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

Renyan Feng 1 , Yisong Wang 2* , Stefan Schlobach 3† and Erman Acar 4

^{1,2}GuiZhou University ^{3,4}Vrije Universiteit Amsterdam

fengrenyan@gmail.com, yswang@gzu.edu.cn, {k.s.schlobach, erman.acar}@vu.nl

Abstract

This article describes a method to compute the SNC (WSC) of a given property (a CTL formula) under a given transition system by proposing a semantic forgetting for CTL. We show that a CTL system is closed under our definition of forgetting, and this definition satisfies those four postulates of forgetting at first. And then we show the SNC and WSC can be computed by using the technology of forgetting. Besides, an algorithm model-based has been put forward to compute forgetting in CTL.

1 Introduction

Weakest precondition, we also called weakest sufficient condition (WSC), is introduced by Dijkstra in [Dijkstra, 1978]. Strongest postcondition (we also called 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. It has been extensively studied in active system [Mraihi *et al.*, 2011; Bruggink *et al.*, 2011].

Model checking is an automatic, model-based, propertyverification approach. It is intended to be used for concurrent, reactive systems and originated as a post-development methodology. In model checking, specification is mainly used to specify the properties that should be satisfied by a system. CTL, one of the main logical formalisms for program specification and verification, has been proposed by Clarke and Emerson [Clarke and Emerson, 1981]. Given a transition system \mathcal{M} with s_0 as an initial state and a specification φ (a CTL formula), we should decide wether $(\mathcal{M}, s_0) \models \varphi$. It is a good thing if $(\mathcal{M}, s_0) \models \varphi$ indeed. However, if $(\mathcal{M}, s_0) \nvDash \varphi$, how can we find the WCS ψ under a given set of atoms such that $(\mathcal{M}, s_0) \models \psi \supset \varphi$. This means that how can we find such a condition such that $(\mathcal{M}, s_0) \models \varphi$ after we repairing the system by adding this condition to consider. It can be show 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 $S = \{s_0, s_1, s_2\}, R = \{(s_0, s_1), (s_1, s_2), (s_1, s_0), (s_2, s_0)\}, L(s_0) = \{select\}, L(s_1) = \{pay, soda\}, L(s_2) = \{pay, beer\}$ 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 then we change 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, pay, pay, pay, pay, pay, pay, pay,

We can decide $(\mathcal{M}, s_0) \nvDash \varphi$ 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 add to (\mathcal{M}, s_0) . As we know that if this condition exists, there are many, 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 property. 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 these properties is the same as discovering theorems said by Lin in [Lin, 2018].

However, as we have said that \mathcal{M} is a model structure, which needs to be converted into a logical formula (theory), that is the characteristic formula - a CTLformula - we will introduce in the paper. At the same time the property what we need is a formula defined on the specified set V. Inspired by the forgetting-based method to compute SNC (WSC) [Lin, 2001], in this paper, we tackle those problem by proposing a semantic forgetting for CTL.

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, 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], forget-

^{*}Contact Author

[†]Contact Author

ting 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.

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

- \bullet 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 $S \to 2^{\mathcal{A}}$.

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 $j \ (0 \le i \le j), \ (s_j, s_{j+1}) \in R$. 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 $\mathrm{Tr}_n^{\mathcal{M}}(s)$ of $\mathcal{M}(\mathrm{or\ simply\ }\mathrm{Tr}_n(s))$, that has

depth n and is rooted at s, is recursively defined as [Browne *et al.*, 1988], for $n \ge 0$,

- $Tr_0(s)$ consists of a single node s with label s.
- $\operatorname{Tr}_{n+1}(s)$ has as its root a node m with label s, and if $(s,s')\in R$ then the node m has a subtree $\operatorname{Tr}_n(s')$.

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

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 *Computation Tree Logic* (CTL in short) [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 , \lor 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:

$$\phi ::= \bot \mid p \mid \neg \phi \mid \phi \lor \phi \mid \mathsf{EX}\phi \mid \mathsf{EG}\phi \mid \mathsf{E}[\phi \cup \phi] \tag{1}$$

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):

$$\neg$$
, EX, EF, EG, AX, AF, AG $\prec \land \prec \lor \prec$ EU, AU, EW, AW, \rightarrow .

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:

- $(\mathcal{M}, s) \not\models \bot$;
- $(\mathcal{M}, s) \models p \text{ iff } p \in L(s);$
- $(\mathcal{M}, s) \models \phi_1 \lor \phi_2$ iff $(\mathcal{M}, s) \models \phi_1$ or $(\mathcal{M}, s) \models \phi_2$;
- $(\mathcal{M}, s) \models \neg \phi \text{ iff } (\mathcal{M}, s) \not\models \phi;$
- $(\mathcal{M}, s) \models \text{EX}\phi \text{ iff } (\mathcal{M}, s_1) \models \phi \text{ for some } s_1 \in S \text{ and } (s, s_1) \in R;$
- $(\mathcal{M}, s) \models EG\phi$ iff \mathcal{M} has a path $(s_1 = s, s_2, ...)$ such that $(\mathcal{M}, s_i) \models \phi$ for each $i \geq 1$;
- $(\mathcal{M}, s) \models E[\phi_1 U \phi_2]$ iff \mathcal{M} has a path $(s_1 = s, s_2, ...)$ 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

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

 $\mathcal{K} \models \phi$. Let Π be a set of formulae, $\mathcal{K} \models \Pi$ if for each $\phi \in \Pi$ there is $\mathcal{K} \models \phi$. We denote $Mod(\phi)$ ($Mod(\Pi)$) the set of models of ϕ (Π). The formula ϕ (set Π of formulae) is *satisfiable* if $Mod(\phi) \neq \emptyset$ ($Mod(\Pi) \neq \emptyset$). Since signatures (atoms) in model structure is finite, $Mod(\phi)$ ($Mod(\Pi)$) is finite for any formula ϕ (set Π of formulae).

Let ϕ_1 and ϕ_2 be two formulas or set of formulas. By $\phi_1 \models \phi_2$ we denote $Mod(\phi_1) \subseteq Mod(\phi_2)$. 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 $Var(\phi_1)$ we mean the set of atoms occurring in ϕ_1 . ϕ_1 is V-irrelevant, written $IR(\phi_1, V)$, if there is a formula ψ with $Var(\psi) \cap V = \emptyset$ 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 knowledge forgetting, we need the concepts of V-bisimulation, which is similar with that in [Zhang and Zhou, 2009]. 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 $\mathcal{K}_i=(\mathcal{M}_i,s_i)$ and $\mathcal{M}_i=\langle S_i,R_i,L_i,s_0^i\rangle$ with $i\in\{1,2\}$,

- $(K_1, K_2) \in \mathcal{B}_0$ 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 $V \subseteq A$. The *V*-bisimular relation 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, K_1 and K_2 are called V-bisimular. 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.

Proposition 1 Let $V \subseteq A$ and $K_i = (M_i, s_i)$ (i = 1, 2) be K-structures. Then $(K_1, K_2) \in \mathcal{B}$ 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 vice versa.

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

$$(\mathcal{K}_{1,j},\mathcal{K}_{2,j}) \in \mathcal{B}$$
 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 $i \in \{1,2\}$, $V_1, V_2 \subseteq \mathcal{A}$, $s_i's$ be two states and $\pi_i's$ 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) $s'_1 \leftrightarrow_{V_i} s'_2 \ (i=1,2) \text{ implies } s'_1 \leftrightarrow_{V_1 \cup V_2} s'_2;$
- (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 $\mathcal{K}_1 \leftrightarrow_{V_2} \mathcal{K}_2$

Let $\mathcal{M}=(S,R,L,s_0)$ be a model structure on a finite set \mathcal{A} of atoms, $V\subseteq\mathcal{A}$ and $\mathcal{B}=\{(\mathcal{M},s)\leftrightarrow_V(\mathcal{M},s')|s,s'\in S\}$. For $s\in S$, $[s]_{\mathcal{B}}$ denotes the equivalence class of state s under \mathcal{B} , i.e. , $[s]_{\mathcal{B}}=\{s'\in S|(s,s')\in \mathcal{B}\}$. Note that for $s'\in [s]_{\mathcal{B}}$ we have $[s']_{\mathcal{B}}=[s]_{\mathcal{B}}$. The set $[s]_{\mathcal{B}}$ is referred to as the \mathcal{B} -equivalence class of s. The V-quotient space of S under \mathcal{B} , denoted by $S/\mathcal{B}=\{[s]_{\mathcal{B}}|s\in S\}$, is the consisting of all \mathcal{B} -equivalence classes.

Definition 2 For K-structure $\mathcal{K} = (\mathcal{M}, s_0)$ with $\mathcal{M} = (S, R, L, s_0)$ a model structure on a finite set \mathcal{A} of atoms, $V \subseteq \mathcal{A}$ and $\mathcal{B} = \{(\mathcal{M}, s) \leftrightarrow_{V'} (\mathcal{M}, s') | s, s' \in S\}$ where $V' = \mathcal{A} - V$, the V-quotient structure is $\mathcal{K}_{|V} = (\mathcal{M}^*, s_0^*)$ with $\mathcal{M}^* = (S^*, R^*, L^*, s_0^*)$ on V, where

- $s_0^* = [s_0]_{\mathcal{B}}, S^* = S/\mathcal{B}, L^*([s]_{\mathcal{B}}) = L(s) \cap V$, and
- $R^* = \{([s]_{\mathcal{B}}, [s']_{\mathcal{B}}) | \exists s_1 \in [s]_{\mathcal{B}} \text{ s.t. } \exists s_2 \in [s']_{\mathcal{B}} \text{ and } (s_1, s_2) \in R\}.$

Example 2 Let \mathcal{K} be the initial structure in Example 1 and $V = \{p, r\}$ ($V' = V_a \setminus V$), then we can obtain $\mathcal{K}_V = (\mathcal{M}^*, s_0^*)$ with $\mathcal{M}^* = (S^*, R^*, L^*, s_0^*)$ due to $\mathcal{B}_V = \{(s_1, s_2)\}$, where $S^* = \{s_0' = [s_0]_{\mathcal{B}_V}, s_1' = [s_1]_{\mathcal{B}_V}\}$, $s_0^* = s_0'$, $R^* = \{(s_0', s_1'), (s_1', s_0')\}$, and $L^*(s_0') = \{\}$, $L^*(s_1') = \{p\}$.

Proposition 3 For any K-structure $\mathcal{K} = (\mathcal{M}, s_0)$ with $\mathcal{M} = (S, R, L, s_0)$ a model structure on a finite set \mathcal{A} of atoms and $V \subseteq \mathcal{A}$, it holds that $\mathcal{K} \leftrightarrow_{V'} \mathcal{K}_{|V|}$ where $V' = \mathcal{A} - V$.

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. $\mathrm{IR}(\varphi,V)$.

Theorem 1 Let $V \subseteq A$, K_i (i = 1, 2) be two K-structures such that $K_1 \leftrightarrow_V K_2$ and ϕ a formula with $IR(\phi, V)$. Then $K_1 \models \phi$ if and only if $K_2 \models \phi$.

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

- $L_1(s_1) V = L_2(s_2) V$,
- for every subtree $\operatorname{Tr}_{n-1}(s_1')$ of $\operatorname{Tr}_n(s_1)$, $\operatorname{Tr}_n(s_2)$ has a subtree $\operatorname{Tr}_{n-1}(s_2')$ such that $\operatorname{Tr}_{n-1}(s_1') \leftrightarrow_V \operatorname{Tr}_{n-1}(s_2')$, and vice verse.

Please note that the last two conditions in the above definition hold trivially for n=0.

Proposition 4 Let $V \subseteq A$ and (M_i, s_i) (i = 1, 2) be two K-structures. Then

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

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

Proposition 5 Let $V \subseteq A$, M be a model structure and $s, s' \in S$ such that $(s, s') \notin B$. 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 $\operatorname{dis}_V(\mathcal{M},s,s',k)$. It is evident that $\operatorname{dis}_V(\mathcal{M},s,s',k)$ implies $\operatorname{dis}_V(\mathcal{M},s,s',k')$ whenever $k' \geq k$. The V-characterization number of \mathcal{M} , written $\operatorname{ch}(\mathcal{M},V)$, is defined as

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

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

Definition 3 (Forgetting) *Let* $V \subseteq \mathcal{A}$ *and* ϕ *a formula. A formula* ψ *with* $Var(\psi) \cap V = \emptyset$ *is a* result of forgetting V from ϕ , *if*

$$Mod(\psi) = \{ \mathcal{K} \text{ is initial } | \exists \mathcal{K}' \in Mod(\phi) \& \mathcal{K}' \leftrightarrow_V \mathcal{K} \}.$$
 (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, 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 $\mathit{Var}(\varphi)\subseteq V$) in CTL to equivalent uniquely describe a computation tree.

Definition 4 Let $V \subseteq A$, $M = (S, R, L, s_0)$ 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:

$$\mathcal{F}_{V}(\mathit{Tr}_{0}(s)) = \bigwedge_{p \in V \cap L(s)} p \wedge \bigwedge_{q \in V - L(s)} \neg q,$$

$$\mathcal{F}_{V}(\mathit{Tr}_{k+1}(s)) = \bigwedge_{(s,s') \in R} \mathsf{E} \mathsf{X} T(s') \wedge \mathsf{A} \mathsf{X} \bigvee_{(s,s') \in R} T(s') \wedge \mathcal{F}_{V}(\mathit{Tr}_{0}(s))$$

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 can 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 $\operatorname{Tr}_n(s)$ then $\varphi \equiv \mathcal{F}_V(\operatorname{Tr}_n(s))$.

Let $V\subseteq \mathcal{A}$, $\mathcal{K}=(\mathcal{M},s_0)$ be an initial K-structure and $T(s')=\mathcal{F}_V(\operatorname{Tr}_c(s'))$. 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:

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

$$\operatorname{AG}\left(\mathcal{F}_{V}(\operatorname{Tr}_{c}(s)) \to \bigwedge_{(s,s') \in R} \operatorname{Ex}T(s') \wedge \operatorname{Ax} \bigvee_{(s,s') \in R} T(s')\right)$$

where $c = ch(\mathcal{M}, V)$. It is apparent that $IR(\mathcal{F}_V(\mathcal{M}, s_0), \overline{V})$. We will give an example to show the computing of characterizing formula:

Example 3 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 $\operatorname{Tr}_0(s_0) \leftrightarrow_{\overline{V}} \operatorname{Tr}_0(s_1)$ due to $L(s_0) - \overline{V} = L(s_1) - \overline{V}$, $\operatorname{Tr}_1(s_0) \not \leftrightarrow_{\overline{V}} \operatorname{Tr}_1(s_1)$ 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.* $\operatorname{dis}_{\overline{V}}(\mathcal M, s_0, s_1, 1)$. Similarly, we have $\operatorname{dis}_{\overline{V}}(\mathcal M, s_0, s_2, 0)$ and $\operatorname{dis}_{\overline{V}}(\mathcal M, s_1, s_2, 0)$. Therefore, $\operatorname{ch}(\mathcal M, \overline{V}) = \max\{k \mid s, s' \in S \ \& \operatorname{dis}_{\overline{V}}(\mathcal M, s, s', k)\} = 1$. Then we have:

$$\mathcal{F}_V(\operatorname{Tr}_0(s_0)) = a \wedge \neg b,$$

$$\mathcal{F}_V(\operatorname{Tr}_0(s_1)) = a \wedge \neg b,$$

$$\mathcal{F}_V(\operatorname{Tr}_0(s_2)) = b \wedge \neg a,$$

$$\mathcal{F}_V(\operatorname{Tr}_1(s_0)) = \operatorname{EX}(a \wedge \neg b) \wedge \operatorname{EX}(b \wedge \neg a) \wedge \operatorname{AX}((a \wedge \neg b) \vee (b \wedge \neg a)) \wedge (a \wedge \neg b),$$

$$\mathcal{F}_V(\operatorname{Tr}_1(s_1)) = \operatorname{EX}(a \wedge \neg b) \wedge \operatorname{AX}(a \wedge \neg b) \wedge (a \wedge \neg b),$$

$$\mathcal{F}_V(\operatorname{Tr}_1(s_2)) = \operatorname{EX}(a \wedge \neg b) \wedge \operatorname{AX}(a \wedge \neg b) \wedge (b \wedge \neg a).$$

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

Lemma 3 Let φ be a formula. We have

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

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 $V \subseteq A$, $M = (S, R, L, s_0)$ be a model structure with initial state s_0 and $M' = (S', R', L', s'_0)$ 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 3, the result ψ of ϕ forget the set V of atoms always exists, which is equivalent to

$$\bigvee$$
 $\mathcal{F}_{\overline{V}}(\mathcal{K}).$

 $\mathcal{K} \in \{\mathcal{K}' \text{ is an initial interpretation} | \exists \mathcal{K}'' \in Mod(\phi) \& \mathcal{K}'' \leftrightarrow_V \mathcal{K}' \}$

For this reason, the forgetting result is denoted by $F_{CTL}(\phi, V)$. 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 $V \subseteq A$. Then the following statements are equivalent:

- (i) $\psi \equiv \mathbf{F}_{\mathrm{CTL}}(\varphi, V)$,
- (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 closed, *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 $F_{CTL}(\varphi \cup \{q \leftrightarrow \alpha\}, q) \equiv \varphi$.

Proposition 6 Let φ be a formula, V a set of atoms and p an atom such that $p \notin V$. Then:

$$F_{CTL}(\varphi, \{p\} \cup V) \equiv F_{CTL}(F_{CTL}(\varphi, p), V).$$

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:

$$F_{\text{CTL}}(\varphi, V_1 \cup V_2) \equiv F_{\text{CTL}}(F_{\text{CTL}}(\varphi, V_1), V_2).$$

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

Proposition 7 Let φ , φ_i , ψ_i (i = 1, 2) be formulas and $V \subseteq \mathcal{A}$. We have

- (i) $F_{CTL}(\varphi, V)$ is satisfiable iff φ is;
- (ii) If $\varphi_1 \equiv \varphi_2$, then $F_{CTL}(\varphi_1, V) \equiv F_{CTL}(\varphi_2, V)$;
- (iii) If $\varphi_1 \models \varphi_2$, then $F_{CTL}(\varphi_1, V) \models F_{CTL}(\varphi_2, V)$;
- (iv) $F_{CTL}(\psi_1 \vee \psi_2, V) \equiv F_{CTL}(\psi_1, V) \vee F_{CTL}(\psi_2, V)$;
- (v) $F_{CTL}(\psi_1 \wedge \psi_2, V) \models F_{CTL}(\psi_1, V) \wedge F_{CTL}(\psi_2, V)$;

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

Proposition 8 *Let* $V \subseteq A$ *and* ϕ *a formula.*

- (i) $F_{CTL}(AX\phi, V) \equiv AXF_{CTL}(\phi, V)$.
- (ii) $F_{CTL}(EX\phi, V) \equiv EXF_{CTL}(\phi, V)$.
- (iii) $F_{CTL}(AF\phi, V) \equiv AFF_{CTL}(\phi, V)$.
- (iv) $F_{CTL}(EF\phi, V) \equiv EFF_{CTL}(\phi, V)$.

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

- ψ is a necessary condition (NC in short) of q on V under φ if φ |= q → ψ.
- 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 $\phi \models \psi' \to \psi$ 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 5.

 ψ is a SNC of q on V under φ iff ¬ψ is a WSC of q on V under φ. (ii) ψ is a WSC of q on V under φ iff $\neg \psi$ is a SNC of 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 6 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 6. 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 change to the SNC (WSC) of an atom by Proposition 10.

Theorem 6 Let φ be a formula, $V \subseteq Var(\varphi)$ and $q \in Var(\varphi) - V$.

- (i) $F_{CTL}(\varphi \wedge q, (Var(\varphi) \cup \{q\}) V)$ is a SNC of q on V under φ .
- (ii) $\neg F_{CTL}(\varphi \wedge \neg q, (Var(\varphi) \cup \{q\}) V)$ 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 7 Let K = (M, s) be a initial K-structure with M = (S, R, L) on the finite set A of atoms, $V \subseteq A$ and $q \in V'$ (V' = A - V). Then:

- (i) the SNC of q on V under K is $F_{CTL}(\mathcal{F}_{V \cup \{q\}}(K_{|V \cup \{q\}}) \land q, q)$.
- (ii) the WSC of q on V under \mathcal{K} is $\neg F_{CTL}(\mathcal{F}_{V \cup \{q\}}(\mathcal{K}_{|V \cup \{q\}}) \land \neg q, q)$.

Example 4 For the Example 1, the WSC of φ on V under $\mathcal{K} = (\mathcal{M}, s_0)$ is $\neg F_{\text{CTL}}(\mathcal{F}_V(\mathcal{K}_{|V}) \land (q \equiv \varphi) \land \neg q, q)$.

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 let φ be a CTL formula, the $|\mathit{Var}(\varphi)| = m$ is a positive integer, we have the following theorem:

Theorem 8 Let φ be a CTL formula, $V = Var(\varphi)$, |V| = m, and $\mathcal{K} = (\mathcal{M}, s_0)$ with $\mathcal{M} = (S, R, L, s_0)$ be a initial K-structure. If $(\mathcal{M}, s_0) \models \varphi$, then there is an initial K-structure $\mathcal{K}' = (\mathcal{M}', s_0')$ with $\mathcal{M}' = (S', R', L', s_0')$ that satisfy:

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Algorithm 1: Model-based: Computing forgetting
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Input: A CTL formula \varphi and a set V of atoms Output: F_{CTL}(\varphi, V)

1 T = \emptyset // the set of models of \varphi;

2 T' = \emptyset // the set of models of F_{CTL}(\varphi, V);

3 m = |Var(\varphi)|;

4 n = |V|;

5 for i = 1, ..., 2^m do

6 | Enumerating all possible initial K-structures (\mathcal{M}, s_0) with \mathcal{M} = (S, R, L, s_0) and |S| = i;

7 | For all initial K-structures (\mathcal{M}, s_0) if (\mathcal{M}, s_0) \models \varphi then

8 | T = T \cup \{(\mathcal{M}, s_0)\};

9 | end

10 end

11 return \bigvee_{(\mathcal{M}', s_0') \in T} \mathcal{F}_{\overline{V}}(\mathcal{M}', s_0').
```

- (i) |S'| is at most 2^m ,
- (ii) |R'| is at most $2^m * 2^m$,
- (iii) |L'| is at most $2^m * 2^m$,
- (iv) $\mathcal{K} \leftrightarrow_{\mathcal{A} \setminus V} \mathcal{K}'$ and $(\mathcal{M}', s_0') \models \varphi$.

By this theorem, we can see that any initial K-structures $\mathcal K$ that satisfy φ can be transformed to an initial K-structure $\mathcal K'$ such that $\mathcal K \leftrightarrow_{\mathcal A\setminus V} \mathcal K'$ and $\mathcal K\models \varphi$ iff $\mathcal K'\models \varphi$ due to $\mathrm{IR}(\varphi,\mathcal A\setminus V)$, in which $V=\mathit{Var}(\varphi)$. Therefore, the size of the model of φ is at most 2^m by Theorem 3(we only consider the number of states of this model).

Then we have the following model-based Algorithm 1 to computing the forgetting under CTL:

Example 5 Let $\varphi = \text{AGAF}(p \lor r)$ 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 $\mathcal{M}_1 = (\{\{p,r\}\}, \{(\{p,r\}, \{p,r\})\}, \{p,r\}), \mathcal{M}_2 = (\{\{\}, \{p,r\}\}, \{(\{\}, \{p,r\}\}, (\{p,r\}, \{p,r\})\}, \{\}), \mathcal{M}_3 = (\{\{\}, \{p\}, \{p,r\}\}, \{(\{\}, \{p\}), (\{p\}, \{p,r\}), (\{p,r\}, \{\})\}, \{\}), \mathcal{M}_3 = (\{\{p\}, (\mathcal{M}_3, \{\}\}, \dots\})$ Let $\mathcal{M}_1' = (\{\{p\}\}, \{(\{p\}, \{p\})\}, \{p\})$ and $\mathcal{M}_2' = (\{\{p\}\}, \{p\}), \{p\})$. Then we can obtain

Let $\mathcal{M}_1' = (\{\{p\}\}, \{(\{p\}, \{p\})\}, \{p\}) \text{ and } \mathcal{M}_2' = (\{\{\}, \{p\}\}, \{(\{\}, \{p\}), (\{p\}, \{p\})\}, \{\}) \text{ Then we can obtain all the possible initial structure that is a model of } F_{CTL}(\varphi, V), i.e. <math>Mod(F_{CTL}(\varphi, V)) = \{\mathcal{K}_1 = (\mathcal{M}_1', \{p\}), \mathcal{K}_2 = (\mathcal{M}_2', \{\}), \ldots\}.$

Then the characteristic formula of \mathcal{K}_1 is $\psi_1 = p \wedge \mathrm{AG}(p \supset \mathrm{EX} p \wedge \mathrm{AX} p)$, and the characteristic formula of \mathcal{K}_2 is $\psi_2 = \neg p \wedge \mathrm{AG}(p \supset \mathrm{EX} \neg p \wedge \mathrm{AX} \neg p) \wedge \mathrm{AG}(\neg p \supset \mathrm{EX} p \wedge \mathrm{AX} p)$. Similarly, we can obtain the characteristic formula of other models. Then $\mathrm{F}_{\mathrm{CTL}}(\varphi, V)$ is the disjunction of all the characteristic formulae of those models.

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

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