

Technische Universität Dresden • Fakultät Informatik

Data Preperation for PMC-Visualization

Bachelorarbeit zur Erlangung des ersten
Hochschulgrades

Bachelor of Science (B.Sc.)

vorgelegt von

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(geboren am 7. April 1999 in HALLE (SAALE))

Tag der Einreichung: July 28, 2023

Dipl. Inf Max Korn (Theoretische Informatik)

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Abstract

Lorem ipsum

1 Introduction

PMC is short for Probabilistic Model Checking.

Model Checking is... Probabilistic means...

When analyzing PMCs a result table is generated which consists of result vectors. These result vectors for example contain probability values for temporal events, expected values of random variables and other qualitative and quantitative results. For central measure so called schedulers can be used to resolve non-determinism in the MDPs which in addition to the result provide one or more optimal actions per state.

Already simple system models can lead to complex system behavior and an immense amount of states. An interactive visualization of an MDP can be used in support of understanding, analyzing and reviewing these complex systems. It can even be made use of to achieve further goals such as debugging or model repair. However, the pure data volume remains unaffected. The complexity is shifted to the visualization which has some general techniques to represent large amounts of data in a suitable way.

WHAT EXACTLY CAN BE DONE/HAS BEEN DONE —; INSERT HERE

Apart from the methods available in visualization there exist domain specific abstraction methods for transition systems. These are based on relations such as simulation order, simulation equivalence, bisimulation equivalence as well as trace equivalence, stutter linear time relations and stutter bisimulation equivalences. These relations can be based on the set of all possible transition systems or the set of states of one single transition system. Whereas the first declares some similarity of transition systems with regard to model check the latter one can be used to obtain quotient systems. The set of states of such quotient system coincides with the the respective equivalence relation.

- all relations declare that at least one of the two systems being in relation can mimic the behavior of the other one in some way. In case of equivalence the systems can mimic each other.
- except for trace inclusion and trace equivalence whether such mimicking is possible depends on the existence of a specific relation $\mathcal{R} \subseteq S_1 \times S_2$ where S_1 and S_2 are the state sets of two corresponding transition systems TS_1 and TS_2 . The symbol \mathcal{R} is included in each definition of a relation on the set of all possible transition systems. In each definition it receives a different name and represents a different relation between the sets S_1 and S_2 .
- Depending on the type of mimicking (simulation order, bisimulation equivalence, ...) different requirements are made on \mathcal{R} . It is dependent on these requirements whether \mathcal{R} exists for a given pair of transition systems TS_1 and TS_2 and hence also whether or not these two transition systems are in relation.
- all relations have three requirements made on \mathcal{R} in common.
- firstly, there has to be some relationship between the initial states. That is, at least $\exists(s_1, s_2) \in \mathcal{R}$ with $s_1 \in I_1, s_2 \in I_2$ where I_1 and I_2 are the respective sets of initial states of the two transition systems TS_1 and TS_2 in question

- Secondly for each $(s_1, s_2) \in \mathcal{R}$ the two states must have the same set of atomic propositions. That is $L(s_1) = L(s_2)$.
- lastly for all s_1, s_2 with $(s_1, s_2) \in \mathcal{R}$ some requirement is made about their (not necessarily direct) successors
- the mimicking types differ in the requirements they make on the relation \mathcal{R} with respect to initial states (1) of the two transition systems and the requirements made on the successors (2).
- when one transition system can mimic another it is only required that that each initial state $s_1 \in I_1$ there is an initial state $s_2 \in I_2$ with $(s_1, s_2) \in \mathcal{R}$
- when two systems mimic each other it is as well required that for each initial state $s_1 \in I_1$ it is $(s_1, s_2) \in \mathcal{R}$ with $s_2 \in I_2$ but additionally that for each $s_2 \in I_2$ it is $(s_2, s_1) \in \mathcal{R}$ with $s_1 \in I_1$.
- where all of them differ is the requirements made on the successors of the states being in relation
- bisimulation equivalence (\sim) declares that two transition systems can mimic each other. Every two states with $(s_1, s_2) \in \mathcal{R}$ one must have an successor s' for each successor \tilde{s}' of the other state so that the two successors stand in relation with respect to \mathcal{R} . That is the successors standing in relation again must have the same set of atomic propositions and meet the requirement regarding their successors. If a state s is equal in its set of atomic propositions to another state \tilde{s} and meets a respective requirement on a selection of the successors of \tilde{s} we say it state-mimics it. Note that here state-mimicking is mutual. s_1 state-mimics s_2 and vice versa. **definition?** bisimulation equivalence is, as the name suggest, an equivalence relation on the set of all possible transition systems.
- simulation order (\preceq) is weaker than bisimulation. It only declares that one system can state-mimic the other. For every two states with $(s_1, s_2) \in \mathcal{R}$ the second one must state-mimic the first one. That is, for each successor s'_1 of s_1 the state s_2 must have an successor s'_2 so that s_2 *state-mimics* s'_1 . In this case of the simulation order relation s'_2 state-mimicking s'_1 coincides with $(s'_1, s'_2) \in \mathcal{R}$. For the simulation order relation it is not required that s_1 can state-mimic s_2 . Simulation order is a preorder on the set of all possible transition systems.
- Simulation equivalence (\simeq) declares that for two transition systems TS_1, TS_2 it holds that $TS_1 \preceq TS_2$ and $TS_2 \preceq TS_1$. It might seem like it coincides with bisimulation but it does not. For simulation order it has to hold that for a given state s_1 in the mimicked system that there has to exist another state s_2 in the mimicking system, that state-mimics s_1 . Although due to simulation equivalence s_2 now also must have a state s' that mimics s_2 , it does not need to hold that $s' = s_1$. The state s_2 could be mimicked by another state than s_1 . The mimicking between two states does not need to be mutual. Simulation equivalence is an equivalence relation on the set of all possible transition systems.

- in general it holds that for TS_1, TS_2 being transition systems $TS_1 \sim TS_2$ implies $TS_1 \simeq TS_2$ implies $TS_1 \preceq TS_2$.
- All of the mimicking types presented, mimicking is happening in a stepwise manner and atomic propositions are considered as the property of concern. Choosing atomic propositions has the advantage of preserving the logical formulae used in model checking.
- bisimulation equivalence preserves CTL and CTL* formulae for finite TS without terminal states 579
- simulation equivalence/order maintain LTL by trace inclusion and universally and existentially quantified fragments of CTL* **correct?** 579
- The condition of stepwise mimicking is softened with stutter simulation and bisimulation equivalence expands the concept of simulation equivalence and stutter bisimulation equivalence. Instead of requiring for each successor s'_1 of a state s_1 to be state-mimicked by a direct successor of s_2 it suffices if there is path to such a state s'_2 , that state-mimics s'_1 . All states on the path must then also state-mimic s_1 . Moreover this requirement is only made for successors of s_1 if the successor s'_1 is not in relation to s_2 .
- There also is another variant of stutter bisimulation equivalence which is called stutter bisimulation divergent equivalence which we wont consider here.
- So far relations on the set of all transition systems have been considered.
- All the above mentioned equivalence relations are also defined on the set of states of one single transition system.
- They are defined very similar to the relations on the set of all possible transition systems. Let $\bullet_{TS} \subseteq S \times S$ be some relation on the set of states S of an transition system. Two states $s_1, s_2 \in S$ are in relation $((s_1, s_2) \in \bullet_{TS})$ if there exists a relation $\mathcal{R} \subseteq S \times S$ that meets some requirements. These requirements are the same as the ones used on \mathcal{R} for transition systems except for the condition on the initial states. That is, only the set of atomic propositions of s_1, s_2 has to be equal ($L(s_1) = L(s_2)$) and some requirement on their successors has to be met.
- They are used for abstractions by defining quotient systems of a given transition system. In a quotient systems the set of states consists of the equivalence classes of the equivalence relation.

Definition 1.1. Let $TS = (S, Act, \longrightarrow, I, AP, L)$ be a transition system and \bullet_{TS} and equivalence relation on S . The quotient transition system is defined by $TS/\bullet_{TS} = (S/\bullet_{TS}, \{\tau\}, \longrightarrow', I', AP, L')$ where:

- $I' := \{[s]_{\bullet_{TS}} \mid s \in I\}$
- $\longrightarrow' := \{([s]_{\bullet_{TS}}, \tau, [s']_{\bullet_{TS}}) \mid (s, \alpha, s') \in \longrightarrow\}$

- $L'([s]_{\bullet_{TS}}) := L(s)$
- holds $TS \bullet TS / \bullet_{TS}$
- Similarly an abstract transition system using the simulation relation can be obtained... using an abstraction function
- **def abstraction function, def relation**
- obtain smaller system that has properties with respect to formulas **discuss how much smaller? → would rather not**
- expensive: to calculate with complexities
 - bisimulation:
 - simulation
 - stutter simulation:
 - stutter bisimulation:
- usecase: browsability for humans
 - understand and looking at the system from a human perspective of highest importance
 - less expensive calculation
- views of an MDP aim for high comprehensibility, low performance and will conceptually be very similar to the the notion of an abstract transition system.
- **”Partial order reduction attempts to analyze only a fragment \hat{TS} of the full transition system TS by ignoring several interleavings of independent actions.”**

In alternative to these general approaches domain specific preprocessing can be used to gain certain views on a given MDP that display areas or sets of states in a summarized more compact form. For this purpose structural properties as well as criteria obtained from the result vectors and their containing optimal actions can be utilized.

This thesis is about the development, formalization and implementation of a collection of different *views*. **Views are supposed to be applicable to only a subset of a given MDP and be composable with other views** The input for a view is the result table of the MDP, whereas the output is a new column in the MDP result table that determines which states are to be grouped to a new one. The evaluation is performed within the existing web-based prototype of a PMC visualization platform (PMC-Viz) and on the basis of different MDP models. They are to be analyzed regarding the suitability to facilitate the understanding of complex operational models incl. analysis results and to support processes of debugging, model repair or strategy synthesis.

2 Preliminaries

Views will be defined on MDP. Instead of directly providing the definition of MDP we will consider less powerful classes of models to represent systems that are extended by MDPs.

Firstly we will consider transition systems. Transition systems are basically digraphs consisting of *states* and *transitions* in place of nodes and edges. A state describes some information about a system at a certain moment of its behavior. Considering a traffic light displaying green could be considered as one state whereas displaying red could be another one. Transitions model the progression of the system from one state to another one. Sticking to the example of the traffic light a transition could model the switch from state green light being displayed to the state of red light being displayed. There are several variants of transitions systems. We will use transition systems with *action names* and *atomic propositions* for states as in **BAIER**. Actions are used for communication between processes. We denote them with Greek letters $(\alpha, \beta, \gamma, \dots)$. Atomic propositions are simple facts about states. For instance "x is greater than 20" or "red and yellow light are on" could be atomic propositions. They are assigned to a state by a Labeling function L .

The following definition is directly taken form Principles of Modelchecking, Baier p. 20

Definition 2.1. A *transition system* TS is a tuple $(S, Act, \longrightarrow, I, AP, L)$ where

- S is a set of states,
- Act is a set of actions,
- $\longrightarrow \subseteq S \times Act \times S$ is transition relation,
- $I \subseteq S$ is a set of initial states,
- AP is a set of atomic propositions, and
- $L : S \rightarrow \mathcal{P}(AP)$

A transition system is called *finite* if S , AP and L are finite. The intuitive behavior of transition systems is as follows. The evolution of a transition system starts in some state $s \in I$. If the set of I of initial states is empty the transition system has no behavior at all. From the initial state the transition system evolves according to the transition relation \longrightarrow . The evolution ends in a state that has no outgoing transitions. For every state there may be several possible transitions to be taken. The choice of which one is take is done nondeterministically. That is the outcome of the selection can not be know a priori. It is especially not following any probability distribution. Hence there can not be made any statement about the likelihood of a transition being selected.

In contrast with Markov Chains this nondeterministic behavior is replaced with a probabilistic one. That is for every state there exists a probability distribution that describes the chance of a transition of being selected. It is important to note that Markov Chains are memoryless in the sense that the system evolution is not

dependent on the history of only on the current state. That is the evolution of the system does not depend on the sequence of so far traversed states with the transitions. That the nondeterministic behavior is replaced with a probabilistic one also means that there are no actions in Markov chains.

Definition 2.2. A (*discrete-time*) *Markov chain* is a tuple $\mathcal{M} = (S, \mathbf{P}, \iota_{init}, AP, L)$ where

- S is a countable, nonempty set of states,
- $\mathbf{P} : S \times S \rightarrow [0, 1]$ is the *transition probability function*, such that for all states s :

$$\sum_{s' \in S} \mathbf{P}(s, s') = 1.$$

- $\iota_{init} : S \rightarrow [0, 1]$ is the *initial distribution*, such that $\sum_{s \in S} \iota_{init}(s) = 1$, and
- AP is a set of atomic propositions and,
- $L : S \rightarrow \mathcal{P}(AP)$ a labeling function.

A Markov chain \mathcal{M} is called *finite* if S and AP are finite. For finite \mathcal{M} , the *size* of \mathcal{M} , denoted $size(\mathcal{M})$, is the number of states plus the number of pairs $(s, s') \in S \times S$ with $\mathbf{P}(s, s') > 0$. OMIT?? The probability function \mathbf{P} specifies for each state s the probability $\mathbf{P}(s, s')$ of moving from s to s' in one step. The constraint put on \mathbf{P} in the second item ensures that \mathbf{P} is a probability distribution. The value $\iota_{init}(s)$ specifies the likelihood that the system evolution starts in s . All states s with $\iota_{init}(s) > 0$ are considered *initial states*. States s' with $\mathbf{P}(s, s') > 0$ are viewed as possible successors of the state s . The operational behavior is as follows. A initial state s_0 with $\iota_{init}(s_0) > 0$ is yielded. Afterwards in each state a transition is yielded at random according to the probability distribution \mathbf{P} in that state. The evolution of a Markov chain ends in a state s if and only if $\mathbf{P}(s, s) = 1$

- has no actions
 ”As compositional approaches for Markov models are outside the scope of this monograph, actions are irrelevant in this chapter and are therefore omitted.”

NOTES BEGIN

A Markov chain permits both probabilistic and nondeterministic choices thereby is a model that somewhat merges the concept of transition systems with the concept of Markov chains. They add probabilistic choices to transition systems or add nondeterminism to Markov chains. Probabilistic choices may be used to model the (probabilistic) behavior of the environment. But to do so statistical experiments are required to obtain adequate distributions that model the average behavior of the environment. If this information is not available too hard, too obtain, or a guarantee about system properties is required nondeterminism is the natural option for

modeling in these cases. Further usecases of MDPs are randomized distributed algorithms and abstraction in Markov chains. Randomized distributed algorithms are nondeterministic because there is a nondeterministic choice which process performs the next step and randomized because they have rather restricted set of actions that have a random nature. Abstraction in Markov chains can be feasible if states have been grouped by AP and have a wide range of transition probabilities. If that is the case selection of a transition is rather non deterministic and thereby can be abstracted with an MDP by replacing the probabilities with nondeterminism.

Definition 2.3. A *Markov decision process* is a tuple $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ where

- S is a countable set of states,
- Act is a set of actions,
- $\mathbf{P} : S \times Act \times S \rightarrow [0, 1]$ is the transition probability function such that for all states $s \in S$ and actions $\alpha \in Act$:

$$\sum_{s' \in S} \mathbf{P}(s, \alpha, s') \in \{0, 1\},$$

- $\iota_{init} : S \rightarrow [0, 1]$ is the initial distribution such that $\sum_{s \in S} \iota_{init}(s) = 1$,
- AP is a set of atomic propositions and
- $L : S \rightarrow \mathcal{P}(AP)$ a labeling function.

An action α is *enabled* in state s if and only if $\sum_{s' \in S} \mathbf{P}(s, \alpha, s') = 1$. Let $Act(s)$ denote the set of enabled actions in s . For any state $s \in S$, it is required that $Act(s) \neq \emptyset$. Each state s' for which $\mathbf{P}(s, \alpha, s') > 0$ is called an α -*successor* of s .

An MDP is called *finite* if S , Act and AP are finite. The transition probabilities $\mathbf{P}(s, \alpha, t)$ can be arbitrary real numbers in $[0, 1]$ (that sum up to 1 or 0 for fixed s and α). For algorithmic purposes they are assumed to be rational. The unique initial distribution ι_{init} could be generalized to set of ι_{init} with nondeterministic choice at the beginning. For sake of simplicity there is just one single distribution. The operational behavior is as follows. A starting state s_0 yielded by ι_{init} with $\iota_{init}(s_0) > 0$. From there on nondeterministic choice of enabled action takes place followed by a probabilistic choice of a state. The action fixed in the step of nondeterministic selection. Any Markov chain is an MDP in which for every state s , $Act(s)$ is a singleton set. Conversely an MDP with the property of $Act(s) = 1$ is a Markov chain. Thus Markov chains are a proper subset of Markov decision processes.

For convenience for $s_1, s_2 \in S$ and $\alpha \in Act$ we will write $(s_1, \alpha, s_2) \in \mathbf{P}$ if and only if $\mathbf{P}(s_1, \alpha, s_2) > 0$, that is to say if there is a non-zero probability of evolving from state s_1 to state s_2 with action α . Analogously we will write $s \in \iota_{init}$ if and only if is an initial state ($\iota_{init}(s) > 0$).

not sure if I should: $I := \iota_{init}$ thereby meaning the underlying set
CONVENTION END TEST

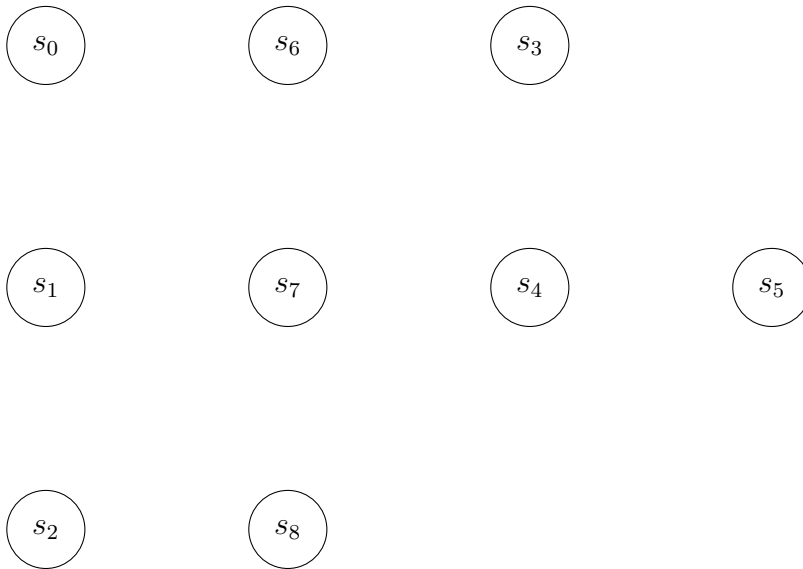


Figure 1: Simplified representation of \mathcal{M} (left) and the view \mathcal{M}_I on it(left)

3 View

Views are the central objective of this thesis. The purpose of a view is to obtain a simplification of a given MDP. It is an independent MDP derived from a given MDP and represents a (simplified) view on the given one - hence the name. Thereby the original MDP is retained.

In the preliminaries transition systems and Markov Chains were listed as simpler version of MDPs. Roughly speaking transition systems and MDPs are special MDPs namely that have no probability distribution in each state for each action or no actions. When defining views it seems feasible to do so for the most general system of the ones of consideration. That is why we will define views on MDPs. Only for specific views and their implementation it is to be kept in mind that if they regard an action or the probability distribution of an action in a state it is not applicable to transition systems or MCs respectively. **topic init states**

3.1 Grouping Function

The conceptional idea of a view is to group states by some criteria and structure the rest of the system accordingly. To formalize the grouping we define a dedicated function.

Definition 3.1. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be MDP, \tilde{S} a subset of S and M be an arbitrary set **with** $\bullet \in M$. We call any function $F_\theta : S \rightarrow M$ with $\tilde{F}_\theta(s) = \bullet$ if $s \notin \tilde{S}$.

$$s \mapsto \begin{cases} \tilde{F}_\theta(s), & \text{if } s \in \tilde{S} \\ \bullet, & \text{otherwise} \end{cases}$$

where $\tilde{F}_\theta : S \rightarrow M$.

a *grouping function*. The symbol θ is a unique identifier.

The identifier θ normally hints the objective of the grouping function. Two states are **grouped (should be Definition?)** to a new state if and only if the grouping function maps them to the same value **if that value is not the unique symbol \bullet** . The mapping to the symbol \bullet will be used whenever a state is supposed to be excluded from the grouping. The definition offers an easy way of defining groups of states. It is also very close to the actual implementation later on. The exact mapping depends on the desired grouping. In order to define a new set of states for the view, we define an equivalence relation R based on a given grouping function F_θ .

Definition 3.2. Let F_θ be a grouping function. We define the equivalence relation $R := \{(s_1, s_2) \in S \times S \mid F_\theta(s_1) = F_\theta(s_2) \neq \perp\} \cup \{(s, s) \mid s \in S\}$

R is an equivalence relation because it is reflexive, transitive and symmetric. R is reflexive because $\{(s, s) \mid s \in S\} \subseteq R$. Thus for all states s it is $(s, s) \in R$. Therefore for the properties transitivity and symmetry we only consider distinct $s_1, s_2 \in S$. If $F_\theta(\tilde{s}) = \bullet$ with \tilde{s} being one of states s_1 or s_2 it is $(s_1, s_2) \notin R$. Hence when considering transitivity and symmetry we assume $F_\theta(s_1), F_\theta(s_2) \neq \bullet$. If $F_\theta(s_1) =$

$F_\theta(s_2)$ it is obviously also $F_\theta(s_2) = F_\theta(s_1)$, which means if $(s_1, s_2) \in R$ it follows that $(s_2, s_1) \in R$. Hence R is symmetric. The property directly conveyed from equality. In the same way transitivity directly conveys from the equality relation to R .

We observe that two states s_1, s_2 are grouped to a new state if and only if $(s_1, s_2) \in R$. This is the case if and only if $s_1, s_2 \in [s_1]_R = [s_2]_R$ where $[s_i]_R$ for $i \in \mathbb{N}$ denotes the respective equivalence class of R , i.e. $[\tilde{s}]_R := \{s \in S \mid (s, \tilde{s}) \in R\}$. $[S]_R$ denotes the set of all equivalence classes of R , i.e. $[S]_R := \bigcup_{s \in S} \{[s]_R\}$.

3.2 Formal Definition

The definition of a view is dependent on a given MDP and a grouping function F_θ . We derive the equivalence relation R as in Definition 3.2 and use its equivalence classes $[s]_R$ ($s \in S$) as states for the view. The rest of the MDP is structured accordingly.

Definition 3.3. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be MDP and F_θ a grouping function. A *view* is MDP $\mathcal{M}_\theta = (S', Act', \mathbf{P}', \iota_{init}', AP', L')$ that is derived from \mathcal{M} with the grouping function F_θ where

- $S' = \{[s]_R \mid s \in S\} = [S]_R$
- $Act' = Act$
- $\mathbf{P} : [S]_R \times Act \times [S]_R \rightarrow [0, 1]$ with

$$\mathbf{P}'([s_1]_R, \alpha, [s_2]_R) = \frac{1}{|[s_1]_R|} \sum_{\substack{s_a \in [s_1]_R, \\ s_b \in [s_2]_R}} \mathbf{P}(s_a, \alpha, s_b)$$

- $\iota_{init}' : [S]_R \rightarrow [0, 1]$ with

$$\iota_{init}'([s]_R) = \sum_{s \in R} \iota_{init}(s)$$

- $L' : S' \rightarrow \mathcal{P}(AP), [s]_R \mapsto \bigcup_{s \in [s]_R} \{L(s)\}$

and R is the equivalence relation according to Definition 3.2.

The identifier θ is inherited from F_θ to \mathcal{M}_θ . The identifier declares that F_θ is the grouping function of \mathcal{M}_θ and thereby also uniquely determines \mathcal{M}_θ . Note that the definition is in a most general form in the sense that if in a view a property accounts to one piece of some entity the whole entity receives the property i.e.

- $(s_1, \alpha, s_2) \in \mathbf{P} \Rightarrow ([s_1]_R, \alpha, [s_2]_R) \in \mathbf{P}'$
- $s \in \iota_{init} \Rightarrow [s]_R \in \iota_{init}'$
- $\forall s \in S : L(s) \in L'([s]_R)$

other versions feasible??

3.3 Composition of Views

In essence views are simplification generated from MDP. It seems rather obvious that the composition of views is a very practical feature. Therefore in this chapter we will introduce, formalize and discuss different notions of compositions. All variants will ensure that the effect caused by each partaking views also takes effect in the composed view. Moreover it is to be generated in a way that the effect of each individual view can be reverted from the composed one. There may be restrictions in the order of removal.

3.3.1 Parallel Composition

One of the most uncomplicated ideas is to group states that match in the function value of all given grouping functions. This idea is parallel in the sense that a set of grouping function is combined to a new grouping function in one single step.

Definition 3.4. Let $\mathcal{M}_{\theta_1}, \mathcal{M}_{\theta_2}, \dots, \mathcal{M}_{\theta_n}$ be views and $F_{\theta_1}, F_{\theta_2}, \dots, F_{\theta_n}$ be its grouping functions. A *parallel composed grouping function* is a grouping function $F_{\theta_1 || \theta_2 || \dots || \theta_n} : S \rightarrow M$ with

$$s \mapsto (F_{\theta_1}(s), F_{\theta_2}(s), \dots, F_{\theta_n}(s))$$

The respective parallel composed view is denoted with $\mathcal{M}_{\theta_1 || \theta_2 || \dots || \theta_n}$.

The operator $||$ used in the definition is inspired from the operator used in electric circuits, when the respective elements are connected in parallel.

If we want to speak about this grouping function in a general way, where it is only of importance that we refer to this type of composition and the given grouping functions are of no importance, we will denote a *parallel composition grouping function* with $F_{||}$.

The operator $||$ in $F_{||}$ is associative with respect to the grouped states. That is exactly the same states are grouped no matter where parenthesis are put.

Proposition 3.1. Let F_u, F_v, F_w be grouping functions and S a set of states. For all $s_1, s_2 \in S$ it holds that

$$F_{(u||v)||w}(s_1) = F_{(u||v)||w}(s_2) \iff F_{u||(v||w)}(s_1) = F_{u||(v||w)}(s_2)$$

Proof. Let F_u, F_v, F_w be grouping functions and S a set of states with $s \in S$. It is

$$\begin{aligned} F_{u||v}(s) &= (F_u(s), F_v(s)) & F_{v||w}(s) &= (F_v(s), F_w(s)) \\ F_{(u||v)||w}(s) &= ((F_u(s), F_v(s)), F_w(s)) & F_{u||(v||w)}(s) &= (F_u(s), (F_v(s), F_w(s))) \end{aligned}$$

Let $s_1, s_2 \in S$ be arbitrary. Moreover let

$$\begin{aligned} F_{(u||v)||w}(s_1) &=: ((a, b), c) \text{ and} & \text{Then it is} & F_{u||(v||w)}(s_1) = (a, (b, c)) \text{ and} \\ F_{(u||v)||w}(s_2) &=: ((x, y), z). & & F_{u||(v||w)}(s_2) = (x, (y, z)). \end{aligned}$$

Then it is

$$F_{(u||v)||w}(s_1) = F_{(u||v)||w}(s_2) \iff F_{u||(v||w)}(s_1) = F_{u||(v||w)}(s_2)$$

This is because

$$\begin{aligned}
& ((a, b), c) = ((x, y), z) \\
\iff & (a, b) = (x, y) \wedge c = z \\
\iff & a = x \wedge b = y \wedge c = z \\
\iff & a = x \wedge (b, c) = (y, z) \\
\iff & (a, (b, c)) = (x, (y, z))
\end{aligned}$$

□

The operator \parallel in F_{\parallel} is also commutative with respect of the states being grouped.

Proposition 3.2. *Let F_u and F_v be grouping functions and S a set of states. For all $s_1, s_2 \in S$ it holds that*

$$F_{u \parallel v}(s_1) = F_{u \parallel v}(s_2) \iff F_{v \parallel u}(s_1) = F_{v \parallel u}(s_2)$$

Proof. Let F_u and F_v be grouping functions, $s_1, s_2 \in S$ and

$$\begin{aligned}
F_{u \parallel v}(s_1) &= (F_u(s_1), F_v(s_1)) =: (a, b) \\
F_{u \parallel v}(s_2) &= (F_u(s_2), F_v(s_2)) =: (x, y)
\end{aligned}$$

Then it is

$$\begin{aligned}
F_{v \parallel u}(s_1) &= (F_v(s_1), F_u(s_1)) = (b, a) \\
F_{v \parallel u}(s_2) &= (F_v(s_2), F_u(s_2)) = (y, x)
\end{aligned}$$

It follows that

$$F_{u \parallel v}(s_1) = F_{u \parallel v}(s_2) \iff F_{v \parallel u}(s_1) = F_{v \parallel u}(s_2)$$

because $(a, b) = (x, y) \iff a = x \wedge b = y \iff (b, a) = (y, x)$.

□

Other variants missing if existent Subsetapplication missing

4 View Examples

In this chapter we will introduce and discuss some view examples created by the author. Their purpose is to understand the idea and concept of a view and get to know some views that might be useful in real world applications.

When considering views we only want into account those that utilize properties of MDPs or that do computations that are also feasible on normal graphs but are of explicit relevance MDPs.

4.1 Views Utilizing MDP Components

In this subsection we will introduce some views the are purely based on the components of an MDP. Their will neither be computations on the graph-structure of an MDP nor computations using the result vector.

4.1.1 Atomic Propositions

One of the least involved approaches to create a view is to base it on the atomic propositions that are assigned to each state by the labeling function. The notion is to group states that were assigned the same set of atomic propositions. **Why useful?, currently no "has AP" → maybe should be added**

Definition 4.1. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. The view \mathcal{M}_{AP} is defined by its grouping function F_{AP} grouping function with $F_{AP} : S \rightarrow M, s \mapsto L(s)$.

The grouping function is exactly the labeling function i.e. for all $s \in S$ it is $F_{AP}(s) := L(s)$. So it is $F_{AP}(s_1) = F_{AP}(s_2) \iff L(s_1) = L(s_2)$. According to Definition 3.2 for $\tilde{s} \in S$ it is $[\tilde{s}]_R = \{s \in S \mid L(s) = L(\tilde{s})\}$.

By this we obtain the view \mathcal{M}_{AP} for a given MDP \mathcal{M} where: $S' = \bigcup_{s \in S} \{[s]_R\} = \bigcup_{a \in AP} \{\{s \in S \mid L(s) = a\}\}$. All other components are constructed as in Definition 3.3.

Although this view might seem rather simple because essentially it only performs $F_{AP} := L$ it is the most powerful one. This is because every view presented in the following is reducible to this one. That is because a grouping function essentially asserts an atomic proposition to every state, namely the value in respect to the considered property of a given view. The reduction can be realized by replacing the labeling function with the grouping function of the resepective view. That is $L := F_\theta$ and $AP := F_\theta(S)$ for some grouping function F_θ . While this works it alters the underlying MDP.

4.1.2 Initial States

An a little more involved idea than directly using a given function is to utilize the set of initial states. We can group states that have a probability greater zero, that they are started from. In practice this might be useful to quickly find all initial states.

Definition 4.2. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $I := \{s \in S \mid s \in \iota_{init}\}$. The view \mathcal{M}_I is defined by its grouping function with $F_I : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } s \in I \\ \bullet, & \text{otherwise} \end{cases}$$

and $M := \{\top, \bullet\}$.

For $s_1, s_2 \in S$ it is $F_I(s_1) = F_I(s_2)$ if and only if $s_1, s_2 \in I$ or $s_1 = s_2$. According to Definition 3.2 it is

$$\begin{aligned} [s]_R &= \{s \in S \mid F_I(s) = \top\} && \text{for } s \in I \text{ and} \\ [s]_R &= \{s \in S \mid F_I(s) = \{s\}\} = \{s\} && \text{for } s \notin I. \end{aligned}$$

By this we obtain the view \mathcal{M}_I for a given an MDP \mathcal{M} where: $S' = \bigcup_{s \in S} \{[s]_R\} = \{s \in S \mid s \in I\} \cup \bigcup_{s \in S \setminus I} \{\{s\}\}$.

All other components are constructed as in Definition 3.3.

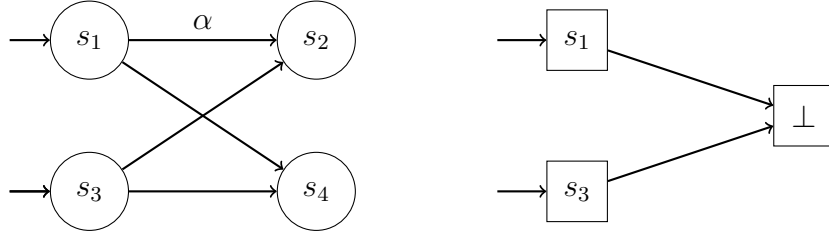


Figure 2: Simplified representation of \mathcal{M} (left) and the view \mathcal{M}_I on it (right)

4.1.3 Outgoing Actions

Another crucial component of an MDP is its set of actions Act . Actions are used for interprocess communication und synchronization. In this subsection we will provide and dicuss some views utilizing actions on transitions that are outgonig from a state. We will write a state s has an outgoing action α if there exists a state s' with $(s, \alpha, s') \in \mathbf{P}$. For a state s , ingoing action α we say it has an ingoing action α if there exist a state s' with $(s', \alpha, s) \in \mathbf{P}$.

The most obvious variant to group states is to group states that *have* a given outgoing action.

Definition 4.3. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\exists \alpha}$ is defined by its grouping function $F_{\exists \alpha} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s' \in S : (s, \alpha, s') \in \mathbf{P} \\ \bullet, & \text{otherwise} \end{cases}$$

and $M := \{\top, \bullet\}$.

For $s_1, s_2 \in S$ it is $F_{\exists\vec{\alpha}}(s_1) = F_{\exists\vec{\alpha}}(s_2)$ if and only if there exist $s_a, s_b \in S$ with $(s_1, \alpha, s_a), (s_2, \alpha, s_b) \in \mathbf{P}$ (i.e. they have the same outgoing action α) or $s_1 = s_2$. In accordance with Definition 3.2 it is

$$\begin{aligned} [s]_R &= \{s \in S \mid F_{\exists\vec{\alpha}}(s) = \alpha\} & \exists s' \in S : (s, \alpha, s') \in \mathbf{P} \\ [s]_R &= \{s \in S \mid F_{\exists\vec{\alpha}}(s) = s\} = \{s\} & \text{otherwise} \end{aligned}$$

Thereby we obtain the view $\mathcal{M}_{\exists\vec{\alpha}}$ for a given MDP \mathcal{M} where $S' = \bigcup_{s \in S} \{[s]_R\} =: S_1 \cup S_2$ where

$$\begin{aligned} S_1 &:= \{s \in S \mid \exists s' \in S : (s, \alpha, s') \in \mathbf{P}\} = \{s \in S \mid s \text{ has outgoing action } \alpha\} \\ S_2 &:= \bigcup_{s \in S \setminus S_1} \{\{s\}\}. \end{aligned}$$

Since actions are a very important part of MDPs as well as of its more powerful siblings MDPs and MCs it seems useful to further enhance this view and look at variants of it. Instead of only grouping states that only *have* outgoing actions we could also quantify the amount of times that action should be outgoing.

For example we could require that a given action has to be outgoing a minimum amount of times.

Definition 4.4. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{n \leq \vec{\alpha}}$ is defined by its grouping function $F_{n \leq \vec{\alpha}} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s_1, \dots, s_n \in S : Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$, $n \in \mathbb{N}$ is the minimum amount of times a transition with action α has to be outgoing in order to be grouped with the other states and

$$Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n) := ((s, \alpha, s_1), \dots, (s, \alpha, s_n) \in \mathbf{P}) \wedge |\{s_1, \dots, s_n\}| = n$$

is a first order logic predicate.

The number and The predicate $Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n)$ requires that there are transitions with action α to n distinct states.

For $s_1, s_2 \in S$ it is $F_{n \leq \vec{\alpha}}(s_1) = F_{n \leq \vec{\alpha}}(s_2)$ if and only if there exist distinct $s_{a_1}, \dots, s_{a_n} \in S$ and distinct $s_{b_1}, \dots, s_{b_n} \in S$ so that $(s_1, \alpha, s_{a_1}), \dots, (s_1, \alpha, s_{a_n}) \in \mathbf{P}$ and $(s_2, \alpha, s_{b_1}), \dots, (s_2, \alpha, s_{b_n}) \in \mathbf{P}$ or $s_1 = s_2$. According to Definition 3.2 it is

$$\begin{aligned} [s]_R &= \{s \in S \mid F_{n \leq \vec{\alpha}}(s) = \alpha\} & \text{if } Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n) \\ [s]_R &= \{s \in S \mid F_{n \leq \vec{\alpha}}(s) = s\} = \{s\} & \text{otherwise} \end{aligned}$$

By this we obtain the view $\mathcal{M}_{n \leq \vec{\alpha}}$ for a given MDP \mathcal{M} where $S' = \bigcup_{s \in S} \{[s]_R\} =: S_1 \cup S_2$ where

$$\begin{aligned} S_1 &:= \{s \in S \mid \exists s_1, \dots, s_n \in S : Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n)\} \\ &= \{s \in S \mid \text{the action } \alpha \text{ is outgoing at least } n \text{ times}\} \text{ and} \\ S_2 &:= \bigcup_{s \in S \setminus S_1} \{\{s\}\}. \end{aligned}$$

In a similar fashion we define view that groups states where at most a certain number of times a given action is outgoing.

Definition 4.5. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\vec{\alpha} \leq n}$ is defined by its grouping function $F_{\vec{\alpha} \leq n} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \forall s_1, \dots, s_{n+1} \in S : Q_{\vec{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$, is the maximal number of times a transition with action α may be outgoing and

$$Q_{\vec{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) := ((s, \alpha, s_1), \dots, (s, \alpha, s_{n+1}) \in \mathbf{P}) \implies \bigvee_{\substack{i,j \in \{1, \dots, n+1\} \\ i < j}} s_i = s_j$$

is a first order logic predicate.

It ensures that if there are one more than n outgoing transitions with an action α at least two of the states where the transitions **red** are in fact the same. Since this is required for all possible combinations of $n + 1$ states by the grouping function, only states that have at most n outgoing actions will be assigned with α by the grouping function. The reasoning about the equality of the grouping function values, the obtained equivalence classes and the resulting set of states S' of the view is analogous to $\mathcal{M}_{n \leq \vec{\alpha}}$.

Since we already defined grouping functions and hence views for a required minimal and maximal amount of times an action has to be outgoing it is now easily possible to define a view that groups states where the amount of outgoing actions is at least n and at most m .

Definition 4.6. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{m \leq \vec{\alpha} \leq n}$ is defined by its grouping function $F_{m \leq \vec{\alpha} \leq n} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s_1, \dots, s_m \in S : Q_{m \leq \vec{\alpha}}(s, s_1, \dots, s_m) \\ & \text{and } \forall s_1, \dots, s_{n+1} \in S : Q_{\vec{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$ and $m, n \in \mathbb{N}$ are the minimal and maximal number of transitions with action α in order for state to be grouped. The predicates $Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n)$ and $Q_{\vec{\alpha} \leq n}(s, s_1, \dots, s_{n+1})$ are the predicates from Definition 4.4 and Definition 4.5 respectively.

We already know that for a given $s \in S$ the expressions $\exists s_1, \dots, s_n \in S : Q_{n \leq \vec{\alpha}}(s, s_1, \dots, s_n)$ and $\forall s_1, \dots, s_{m+1} \in S : Q_{\vec{\alpha} \leq m}(s, s_1, \dots, s_{m+1})$ from Definition 4.4 and Definition 4.5 require that s has minimal and maximal amount of outgoing transitions with an action α respectively. Hence the conjunction will be true for states where the amount of outgoing transitions with action α is element of the set $\{m, n + 1, \dots, n - 1, n\}$. We will write for this that the number of outgoing actions is *in the span*.

For a given state s and action α we set

$$C_{s,\vec{\alpha}} := \exists s_1, \dots, s_m \in S : Q_{m \leq \vec{\alpha}}(s, s_1, \dots, s_m) \wedge \forall s_1, \dots, s_{n+1} \in S : Q_{\vec{\alpha} \leq n}(s, s_1, \dots, s_{n+1})$$

for convenience. $C_{s,\vec{\alpha}}$ is true if and only if the number of outgoing actions is in the span. For $s_1, s_2 \in S$ it is $F_{m \leq \vec{\alpha} \leq n}(s_1) = F_{m \leq \vec{\alpha} \leq n}(s_2)$ if and only if $C_{s_1,\vec{\alpha}} \wedge C_{s_2,\vec{\alpha}}$ or $s_1 = s_2$. Then its equivalence classes are

$$\begin{aligned} [s]_R &= \{s \in S \mid F_{m \leq \vec{\alpha} \leq n}(s) = \alpha\} & C_{s,\vec{\alpha}} \text{ true} \\ [s]_R &= \{s \in S \mid F_{m \leq \vec{\alpha} \leq n}(s) = s\} = \{s\} & \text{otherwise} \end{aligned}$$

The new set of states S' of the view $\mathcal{M}_{m \leq \vec{\alpha} \leq n}$ is the union of the equivalence classes of equivalence relation R on the set of states S of the original MDP. Hence it is $S' = \bigcup_{s \in S} [s]_R =: S_1 \cup S_2$ where

$$\begin{aligned} S_1 &:= \{s \in S \mid F_{m \leq \vec{\alpha} \leq n}(s) = \alpha\} \\ &= \{s \in S \mid C_{s,\vec{\alpha}} \text{ true}\} \\ &= \{s \in S \mid \text{the action } \alpha \text{ is outgoing } m \text{ to } n \text{ times}\} \text{ and} \\ S_2 &:= \bigcup_{s \in S \setminus S_1} \{\{s\}\}. \end{aligned}$$

The views above can be combined with parallel composition. The thereby obtained view requires that the respective conditions of all the combined views are met. In this sense it is a conjunctive combination.

Instead of making requirements about states and group them based on whether they meet these requirements it also possible to group states that are very similar or even identical in regard to their outaction. We consider this idea with the *OutActionsIdent View* in two variants: strong and weak (identitiy). Firstly we will consider the variant of strong identity.

Definition 4.7. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\vec{Act}(s)=}$ is defined by its grouping function $F_{\vec{Act}(s)=} : S \rightarrow M$ with

$$s \mapsto \{(\alpha, n) \mid \alpha \in Act, n \text{ is the number of times that } \alpha \text{ is outgoing from } s\}$$

and $M := Act \times \mathbb{N}_0 \cup \{\perp\}$.

The grouping function asserts to each state a set of pairs. Note that a pair is contained into the set for each action $\alpha \in Act$. In case there is no outgoing transition from state s with an action α it is $(\alpha, 0) \in F_{\vec{Act}(s)=}$. For $s_1, s_2 \in S$ it is $F_{\vec{Act}(s)=}(s_1) = F_{\vec{Act}(s)=}(s_2)$ if and only if s_1 and s_2 are mapped to the same set of pairs. By Definition 3.2 the obtained equivalence classes of R are

$$[s]_R := \{s \in S \mid F_{\vec{Act}(s)=}(s) = \{(\alpha_1, n_1), \dots, (\alpha_l, n_l)\}, l = |Act|\}$$

According to Definition 3.3 the set $S' := \bigcup_{s \in S} [s]_R$ is the set of states of $\mathcal{M}_{\vec{Act}(s)=}$. All other components of $\mathcal{M}_{\vec{Act}(s)=}$ are as usual structured in accordance with the Definition 3.3. As mentioned earlier a weak variant of the OutActionsIdent view is also conceivable.

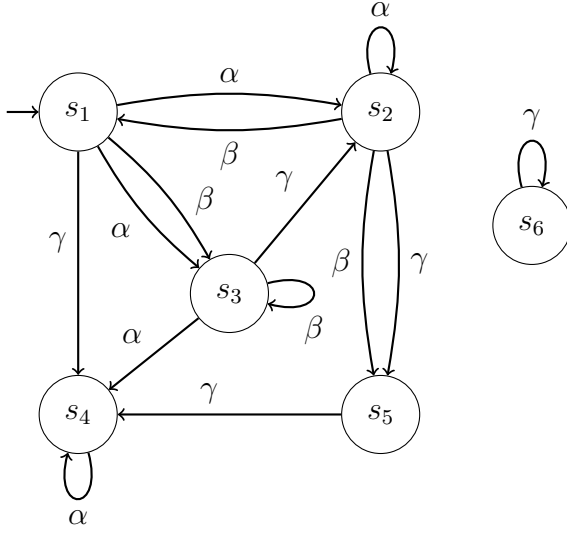


Figure 3: Simplified representation of \mathcal{M} (left) and the view \mathcal{M}_I on it (left)

Definition 4.8. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\vec{Act(s)} \approx}$ is defined by its grouping function $F_{\vec{Act(s)} \approx} : S \rightarrow M$ with

$$s \mapsto \{\alpha \in Act \mid \exists s' \in S : (s, \alpha, s') \in \mathbf{P}\}$$

and $M := Act \cup \{\perp\}$.

condition of image-set written in inconsistent style to strong identity. Swap strong and weak (order)?

The grouping function maps to the set of outgoing actions of a state and thereby discards information about the number of times an actions is outgoing. If an action is not outgoing from a state it is not contained in the set.

For $s_1, s_2 \in S$ it is $F_{\vec{Act(s)} \approx}(s_1) = F_{\vec{Act(s)} \approx}(s_2)$ if and only if they are mapped to the same set of actions. Hence the equivalence classes of R are

$$[\tilde{s}]_R = \{s \in S \mid F_{\vec{Act(s)} \approx}(s) = F_{\vec{Act(s)} \approx}(\tilde{s}) =: \{\alpha_1, \dots, \alpha_l\}, l \in \mathbb{N}\}.$$

According to Definition 3.3 the set $S' := \bigcup_{s \in S} [s]_R$ is the set of states of $\mathcal{M}_{\vec{Act(s)} \approx}$.

Apart from the option of directly considering one or a set of outgoing actions with possible quantities it is also possible to only consider the quantity of outgoing actions without regarding the any specific action.

Definition 4.9. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{|Act(s)|}$ is defined by its grouping function $F_{|Act(s)|} : S \rightarrow M$ with

$$s \mapsto |\{\alpha \in Act \mid \exists s' \in S : (s, \alpha, s') \in \mathbf{P}\}|$$

and $M := \mathbb{N} \cup \{\perp\}$.

The notion of the view is similar to utilize the outdegree, with the difference being here that outgoing transitions with the same action are considered as one single edge. This reflects on the options available for nondeterminism in this state.

The sepcialcase of there only being one single outgoing action is worth a distinct view, since it hides all nondeterministic choices, but nothing more.

Definition 4.10. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{|\overrightarrow{Act(s)}|_1}$ is defined by its grouping function $F_{|\overrightarrow{Act(s)}|_1} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } |\{\alpha \in Act \mid \exists s' \in S : (s, \alpha, s') \in \mathbf{P}\}| = 1 \\ \bullet, & \text{otherwise} \end{cases}$$

and $M := \mathbb{N} \cup \{\perp\}$.

4.1.4 Ingoing Actions

Analogously to Outgoing Actions views of utilizing ingoing actions are feasible. Since there is no difference apart from the definitions itself, we only provide the definitions.

Definition 4.11. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\exists \overleftarrow{\alpha}}$ is defined by its grouping function $F_{\exists \overleftarrow{\alpha}} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s' \in S : (s', \alpha, s) \in \mathbf{P} \\ \bullet, & \text{otherwise} \end{cases}$$

and $M := \{\top, \bullet\}$.

Definition 4.12. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{n \leq \overleftarrow{\alpha}}$ is defined by its grouping function $F_{n \leq \overleftarrow{\alpha}} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s_1, \dots, s_n \in S : Q_{n \leq \overleftarrow{\alpha}}(s, s_1, \dots, s_n) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$, is the minimum amount of times a transition with action α has to be ingoing in order to be grouped with the other states and

$$Q_{n \leq \overleftarrow{\alpha}}(s, s_1, \dots, s_n) := ((s_1, \alpha, s), \dots, (s_n, \alpha, s) \in \mathbf{P}) \wedge |\{s_1, \dots, s_n\}| = n$$

is a first order logic predicate.

Definition 4.13. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\overleftarrow{\alpha} \leq n}$ is defined by its grouping function $F_{\overleftarrow{\alpha} \leq n} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \forall s_1, \dots, s_{n+1} \in S : Q_{\overleftarrow{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$ is the maximal number of times a transition with action α may be ingoing and

$$Q_{\overleftarrow{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) := ((s_1, \alpha, s), \dots, (s_{n+1}, \alpha, s) \in \mathbf{P}) \implies \bigvee_{\substack{i, j \in \{1, \dots, n+1\} \\ i < j}} s_i = s_j$$

is a first order logic predicate.

Definition 4.14. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{m \leq \overleftarrow{\alpha} \leq n}$ is defined by its grouping function $F_{m \leq \overleftarrow{\alpha} \leq n} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } \exists s_1, \dots, s_m \in S : Q_{m \leq \overleftarrow{\alpha}}(s, s_1, \dots, s_m) \\ & \text{and } \forall s_1, \dots, s_{n+1} \in S : Q_{\overleftarrow{\alpha} \leq n}(s, s_1, \dots, s_{n+1}) \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$ and $m, n \in \mathbb{N}$ are the minimal and maximal number of transitions with action α in order for state to be grouped. The predicates $Q_{n \leq \overleftarrow{\alpha}}(s, s_1, \dots, s_n)$ and $Q_{\overleftarrow{\alpha} \leq n}(s, s_1, \dots, s_{n+1})$ are the predicates from Definition 4.11 and Definition 4.12 respectively.

Definition 4.15. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\overleftarrow{Act}(s)=}$ is defined by its grouping function $F_{\overleftarrow{Act}(s)=} : S \rightarrow M$ with

$$s \mapsto \{(\alpha, n) \mid \alpha \in Act, n \text{ is the number of times that } \alpha \text{ is ingoing from } s\}$$

and $M := (Act \times \mathbb{N}_0) \cup \{\perp\}$.

Definition 4.16. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{\overleftarrow{Act}(s) \approx}$ is defined by its grouping function $F_{\overleftarrow{Act}(s) \approx} : S \rightarrow M$ with

$$s \mapsto \{\alpha \in Act \mid \exists s' \in S : (s', \alpha, s) \in \mathbf{P}\}$$

and $M := Act \cup \{\perp\}$.

Definition 4.17. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\alpha \in Act$. The view $\mathcal{M}_{|\overleftarrow{Act}(s)|}$ is defined by its grouping function $F_{|\overleftarrow{Act}(s)|} : S \rightarrow M$ with

$$s \mapsto |\{\alpha \in Act \mid \exists s' \in S : (s', \alpha, s) \in \mathbf{P}\}|$$

and $M := \mathbb{N} \cup \{\perp\}$.

4.1.5 Variables

The concept of variables is not part of the definitions of neither TS, MCs or MDPs even though, it is of great importance in practice. We will introduce and discuss this concept before utilizing it for views.

Since states describe some information about a system at a certain moment of its behavior the information carried is usually not atomic but rather consists of several pieces. For instance when considering a computer program at a given moment during execution all its currently available variables will have some value, the stack will have a certain structure and the program counter points to a specific instruction. Systems in general at a certain moment of their behavior have several properties that in total pose the current state of the system. That is in practice each state is actually derived from a possible variable assertion. Choosing the respective variable assertion as state representation would result in rather complex state objects. Therefore the values of the variables are usually stored in a separate data structure and a simple identifier like an integer represents the actual state. When formalizing this in practice used

approach several options com into consideration. Available MDP components like atomic propositions or the set of states could be used to contain this information. There could be a subset of atomic propositions that declares that a certain variable has a certain value. There had to be taken care that for each variable in each state there is only one atomic proposition declaring its value. The set of states could be used in the sense that the states itself are complex objects containing the information. Both of these options are rather tedious but possible. A third option is to define a set of variables and an evaluation function that are induced by the MDP.

Definition 4.18. Let \mathcal{M} be an MDP. The set $Var_{\mathcal{M}}$ is called *variables* (of \mathcal{M}). It contains all variables induced by \mathcal{M} .

Definition 4.19. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and M be an arbitrary set. The by im induced function $VarEval_{\mathcal{M}} : S \times Var_{\mathcal{M}} \rightarrow M$ is called *variable evaluation function*.

Most of the time we will use $VarEval_{\mathcal{M}}$ to refer to the variable evaluation function. When we speak about the value of a variable in a state we refer to the image of $VarEval_{\mathcal{M}}$ for that state and variable. The set M is arbitrary so that arbitrary values can be assigned to a variable. Speaking in terms of computer science and programming this loosens as an example the restriction of only being able to assign numbers and no booleans.

The most apparent idea for a view utilizing variables is to group states that meet some requirement regarding the values of the variables.

state has variable not here uptil now because probably not used

Definition 4.20. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP, $x \in Var_{\mathcal{M}}$ and $a \in VarEval_{\mathcal{M}}(S, Var_{\mathcal{M}})$. The view $\mathcal{M}_{x=a}$ is defined by its grouping function $F_{x=a} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } VarEval_{\mathcal{M}}(s, x) = a \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$.

The view $\mathcal{M}_{x=a}$ groups states that share the same value for a given variable. For s_1, s_2 it is $F_{x=a}(s_1) = F_{x=a}(s_2)$ if and only if $VarEval_{\mathcal{M}}(s_1, x) = VarEval_{\mathcal{M}}(s_2, x)$ or $s_1 = s_2$. The obtained equivalence classes are

$$\begin{aligned} [s]_R &= \{s \in S \mid VarEval_{\mathcal{M}}(s, x) = a\} \\ [s]_R &= \{s \in S \mid F_{x=a}(s) = s\} = \{s\} \end{aligned}$$

The set of states S' of $\mathcal{M}_{x=a}$ is the union of the equivalence classes of R . It is $S' = \bigcup_{s \in S} [s]_R =: S_1 \cup S_2$ where

$$\begin{aligned} S_1 &:= \{s \in S \mid F_{x=a}(s) = a\} \\ &= \{s \in S \mid VarEval_{\mathcal{M}}(s, x) = a\} \text{ and} \\ S_2 &:= \bigcup_{s \in S \setminus S_1} \{\{s\}\}. \end{aligned}$$

Analogously a view that requires inequality instead of equality is feasible.

Definition 4.21. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP, $x \in Var_{\mathcal{M}}$ and $a \in VarEval_{\mathcal{M}}(S, Var_{\mathcal{M}})$. The view $\mathcal{M}_{x \neq a}$ is defined by its grouping function $F_{x \neq a} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } VarEval_{\mathcal{M}}(s, x) \neq a \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$.

If states are to be grouped with the requirement of several variables equaling or not equaling specified values this can be achieved by using parallel composition.

To allow even more flexibility a view can be used that also allows a combination of requirements on variables in a disjunctive manner. To extend this idea to its full potential we will define a view that allows requirements using a disjunctive normal formal (DNF). To formalize this view more efficiently we will write $x_{s,i}$ short for $VarEval_{\mathcal{M}}(s, x_i)$ where $x_i \in Var_{\mathcal{M}}$ and $i \in \mathbb{N}$. We define the symbol \doteq to be an element of the set $\{=, \neq\}$. That is to say, whenever it is used each time written it is a representative for either the symbol $=$ or \neq . It allows to write one symbol whenever $=$ and \neq could or should be possible. Moreover for this context we consider $(x_{s,i} = a)$ as a literal and $(x_{s,i} \neq a)$ as its negation. We write $(x_{s,i} \doteq a)$ for a literal that could be negated or not negated.

Definition 4.22. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP, $x \in Var_{\mathcal{M}}$, $a \in VarEval_{\mathcal{M}}(S, Var_{\mathcal{M}})$ and

$$d(s) = ((x_{s,11} \doteq a_{11}) \wedge \dots \wedge (x_{s,1l_1} \doteq a_{1l_1})) \vee \dots \vee ((x_{s,k1} \doteq a_{k1}) \wedge \dots \wedge (x_{s,kl_k} \doteq a_{kl_k}))$$

proposition logical formulae in disjunctive normal form where $x_{s,ij}, a_{ij} \in VarEval_{\mathcal{M}}(S, Var_{\mathcal{M}})$. The view \mathcal{M}_{VarDNF} is defined by its grouping function $F_{VarDNF} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } d(s) \text{ is true} \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$.

The DNF allows us to specify a requirement in disjunctive normal form about variables. States are mapped to the same value depending on whether or not they meet this requirement.

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Analogously a view based on a conjunctive normal formal can be defined that may be more convenient, depending on the query.

Definition 4.23. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and

$$\begin{aligned} c(s) = & ((x_{s,i_1} \doteq a_{i_1}) \vee \dots \vee (x_{s,i_{l_1}} \doteq a_{i_{l_1}})) \wedge \\ & ((x_{s,i_{l_1}+1} \doteq a_{i_{l_1}+1}) \vee \dots \vee (x_{s,i_{l_2}} \doteq a_{i_{l_2}})) \wedge \\ & \dots \\ & ((x_{s,i_{(l_{m-1})+1}} \doteq a_{i_{(l_{m-1})+1}}) \vee \dots \vee (x_{s,i_{l_m}} \doteq a_{i_{l_m}})) \end{aligned}$$

proposition logical formula in disjunctive normal form where

- $x_{s,i_k}, a_{i_k} \in \text{VarEval}_{\mathcal{M}}(S, \text{Var}_{\mathcal{M}})$ with $i_k \in \mathbb{N}$ and $k \in \{1, \dots, l_m\}$
- $l_1 < l_2 < \dots < l_m$ and m is the number of clauses in $c(s)$

The view $\mathcal{M}_{\text{VarCNF}}$ is defined by its grouping function $F_{\text{VarCNF}} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } c(s) \text{ is true} \\ \bullet, & \text{otherwise} \end{cases}$$

where $M := \{\top, \bullet\}$.

The only difference from the view $\mathcal{M}_{\text{VarCNF}}$ to the view $\mathcal{M}_{\text{VarDNF}}$ is whether the respective formulae is in conjunctive or disjunctive normal form. **CITATION** Since each formulae in conjunctive normal form can be transformed to a formulae in disjunctive normal form and vice versa neither one of the views can perform an action the other can not. Hence there is no difference in expressivity, but there may be one in size. Therefore both views have been implemented and formalized.

The views discussed before reduce the MDP in a very precise but also manual manner, because it not only dictates the variable but also its value. A more general approach is to stipulate only the variable but not its value. This way states will be grouped that have the same value for that variable with no regard to the actual value of that variable. This idea could be achieved with a view based on the grouping function F_{a_1}

Definition 4.24. Let $\mathcal{M} = (S, \text{Act}, \mathbf{P}, \iota_{\text{init}}, AP, L)$ be an MDP. The view $\mathcal{M}_{\text{VarEval}_{\mathcal{M}}(s,x)}$ is defined by its grouping function $F_{\text{VarEval}_{\mathcal{M}}(s,x)} : S \rightarrow M$ with

$$s \mapsto \text{VarEval}_{\mathcal{M}}(s, x)$$

and $M := \text{VarEval}_{\mathcal{M}}(S, x)$.

With this grouping function we directly map to the value of the variable. Hence for $s_1, s_2 \in S$ it is $F_{\text{VarEval}_{\mathcal{M}}(s,x)}(s_1) = F_{\text{VarEval}_{\mathcal{M}}(s,x)}(s_2)$ if and only if they are mapped to the value $a \in M$. Hence the equivalence classes of R are

$$[\tilde{s}]_R = \{s \in S \mid \text{VarEval}_{\mathcal{M}}(s) = \text{VarEval}_{\mathcal{M}}(\tilde{s})\}.$$

According to Definition 3.3 the set $S' := \bigcup_{s \in S} [s]_R$ is the set of states of $\mathcal{M}_{\text{VarEval}_{\mathcal{M}}(s,x)}$.

4.2 Utilizing the MDP Graphstructure

4.2.1 Distance

Considering distances in a graph can be very helpful to get an overview of a graph. Likewise it helps a lot with understanding the structure of an MDP. In order to consider the distance between nodes we will need to formalize it.

definition is very similar to execution fragments! Citation?

Definition 4.25. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. A (simple and finite) *path* P is a sequence $(s_1, \alpha_1, s_2, \alpha_2, \dots, \alpha_{n-1}, s_n)$ alternating between states and actions where $n \in \mathbb{N}$, $\{s_1, \dots, s_n\} \subseteq S$ is a set of distinct states, $\{\alpha_1, \dots, \alpha_n\} \subseteq Act$ and for all $i \in \{1, \dots, n\}$ it is $(s_i, \alpha_i, s_{i+1}) \in \mathbf{P}$.

It is $first(P) := s_0$ and $last(P) := s_n$ and $P_{\mathcal{M}}$ the set of all paths in \mathcal{M} .

Definition 4.26. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $P = (s_1, \alpha_1, s_2, \alpha_2, \dots, \alpha_{n-1}, s_n)$ be a path in \mathcal{M} . The number $n =: len(P)$ is called *length* of the path P .

Definition 4.27. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. The *distance* from $\tilde{S} \subseteq S$ to $s \in S$ is the length of the shortest path from a state $\tilde{s} \in \tilde{S}$ to s . That is $dist(\mathcal{M}, \tilde{S}, s) := \min\{len(P) \mid P \in P_{\mathcal{M}}, first(P) \in \tilde{S}, last(P) \in s\}$.

For a view that applies to the whole and utilizes distance it only makes sense to consider a given set of states from which one the distance is measured. An intuitive choice for such set is the set of initial states. The following algorithm calculates the distance of each node from a given set $\tilde{S} \subseteq S$ considering granularity n :

Implementation Algorithm

The set $distance(\mathcal{M}, \tilde{S}, n)$ declares the returned set of **ALGORITHM**. The implementation ensures that for every state s there exists a pair (s, d) in $distance(\mathcal{M}, \tilde{S}, n)$. The following view groups states that have the same distance to the set measured with the amounts of transitions necessary to reach the the closest \tilde{S} considering the granularity n .

Definition 4.28. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $\tilde{S} \subseteq S$ arbitrary. **$distance(\mathcal{M}, \tilde{S}, n)$ IMPLEMENTATION IN PSEUDOCODE? DEFINITION BEFORE?** The view $\mathcal{M}_{\leftrightarrow}$ is defined by its grouping function $F_{\leftrightarrow} : S \rightarrow M$ with

$$s \mapsto d \quad \text{where } (s, d) \in distance(\mathcal{M}, \tilde{S}, n)$$

and $M = \mathbb{N} \cup \{\inf\} \cup \{\perp\}$.

Variants of **ALGORITHM** which allow directinoless or only reverse traversal of the MDP are provided in the appendix. The respective views would utilize their returned set in the exact same way. That is, $distance(\mathcal{M}, \tilde{S}, n)$ is replaced with the set of the respective algorithm and nothing else changes.

4.2.2 Double Directed Edges with same Action

NOT SURE IF SHOULD BE INCLUDED

Definition 4.29. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP.

4.2.3 Cycles

A cycle is a structure that can exist in every graph. The concept of cycles is not specific to MDPs or any of its more specialized variants. The purpose of this thesis is to discuss views that utilize domain specific knowledge or if a general concept is

of special relevance when exploring an MDP. The former and the latter apply on cycles.

Formalizing views based on cycles requires some formalization of the concept cycle. We will use a domain specific definition that will serve us the most.

Definition 4.30. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. A (simple) *cycle* C in \mathcal{M} is a sequence $(s_0, \alpha_0, s_1, \alpha_1, \dots, \alpha_{n-1}, s_0)$ alternating between states and actions where $n \in \mathbb{N}$, $\{s_0, \dots, s_{n-1}\} \subseteq S$ is a set of distinct states, $\{\alpha_0, \dots, \alpha_{n-1}\} \subseteq Act$ and for all $i \in \{0, \dots, n-1\}$ it is $(s_i, \alpha_i, s_{i+1 \bmod n}) \in \mathbf{P}$.

When the actions in the cycle are of no further importance, we will omit them only writing a sequence of states. In the following let $C = (s_0, \alpha_0, s_1, \alpha_1, \dots, \alpha_{n-1}, s_0)$ be a cycle. For conveniences we will write $s \in C$ if the state is contained in the cycle C and $\alpha \in C$ if the action is contained in in the cycle C . **only if??** In words we will both write a state or action is *on* or *in* the cycle. Let C_1 and C_2 be cycles. $C_1 \cup C_2 := \{s \in S \mid s \in C_1 \text{ or } s \in C_2\}$.

Definition 4.31. Let C be an cycle in \mathcal{M} and S be the states set of \mathcal{M} . The number $|\{s \in S \mid s \in C\}| =: len(C)$ is called the *length* of a cycle.

Definition 4.32. Let \mathcal{M} be an MDP. The set $C_{\mathcal{M},n} := \{C \mid len(C) \geq n\}$ declares the set of all cycles in \mathcal{M} with a length of at least n .

In practice there exist several cycle finding algorithms. The function $findCycles(\mathcal{M}, n)$ is an abstraction for one of these algorithms being used. The actual implementation relies on algorithms of the java library jGraphT namely the **Algorithm Szwarcfiter and Lauer - $O(V + EC)$ and Tiernan - $O(V \cdot constV)$** CITATION!!

With the formalization done we will introduce some views utilizing the concept cycles. For one we will combine the notion cycle with domain specific knowledge for the other we will consider how and why cycles in MDPs are in general of relevance. We will begin with the latter. Cycles in general are of interest because they pose the risk of getting stuck in endless loops, when performing actions on the MDP. **Model checking is one of the most relevant actions on MDPs that are vulnerable to loops. They are performed commonly on them.** Therefore a view that simply finds existing cycles is feasible.

Definition 4.33. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $n \in \mathbb{N}$. The view $\mathcal{M}_{\exists C}$ is defined by its grouping function $F_{\exists C} : S \rightarrow M$ with

$$s \mapsto \begin{cases} \top, & \text{if } C_{\mathcal{M},n} \neq \emptyset \\ \bullet, & \text{otherwise} \end{cases}$$

and $M = \{\top, \bullet\}$.

This view groups all states that are contained in cycle with a length of at least n . It is to note that the affinity of a state to one or several respective cycles is lost.

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Definition 4.34. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $n \in \mathbb{N}$. The view $\mathcal{M}_{\{C_n\}}$ is defined by its grouping function $F_{\{C_n\}} : S \rightarrow M$ with

$$s \mapsto \{C \in C_{\mathcal{M},n} \mid s \in C\}$$

and $M = \mathcal{P}(C_{\mathcal{M},n}) \cup \{\perp\}$.

This view groups states that have the same set of cycles they are contained in. Thus if C_1 and C_2 are distinct cycles and $s_1, s_2 \in C_1$ and $s_1 \in C_2$ but $s_2 \notin C_2$ they are not grouped. It suffices that a state is on one cycle that the other one is not on, in order for the states not being grouped. In graphs with many cycles this can lead to little grouping.

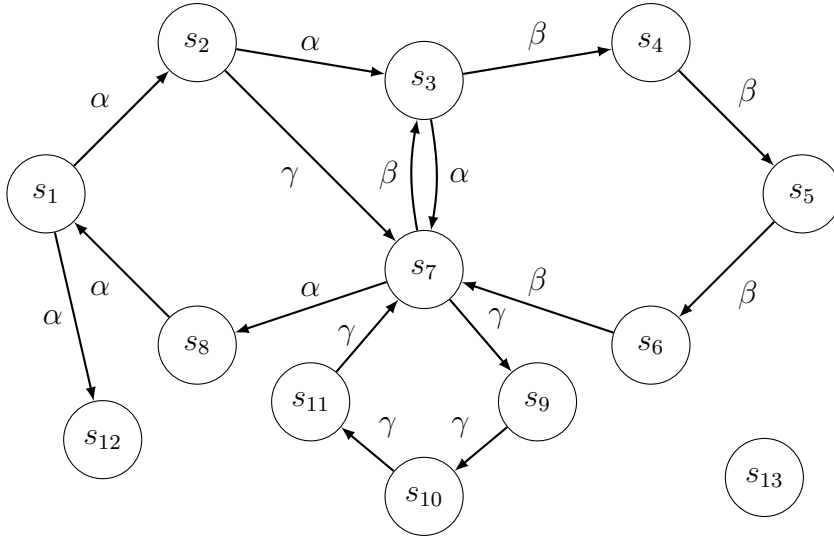


Figure 4: Simplified representation of \mathcal{M} (left) and the view \mathcal{M}_I on it (left)

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GREEDY APPROACHES (STATE ON SINGLE CYCLE), AND SET APPROACH $s \rightarrow C_1 \cup C_2$ omitted because not sure if needed

The view above might be useful when having found cycles to see what cycles specifically exist. Often it is interesting to find cycles that consist only of transitions of the same action. The following view accomplishes that.

Definition 4.35. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $n \in \mathbb{N}$. The view $\mathcal{M}_{\{C_{\alpha,n}\}}$ is defined by its grouping function $F_{\{C_{\alpha,n}\}} : S \rightarrow M$ with

$$s \mapsto \{C \in C_{\mathcal{M},n} \mid s \in C, \tilde{\alpha} \in C, \forall \alpha \in C : \alpha = \tilde{\alpha}\}$$

and $M = \mathcal{P}(C_{\mathcal{M},n}) \cup \{\perp\}$.

The view specializes the view from Definition 4.34 in the sense that it additionally requires for all $C \in F_{\{C_n\}}$ that all actions occurring in the cycle are the same. doubling to above definition?

NO DISCUSSION IF PREVIOUSLY

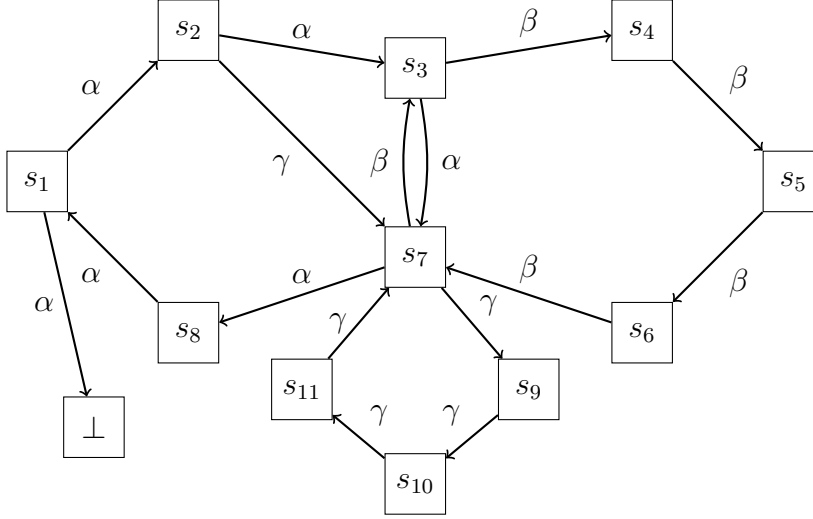


Figure 5: Simplified representation of \mathcal{M} (left) and the view \mathcal{M}_I on it (left)

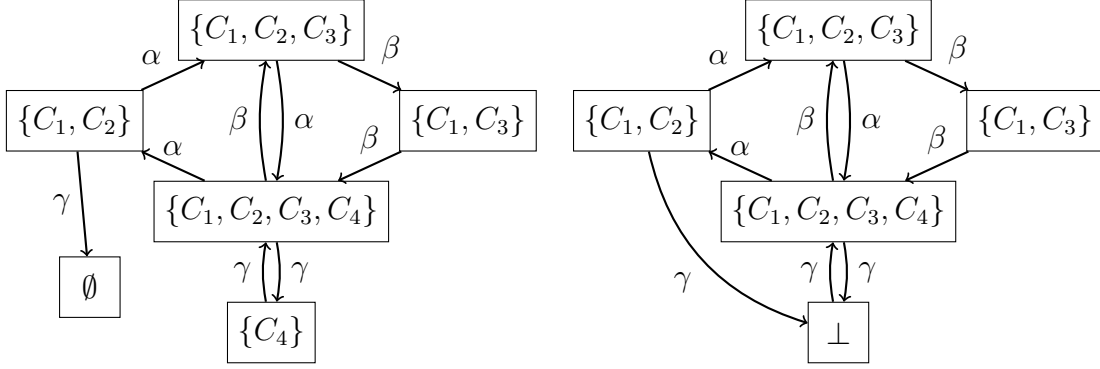


Figure 6: View $\mathcal{M}_{\{C_{\alpha,0}\}}$ (left) and view $\mathcal{M}_{\{C_{\alpha,5}\}}$ the right

In definition 4.34 we stated that little grouping can occur when mapping on the set of cycles in all of which the state is contained. Hence we provide the definition and implementation of a view that groups states even though their set of cycles is not equal, but there is sufficient similarity in the cycles they are on.

Definition 4.36. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP and $n \in \mathbb{N}$. The view $\mathcal{M}_{\cup\{s\}_C}$ is defined by its grouping function $F_{\cup\{s\}_C} : S \rightarrow M$ with

$$s \mapsto \{\tilde{s} \in S \mid s, \tilde{s} \in C \in C_{\mathcal{M},n}\}$$

and $M = S \cup \{\perp\}$.

Each state is mapped to the set of states resulting from joining the state sets of all the cycles the state is on.

4.2.4 Strongly Connected Components

Strongly connect components (SCC) are of major importance in model checking **I think so because I found the terms many times in the book. Remaining question: really? why?** Therefore it is feasible to consider a view utilizing strongly connected

components. Deviating from most definitions we will define SCC not as a subgraph but only as the set of its nodes. Moreover the definition is written in the terms of an MDP.

Definition 4.37. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. A set of the form

$$T = \{s \in S \mid \forall s' \in T : (\exists P \in P_{\mathcal{M}} \text{ with } first(P) = s, last(P) = s') \text{ or } s = s'\}$$

is called *strongly connected component* of \mathcal{M} . That is for every state in the set there has to exist a sequence of transitions to every other state of the set. The state itself is always considered reachable by itself, i.e. there does not need to exist a self loop. In consequence, each state is contained in strongly connected component namely at least in the one that only contains itself. The set of strongly connected components of \mathcal{M} is denoted with $SCC(\mathcal{M})$.

Since the strong connection is an equivalence relation the SCCs are equivalence classes and hence disjoint. To find SCCs Tarjans algorithm is the classic. In the implementation an improved variant from Gabow is used supplied by the jGraphT library. **CITATION**

Definition 4.38. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. The view \mathcal{M}_{scc} is defined by its grouping function $F_{scc} : S \rightarrow M$ with

$$s \mapsto T \quad \text{where } s \in T \in SCC(\mathcal{M})$$

and $M = SCC(\mathcal{M}) \cup \{\perp\}$.

This view groups all states together that are in the same SCC. Because SCCs are disjoint each s will be mapped to its one and only SCC. This is because... **DISCUSSION OF EQUALITY, EQ CLASSES AND RESULTING STATES MISSING**

A special kind of strongly connected components is the the bottom strongly connected component.

Definition 4.39. Let T be a SCC. A *bottom strongly connected component* (BSCC) is a SCC where it holds that: **concurrent versions**

$$\forall s \in T : \forall (s, \alpha, s') \in \mathbf{P} : s' \in T$$

$$\forall s \in T : \sum_{t \in T} \mathbf{P}(s, t) = 1$$

That is from T there is no state reachable outside of T . The set of bottom strongly connected components of \mathcal{M} is denoted with $BSCC(\mathcal{M})$.

These are of special relevance because..

Definition 4.40. Let $\mathcal{M} = (S, Act, \mathbf{P}, \iota_{init}, AP, L)$ be an MDP. The view \mathcal{M}_{bscc} is defined by its grouping function $F_{bscc} : S \rightarrow M$ with

$$s \mapsto T \quad \text{where } s \in T \in BSCC(\mathcal{M})$$

and $M = BSCC(\mathcal{M}) \cup \{\perp\}$.

In the implementation strongly connected components are determined using the algorithm of Gabow, afterwards filtering those SCCs that have only transitions to states within the SCC. Equality, equivalence classes and the new state set are constructed analogously to the view \mathcal{M}_{scc} .

4.3 Utilizing the MDP Result Table

In this section we will discuss views utilizing the result table. The actual implementation relies on one powerful view that can set to perform arbitrary actions using model checking results.

[needhelp Understanding Cluster](#)

5 View Implementation

supposed to give a short overview how view are implemented from a conceptional perspective. No more than 1 to 2 pages.

\hat{F}

- implementation in web application pmc-vis that aims for ...
- the general project structure is ... prism model, database, java, json + roughly were clusters are located in the project
- implementation of views rely on an internal graph structure that was necessary for the views that implement grouping based on structural properties of the MDP graph. Views based on MDP components firstly were implemented with direct accesses on the database but later on also adopted to the internal graph structure for performance reasons
- internal graph structure the jGraphT library was used. It supplies graph structures as well as common algorithms performed on them. It was chosen because it is the most common, most up to date java library for graphs with the best documentation and broadest functionality. It is developed by ... has a java style documentation and is open source. The broad functionality assured that when implementing a view it is most certainly assured that if necessary a respective algorithm is available.
- The MDP was realized as an directed weighted pseudograph. The transition probabilities are stored as weights. A pseudograph has been chosen because it allows double edges as well as loops.
- in order of keeping the graph lean nodes and vertices are long values. They refer to the state id and the transition id respectively. The state id matches the one in the db whereas the transition id is newly generated because in the database a transition is an action with its probability distribution. This difference to the MDP definition is reversed with the internal graph structure for predefined algorithms to work properly on the MDP graph.
- For access to atomic propositions, actions and alike the graph maintains two hashmaps that map the id to the actual transition or state object respectively
- depending on enum value according cluster is created
- normally each view is a dedicated java class, sometimes views are merged into one class that realizes combining behavior that otherwise would have been realized by composition (why? → was easy, allows more flexibility in functionality)
- when implementing a view only the grouping function is implemented. This makes sense because the process of actually generating the view is always the same as it was already notable in the definitions of a view, where every view is defined by its grouping function.

- the grouping function is calculated and its function values to the result table in a new column.
- visualization in general and also visualization after grouping was not part of the objective of the thesis and already present. As within the thesis there is one general approach how the resulting new graph is generated.
- every view class is registered in a public enum used to select which view is to be generated
- every view class is derived from general view class.
- in the implementation view still have the deprecated name cluster
- every view class has some private attributes to store information like
- every view class has the methods
- the most relevant one is build cluster. It consist of 4 parts. 1. Checking if it is already build aborting if so. 2.create a new column in the database where the results of the grouping function are to be saved. 3. The actual computing part of the grouping function, where for every state an SQL Query adding the respective grouping function value is added to a list of SQL Query strings. 4. all the SQL queries that contain the insertion operation of the grouping function value being executed.
- We will take a look at this in the example..

6 Comparison and Evaluation

Performance Cycles - Selection of states and induced subgraph easily possible perform cycle search only on subset of graph - in general generating views on subgraphs easily possible (only when needed for performance) Clustering exact Cycles when clustering exact cycles

7 Outlook

Many other views accomplishable with the properties view. ie Initcluster AP Cluster has Cycle All implemented Clusters in Appendix (only definition) $\mathbb{Z}_n[s]_{RS}$ \mathbb{N} in symbol list Implementation of own algorithms Distance Cluster Forward backward both \circ 870 End Components possible probabilities = 1