(i = 1, 2)$ *implies* $s'_1 \leftrightarrow_{V_1 \cup V_2} s'_2$;

(ii) $\pi'_1 \leftrightarrow_{V_i} \pi'_2$ $(i = 1, 2)$ *implies* $\pi'_1 \leftrightarrow_{V_1 \cup V_2} \pi'_2$;

(iii) for each path π_{s_1} of $\mathcal{M}_1$ there is a path π_{s_2} of $\mathcal{M}_2$ such that $\pi_{s_1} \leftrightarrow_{V_1} \pi_{s_2}$, and vice versa;

(iv) $\mathcal{K}_1 \leftrightarrow_{V_1 \cup V_2} \mathcal{K}_3$;

(v) If $V_1 \subseteq V_2$ then $\mathcal{K}_1 \leftrightarrow_{V_2} \mathcal{K}_2$.

Intuitively, if two K-structures are V-bisimular, then they satisfy the same formula φ that dose not contain any atoms in V, *i.e.* IR(φ, V).

Theorem 1 *Let $V \subseteq \mathcal{A}$, $\mathcal{K}_i$ $(i = 1, 2)$ be two* K*-structures such that $\mathcal{K}_1 \leftrightarrow_V \mathcal{K}_2$ and ϕ a formula with IR(ϕ, V). Then $\mathcal{K}_1 \models \phi$ if and only if $\mathcal{K}_2 \models \phi$.*

Let $V \subseteq \mathcal{A}$, $\mathcal{M}_i$ $(i = 1, 2)$ be model structures. A computation tree $\text{Tr}_n(s_1)$ of $\mathcal{M}_1$ is V*-bisimular* to a computation tree $\text{Tr}_n(s_2)$ of $\mathcal{M}_2$, written $(\mathcal{M}_1, \text{Tr}_n(s_1)) \leftrightarrow_V (\mathcal{M}_2, \text{Tr}_n(s_2))$ (or simply $\text{Tr}_n(s_1) \leftrightarrow_V \text{Tr}_n(s_2)$), if

- $L_1(s_1) - V = L_2(s_2) - V$,
- for every subtree $\text{Tr}_{n-1}(s'_1)$ of $\text{Tr}_n(s_1)$, $\text{Tr}_n(s_2)$ has a subtree $\text{Tr}_{n-1}(s'_2)$ such that $\text{Tr}_{n-1}(s'_1) \leftrightarrow_V \text{Tr}_{n-1}(s'_2)$, and

Please note that the last condition in the above definition hold trivially for $n = 0$.

Proposition 3 *Let $V \subseteq \mathcal{A}$ and $(\mathcal{M}_i, s_i)$ $(i = 1, 2)$ be two* K*-structures. Then*

$(s_1, s_2) \in \mathcal{B}_n$ *iff* $Tr_j(s_1) \leftrightarrow_V Tr_j(s_2)$ *for every* $0 \leq j \leq n$.

This means that $\text{Tr}_j(s_1) \leftrightarrow_V \text{Tr}_j(s_2)$ for all $j \geq 0$ if $s_1 \leftrightarrow_V s_2$, otherwise there is some number k such that $\text{Tr}_k(s_1)$ and $\text{Tr}_k(s_2)$ are not V-bisimular.

Proposition 4 *Let $V \subseteq \mathcal{A}$, $\mathcal{M}$ be a model structure and $s, s' \in S$ such that $(s, s') \notin \mathcal{B}$. There exists a least number k such that $Tr_k(s)$ and $Tr_k(s')$ are not V-bisimular.*

In this case the model structure $\mathcal{M}$ is called V*-distinguishable* (by states s and s' at the least depth k), which is denoted by $\text{dis}_V(\mathcal{M}, s, s', k)$. It is evident that $\text{dis}_V(\mathcal{M}, s, s', k)$ implies $\text{dis}_V(\mathcal{M}, s, s', k')$ whenever $k' \geq$

k. The V-*characterization number* of $\mathcal{M}$, written $ch(\mathcal{M}, V)$, is defined as

$$ch(\mathcal{M},V)=\begin{cases}\max\{k \mid s,s'\in S \ \&\ \mathrm{dis}_V(\mathcal{M},s,s',k)\},\\ \qquad\qquad \mathcal{M} \text{ is } V\text{-distinguishable};\\ \min\{k \mid \mathcal{B}_k=\mathcal{B}\}, \qquad\qquad \text{otherwise.}\end{cases}$$

Now we give the formal definition of forgetting in CTL from the semantic forgetting point view.

Definition 2 (Forgetting) *Let $V \subseteq \mathcal{A}$ and ϕ a formula. A formula ψ with Var(ψ) $\cap$ $V = \emptyset$ is a* result of forgetting V from ϕ, *if*

$$Mod(\psi) = \{\mathcal{K} \text{ is initial} \mid \exists \mathcal{K}' \in Mod(\phi) \ \&\ \mathcal{K}' \leftrightarrow_V \mathcal{K}\}. \quad (2)$$

Note that if both ψ and ψ' are results of forgetting V from ϕ then $Mod(\psi) = Mod(\psi')$, *i.e.* , ψ and ψ' have the same models. In the sense of equivalence the forgetting result is unique (up to equivalence).

Intuitively, forgetting an atom results in a weaker theory which entails the same set of formulae that are irrelevant to the atom. To present the representation property of forgetting in CTL and compute WSC (SNC) under an initial K-structure, we will give the Characterize formula of an initial K-structure on V in the next subsection.

3.2 Characterize formula of initial K-structure

Given a set $V \subseteq \mathcal{A}$, we can define a formula φ of V (that is $Var(\varphi) \subseteq V$) in CTL to equivalent uniquely describe a computation tree.

Definition 3 *Let $V \subseteq \mathcal{A}$, $\mathcal{M} = (S, R, L, s_0)$ be a model structure and $s \in S$. The* characterize formula *of the computation tree $Tr_n(s)$ on V, written $\mathcal{F}_V(Tr_n(s))$, is defined recursively as:*

$$\mathcal{F}_V(Tr_0(s)) = \bigwedge_{p\in V\cap L(s)} p \wedge \bigwedge_{q\in V-L(s)} \neg q,$$

$$\mathcal{F}_V(Tr_{k+1}(s)) = \bigwedge_{(s,s')\in R} \mathrm{EX}T(s') \wedge \mathrm{AX} \bigvee_{(s,s')\in R} T(s') \wedge \mathcal{F}_V(Tr_0(s))$$

for $k \geq 0$, where $T(s') = \mathcal{F}_V(Tr_k(s'))$.

The characterize formula of a computation tree formally exhibit the context of each node on V (atoms are true at this node if they are in V, else false) and the temporal relation between states recursively. In this way, we know:

Lemma 2 *Let $V \subseteq \mathcal{A}$, $\mathcal{M} = (S, R, L, s_0)$ and $\mathcal{M}' = (S', R', L', s_0')$ be two model structures, $s \in S$, $s' \in S'$ and $n \geq 0$. If $Tr_n(s) \leftrightarrow_{\overline{V}} Tr_n(s')$, then $\mathcal{F}_V(Tr_n(s)) \equiv \mathcal{F}_V(Tr_n(s'))$.*

Let $s' = s$, it shows that for any formula φ of V, if φ is a characterize formula of $\mathrm{Tr}_n(s)$ then $\varphi \equiv \mathcal{F}_V(\mathrm{Tr}_n(s))$.

Let $V \subseteq \mathcal{A}$, $\mathcal{K} = (\mathcal{M}, s_0)$ be an initial K-structure and $T(s') = \mathcal{F}_V(\mathrm{Tr}_c(s'))$. The *characterizing formula* of $\mathcal{K}$ on V, written $\mathcal{F}_V(\mathcal{M}, s_0)$ (or $\mathcal{F}_V(\mathcal{K})$), is defined as the conjunction of the following formulas:

$\mathcal{F}_V(\mathrm{Tr}_c(s_0))$, and for each $s \in S$

$$\mathrm{AG}\left(\mathcal{F}_V(\mathrm{Tr}_c(s)) \rightarrow \bigwedge_{(s,s')\in R} \mathrm{EX}T(s') \wedge \mathrm{AX} \bigvee_{(s,s')\in R} T(s')\right)$$

where $c = ch(\mathcal{M}, V)$. It is apparent that $\mathrm{IR}(\mathcal{F}_V(\mathcal{M}, s_0), \overline{V})$.

The following example show how to compute characterizing formula:

Example 2 Let $\mathcal{K} = (\mathcal{M}, s_0)$ with $\mathcal{M} = (S, R, L, s_0)$ be a initial K-structure,in which $S = \{s_0, s_1, s_2\}$, $R = \{(s_0, s_1), (s_0, s_2), (s_1, s_0), (s_2, s_0)\}$, $L(s_0) = \{a\}$, $L(s_1) = \{a, c\}$ and $L(s_2) = \{b, c\}$. Let $V = \{a, b\}$, compute the characterizing formula of $\mathcal{K}$ on V.

It is apparent that $\mathrm{Tr}_0(s_0) \leftrightarrow_{\overline{V}} \mathrm{Tr}_0(s_1)$ due to $L(s_0) - \overline{V} = L(s_1) - \overline{V}$, $\mathrm{Tr}_1(s_0) \not\leftrightarrow_{\overline{V}} \mathrm{Tr}_1(s_1)$ due to there is $(s_0, s_2) \in R$ such that for any $(s_1, s') \in R$ (there is only one immediate successor $s' = s_0$) there is $L(s_2) - \overline{V} \neq L(s') - \overline{V}$. Hence, we have that $\mathcal{M}$ is $\overline{V}$-distinguished by state s_0 and s_1 at the least depth 1, *i.e.* $\mathrm{dis}_{\overline{V}}(\mathcal{M}, s_0, s_1, 1)$. Similarly, we have $\mathrm{dis}_{\overline{V}}(\mathcal{M}, s_0, s_2, 0)$ and $\mathrm{dis}_{\overline{V}}(\mathcal{M}, s_1, s_2, 0)$. Therefore, $ch(\mathcal{M}, \overline{V}) = \max\{k \mid s, s' \in S \ \&\ \mathrm{dis}_{\overline{V}}(\mathcal{M}, s, s', k)\} = 1$. Then we have:

$$\begin{aligned}
&\mathcal{F}_V(\mathrm{Tr}_0(s_0)) = a \wedge \neg b,\\
&\mathcal{F}_V(\mathrm{Tr}_0(s_1)) = a \wedge \neg b,\\
&\mathcal{F}_V(\mathrm{Tr}_0(s_2)) = b \wedge \neg a,\\
&\mathcal{F}_V(\mathrm{Tr}_1(s_0)) = \mathrm{EX}(a \wedge \neg b) \wedge \mathrm{EX}(b \wedge \neg a) \wedge \mathrm{AX}((a \wedge \neg b) \vee\\
&\qquad\qquad (b \wedge \neg a)) \wedge (a \wedge \neg b),\\
&\mathcal{F}_V(\mathrm{Tr}_1(s_1)) = \mathrm{EX}(a \wedge \neg b) \wedge \mathrm{AX}(a \wedge \neg b) \wedge (a \wedge \neg b),\\
&\mathcal{F}_V(\mathrm{Tr}_1(s_2)) = \mathrm{EX}(a \wedge \neg b) \wedge \mathrm{AX}(a \wedge \neg b) \wedge (b \wedge \neg a).
\end{aligned}$$

Then it is easy to obtain $\mathcal{F}_V(\mathcal{M}, s_0)$.

Lemma 3 *Let φ be a formula. We have*

$$\varphi \equiv \bigvee_{(\mathcal{M},s_0)\in Mod(\varphi)} \mathcal{F}_{\mathcal{A}}(\mathcal{M}, s_0). \quad (3)$$

It follows that any CTL formula can be described by the disjunction of the characterizing formulas of all the models of itself due to the number of models of a CTL formula is finite.

Theorem 2 *Let $V \subseteq \mathcal{A}$, $\mathcal{M} = (S, R, L, s_0)$ be a model structure with initial state s_0 and $\mathcal{M}' = (S', R', L', s_0')$ be a model structure with initial state s_0'. Then*

$(\mathcal{M}', s_0') \models \mathcal{F}_V(\mathcal{M}, s_0)$ if and only if $(\mathcal{M}, s_0) \leftrightarrow_{\overline{V}} (\mathcal{M}', s_0')$.

By the following theorem we also have that given a set $V \subseteq \mathcal{A}$, the characterizing formula of an initial K-structure is equivalent uniquely describe this initial K-structure on V.

Theorem 3 *Let $V \subseteq \mathcal{A}$, $\mathcal{M} = (S, R, L, s_0)$ a model structure with initial state s_0 and $\mathcal{M}' = (S', R', L', s_0')$ a model structure with initial state s_0'. If $(\mathcal{M}, s_0) \leftrightarrow_{\overline{V}} (\mathcal{M}', s_0')$ then $\mathcal{F}_V(\mathcal{M}, s_0) \equiv \mathcal{F}_V(\mathcal{M}', s_0')$.*

3.3 Semantic properties of forgetting in CTL

In this subsection we study essential semantic properties of forgetting. We will first show that our forgetting satisfy these postulates [Zhang and Zhou, 2009] that precisely characterize the semantics of forgetting. We then discuss other semantic properties of forgetting.

By Lemma 3 and Definition 2, the result ψ of ϕ forget the set V of atoms always exists, which is equivalent to

$$\bigvee_{\mathcal{K}\in\{\mathcal{K}' \text{ is an initial interpretation} \mid \exists \mathcal{K}''\in Mod(\phi)\ \&\ \mathcal{K}''\leftrightarrow_V \mathcal{K}'\}} \mathcal{F}_{\overline{V}}(\mathcal{K}).$$

For this reason, the forgetting result is denoted by $\mathrm{F}_{\mathrm{CTL}}(\phi, V)$.

In the case ψ is a result of forgetting V from ϕ, there are usually some expected properties (called *postulates*: (**W**), (**PP**), (**NP**) and (**IR**)) for it [Zhang and Zhou, 2009].

Theorem 4 *(Representation theorem). Let φ and ψ be two formulas and $V \subseteq \mathcal{A}$. Then the following statements are equivalent:*

(i) $\psi \equiv \mathrm{F}_{\mathrm{CTL}}(\varphi, V)$,

(ii) $\psi \equiv \{\phi | \varphi \models \phi \& IR(\phi, V)\}$,

*(iii) Postulates (**W**), (**PP**), (**NP**) and (**IR**) hold.*

We can see from this theorem that our definition of forgetting under CTL is close, *i.e.* for any CTL formula the result of forgetting is also a CTL formula, and captures the four postulates that forgetting should satisfy.

Lemma 4 *Let φ and α be two CTL formulae and $q \in Var(\varphi \cup \{\alpha\})$. Then $\mathrm{F}_{CTL}(\varphi \cup \{q \leftrightarrow \alpha\}, q) \equiv \varphi$.*

Proposition 5 *Let φ be a formula, V a set of atoms and p an atom such that $p \notin V$. Then:*

$$\mathrm{F}_{\mathrm{CTL}}(\varphi, \{p\} \cup V) \equiv \mathrm{F}_{\mathrm{CTL}}(\mathrm{F}_{\mathrm{CTL}}(\varphi, p), V).$$

This means that the result of forgetting V from φ can be obtained by forgetting atom in V one by one. Similarly, a consequence of the previous proposition is:

Corollary 5 *Let φ be a formula and $V_i \subseteq \mathcal{A}$ $(i = 1, 2)$. Then:*

$$\mathrm{F}_{\mathrm{CTL}}(\varphi, V_1 \cup V_2) \equiv \mathrm{F}_{\mathrm{CTL}}(\mathrm{F}_{\mathrm{CTL}}(\varphi, V_1), V_2).$$

The following results, which are satisfied in both classical proposition logic and modal logic **S5**, further illustrate other essential semantic properties of forgetting.

Proposition 6 *Let φ, φ_i, ψ_i $(i = 1, 2)$ be formulas and $V \subseteq \mathcal{A}$. We have*

(i) $\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$ *is satisfiable iff φ is;*

(ii) If $\varphi_1 \equiv \varphi_2$, *then* $\mathrm{F}_{\mathrm{CTL}}(\varphi_1, V) \equiv \mathrm{F}_{\mathrm{CTL}}(\varphi_2, V)$;

(iii) If $\varphi_1 \models \varphi_2$, *then* $\mathrm{F}_{\mathrm{CTL}}(\varphi_1, V) \models \mathrm{F}_{\mathrm{CTL}}(\varphi_2, V)$;

(iv) $\mathrm{F}_{\mathrm{CTL}}(\psi_1 \vee \psi_2, V) \equiv \mathrm{F}_{\mathrm{CTL}}(\psi_1, V) \vee \mathrm{F}_{\mathrm{CTL}}(\psi_2, V)$;

(v) $\mathrm{F}_{\mathrm{CTL}}(\psi_1 \wedge \psi_2, V) \models \mathrm{F}_{\mathrm{CTL}}(\psi_1, V) \wedge \mathrm{F}_{\mathrm{CTL}}(\psi_2, V)$;

Another interest result is that the forgetting of $PT\varphi$ ($P \in \{\mathrm{E}, \mathrm{A}\}$, $T \in \{\mathrm{F}, \mathrm{X}\}$) on $V \subseteq \mathcal{A}$ can be computed by $PT\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$. This give a convenient method to compute forgetting.

Proposition 7 *Let $V \subseteq \mathcal{A}$ and ϕ a formula.*

(i) $\mathrm{F}_{\mathrm{CTL}}(\mathrm{AX}\phi, V) \equiv \mathrm{AXF}_{\mathrm{CTL}}(\phi, V)$.

(ii) $\mathrm{F}_{\mathrm{CTL}}(\mathrm{EX}\phi, V) \equiv \mathrm{EXF}_{\mathrm{CTL}}(\phi, V)$.

(iii) $\mathrm{F}_{\mathrm{CTL}}(\mathrm{AF}\phi, V) \equiv \mathrm{AFF}_{\mathrm{CTL}}(\phi, V)$.

(iv) $\mathrm{F}_{\mathrm{CTL}}(\mathrm{EF}\phi, V) \equiv \mathrm{EFF}_{\mathrm{CTL}}(\phi, V)$.

3.4 Main complexity

In the following we consider the main complexities of reasoning problems on forgetting in CTL.

Proposition 8 *Let $\mathcal{M}, s_0$ be an I-structure, φ be a CTL formula and V a set of atoms. Deciding whether $\mathcal{M}, s_0$ is a model of $\mathrm{F}_{CTL}(\varphi, V)$ is NP-complete.*

By this proposition, we have:

Theorem 6 *Let φ and ψ be two CTL(*AF*) (a fragment of CTL, in which each formula contains only* AF *temporal connective) formulas and V a set of atoms. Then we have the results:*

(i) deciding whether $\mathrm{F}_{CTL}(\varphi, V) \models \psi$ is co-NP-complete,

(ii) deciding whether $\psi \models \mathrm{F}_{CTL}(\varphi, V)$ is Π_2^P-complete,

(iii) deciding whether $\mathrm{F}_{CTL}(\varphi, V) \models \mathrm{F}_{CTL}(\psi, V)$ is Π_2^P-complete.

The theorem implies:

Corollary 7 *Let φ and ψ be two CTL(*AF*) formulas and V a set of atoms. Then*

(i) deciding if $\psi \equiv \mathrm{F}_{CTL}(\varphi, V)$ is Π_2^P-complete,

(ii) deciding if $\mathrm{F}_{CTL}(\varphi, V) \equiv \varphi$ is co-NP-complete,

(iii) deciding whether $\mathrm{F}_{CTL}(\varphi, V) \equiv \mathrm{F}_{CTL}(\psi, V)$ is Π_2^P-complete.

4 SNC and WSC

In this section, we will give the definition of SNC (WSC) and show that the SNC (WSC) of a specification (a CTL formula) under a given initial K-structure and set V of atoms can be obtained from forgetting in CTL. The SNC (WSC) of a proposition will be given at first:

Definition 4 (sufficient and necessary condition) *Let ϕ be a formula or an initial* K*-structure, ψ be a formula, $V \subseteq Var(\phi)$, $q \in Var(\phi) - V$ and $Var(\psi) \subseteq V$.*

- *ψ is a* necessary condition *(NC in short) of q on V under ϕ if $\phi \models q \rightarrow \psi$.*
- *ψ is a* sufficient condition *(SC in short) of q on V under ϕ if $\phi \models \psi \rightarrow q$.*
- *ψ is a* strongest necessary condition *(SNC in short) of q on V under ϕ if it is a NC of q on V under ϕ and $\phi \models \psi \rightarrow \psi'$ for any NC ψ' of q on V under ϕ.*
- *ψ is a* weakest sufficient condition *(SNC in short) of q on V under ϕ if it is a SC of q on V under ϕ and $\phi \models \psi' \rightarrow \psi$ for any SC ψ' of q on V under ϕ.*

Note that if both ψ and ψ' are SNC (WSC) of q on V under ϕ then $Mod(\psi) = Mod(\psi')$, *i.e.* ψ and ψ' have the same models. In the sense of equivalence the SNC (WSC) is unique (up to equivalence).

Proposition 9 ***(dual)** Let V, q, φ and ψ are the ones in Definition 4.*

(i) ψ is a SNC of q on V under φ iff $\neg\psi$ is a WSC of $\neg q$ on V under φ.

(ii) ψ is a WSC of q on V under φ iff $\neg\psi$ is a SNC of $\neg q$ on V under φ.

This show that the SNC and WSC are in fact dual conditions. Under the dual property, we can consider the SNC party only in sometimes, while the WSC part can be talked similarly.

For the case of formula, we have that the SCN (WSC) of any formula can be defined as follows:

Definition 5 *Let Γ be a formula or an initial* K*-structure, α be a formula and $P \subseteq (Var(\Gamma) \cup Var(\alpha))$. A formula φ of P is said to be a NC (SC) of α on P under Γ iff $\Gamma \models \alpha \rightarrow \varphi$. It is said to be a SNC (WSC) if it is a NC (SC), and for any other NC (SC) φ', we have that $\Gamma \models \varphi \rightarrow \varphi'$ ($\Gamma \models \varphi' \rightarrow \varphi$).*

It is seems that the SNC (WSC) of any formula can be obtained by changing to that of a proposition. Formally:

Proposition 10 *Let Γ be a formula, P, and α be as in Definition 5. A formula φ of P is the SNC (WSC) of α on P under Γ iff it is the SNC (WSC) of q on P under $\Gamma' = \Gamma \cup \{q \equiv \alpha\}$, where q is a new proposition not in Γ and α.*

We propose the theorem of computing the SNC (WSC) of an atom due to the SNC (WSC) of a formula can be changed to the SNC (WSC) of an atom by Proposition 10.

Theorem 8 *Let φ be a formula, $V \subseteq Var(\varphi)$ and $q \in Var(\varphi) - V$.*

(i) $\mathrm{F}_{\mathrm{CTL}}(\varphi \wedge q, (Var(\varphi) \cup \{q\}) - V)$ *is a SNC of q on V under φ.*

(ii) $\neg\mathrm{F}_{\mathrm{CTL}}(\varphi \wedge \neg q, (Var(\varphi) \cup \{q\}) - V)$ *is a WSC of q on V under φ.*

As we have said before that any initial K-structure can be characterized by a CTL formula, we can obtain the SNC (WSC) of an initial K-structure for satisfy some needed property (formula) by forgetting.

Theorem 9 *Let $\mathcal{K} = (\mathcal{M}, s)$ be a initial* K*-structure with $\mathcal{M} = (S, R, L, s_0)$ on the finite set $\mathcal{A}$ of atoms, $V \subseteq \mathcal{A}$ and $q \in V'$ ($V' = \mathcal{A} - V$). Then:*

(i) the SNC of q on V under $\mathcal{K}$ is $\mathrm{F}_{\mathrm{CTL}}(\mathcal{F}_{\mathcal{A}}(\mathcal{K}) \wedge q, V')$.

(ii) the WSC of q on V under $\mathcal{K}$ is $\neg\mathrm{F}_{\mathrm{CTL}}(\mathcal{F}_{\mathcal{A}}(\mathcal{K}) \wedge \neg q, V')$.

Example 3 For the Example 1, the WSC of φ on V under $\mathcal{K} = (\mathcal{M}, s_0)$ is $\neg\mathrm{F}_{\mathrm{CTL}}(\mathcal{F}_{\mathcal{A}}(\mathcal{K}) \wedge (q \equiv \varphi) \wedge \neg q, \mathcal{A} \setminus V)$.

5 Algorithm to compute forgetting

To compute the forgetting in CTL, we propose a model-based method in this part. Literally speaking, the model-based method means that we can obtain the result of forgetting in CTL by obtain all the possible finite models of this result. By the definition of forgetting in CTL, the set of models of the result of forgetting is also a finite set of initial K-structures.

Then we have the following model-based Algorithm 1 to computing the forgetting under CTL. By Lemma 3 and Theorem 3 we can prove the correctness of this algorithm.

Example 4 Let $\varphi = \mathrm{AGAF}(p \wedge r)$, $\mathcal{A} = \{p, r\}$ and $V = \{r\}$. For convenience, we use the label of a state to express the state and then remove the label function in a model structure. Let $\mathcal{M}_1 = (\{\{p, r\}\}, \{(\{p, r\}, \{p, r\})\}, \{p, r\})$ and $\mathcal{M}_2 = (\{\emptyset, \{p, r\}\}, \{(\emptyset, \{p, r\}), (\{p, r\}, \{p, r\})\}, \emptyset)$. The set of models of$\varphi$ is $Mod(\varphi) = \{(\mathcal{M}_1, \{p\}), (\mathcal{M}_2, \emptyset), \dots\}$.

Algorithm 1: Model-based: Computing forgetting

Input: A CTL formula φ and a set V of atoms
Output: $\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$

```
T = ∅ // the set of models of φ ;
T' = ∅ // the set of possible initial K-structures ;
n = |A|;
for i=1, ..., 2^n do
    for s_j ∈ {s_1, ..., s_i} do
        Let s_j be an initial state, construct
        M = (S, R, L, s_j) by the definition of model
        structure with S = {s_1, ..., s_i};
        if for each K ∈ T', (M, s_j) ↮_{Var(φ)} K then
            Let T' := T' ∪ {(M, s_j)};
        end
    end
    for (M, s_0) ∈ T' do
        if (M, s_0) ⊨ φ then
            T := T ∪ {(M, s_0)};
        end
    end
end
return ⋁_{(M', s'_0) ∈ T} F_V̄(M', s'_0).
```

Let $\mathcal{M}'_1 = (\{\{p\}\}, \{(\{p\}, \{p\})\}, \{p\})$ and $\mathcal{M}'_2 = (\{\emptyset, \{p\}\}, \{(\emptyset, \{p\}), (\{p\}, \{p\})\}, \emptyset)$ Then we can obtain all the possible initial K-structure that is a model of $\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$, *i.e.* $Mod(\mathrm{F}_{\mathrm{CTL}}(\varphi, V)) = \{\mathcal{K}_1 = (\mathcal{M}'_1, \{p\}), \mathcal{K}_2 = (\mathcal{M}'_2, \emptyset), \dots\}$.

Let $V' = \{p\}$, then $\mathcal{F}_{V'}(\mathcal{K}_1) = p \wedge \mathrm{AG}(p \supset \mathrm{EX}p \wedge \mathrm{AX}p)$, and $\mathcal{F}_{V'}(\mathcal{K}_2) = \neg p \wedge \mathrm{AG}(p \supset \mathrm{EX}\neg p \wedge \mathrm{AX}\neg p) \wedge \mathrm{AG}(\neg p \supset \mathrm{EX}p \wedge \mathrm{AX}p)$. Similarly, we can obtain the characteristic formula of other models and then the $\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$.

Proposition 11 *Let φ be a CTL formula and $V \subseteq \mathcal{A}$. The time complexity of Algorithm 1 is $O(2^{2^m})$ and the space complexity is $O(2^{2m})$, where $|\mathcal{A}| = m$.*

6 Concluding Remark

Based on the proposed V-bisimulation between K-structures, forgetting in CTL and characteristic formula on V on an initial K-structure $\mathcal{K}$, a method compute the WSC (SNC) of a property φ (a CTL formula) on $\mathcal{K}$ and V has been introduced by computing forgetting in CTL. Besides, we have shown that the CTL system is close under our definition of forgetting, and this definition satisfies those four postulates of forgetting. As we have said the complexity of Algorithm 1 is $O(2^{2^m})$ (very inefficient), a future work is to find an efficient algorithm to compute forgetting in CTL and then WSC (SNC).

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