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A Partial Order Approach to Branching Time Logic Model Checking

by

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A Partial Order Approach to Branching Time Logic Model Checking

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Abstract

Partial order techniques enable reducing the size of the state graph used for model checking, thus alleviating the 'state space explosion' problem. These reductions are based on selecting a subset of the enabled operations from each program state. So far, these methods have been studied, implemented and demonstrated for assertional languages that model the executions of a program as computation sequences, in particular the logic LTL (linear temporal logic). The present paper shows, for the first time, how this approach can be applied to languages that model the behavior of a program as a tree. We study here partial order reductions for branching temporal logics, e.g., the logics CTL and CTL* (all logics with the next-time operator removed) and process algebras such as CCS. Conditions on the subset of successors from each node to guarantee reduction that preserves CTL properties are given. Provided experimental results show that the reduction is substantial.

1 Introduction

Partial order (or more accurately, commutativity-based) methods are useful for tackling the exponential blowup in the memory required for the automated verification by model-checking of concurrent programs. They exploit the fact that many properties are insensitive to the order in which concurrent actions are executed. Fixing one out of many such orders can then be used to reduce the memory and time needed to check such properties. Such methods were studied so far [5, 10, 20, 23, 24] in conjunction with specifications that assert about the set of interleaved executions of the program; e.g., that use linear temporal logic without the next-state operator (LTL-X).

State-based algorithms for model checking a system are patterned after a depth-first search of the system’s configurations or states, thus generating a state graph that allows checking whether a concurrent finite state program P satisfies a temporal logic property \( \varphi \). Partial order reductions are aimed at constructing a reduced state graph, based on exploring for each visited state only a subset of the enabled operations, so that only some of the successors of that state are expanded and, hence, specifications can be verified in less space and time. The correctness of the reduced state graph generation algorithm is based on employing a set of constraints that limit the choice of such subsets of operations to those that guarantee that the evaluation of specifications is preserved.

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The next step is to try to extend these methods to handle other types of specifications. Natural candidates are specification languages based on branching models, in particular, branching time temporal logics. Such logics, as opposed to LTL-X, can distinguish the state where a nondeterministic choice is made in the execution of the program. We are guided by three main reasons for our pursuit of a reduction that preserves branching-time logics. The first one is achieving greater expressiveness, e.g., by using a logic such as CTL*-X, which, besides being able to distinguish the nondeterministic choices, can express all LTL-X-properties. The second one is the existence of some interesting restricted versions of branching time logics such as CTL-X. Although CTL-X does not include LTL-X (and vice versa), it can, by virtue of the branching operators, describe many interesting properties of programs. Moreover, due to its restrictions, it has a model-checking algorithm that is linear in the size of the checked formula [2], as opposed to the exponential algorithm for LTL-X [14]. The third motivation for such a reduction lies within the fact that branching temporal properties are preserved by bisimulation [1]; besides basing our correctness proof on this fact, checking that two states are bisimulation equivalent is itself important for process-algebra style correctness. Thus, our reductions can be used to improve the time required and the size of the state graph and can be used in conjunction with process-algebra based tools such as PSF and AUTO [15, 19].

The paper starts out investigating the proper constraints on the subset that is chosen to be explored at each visited state. Not unexpectedly, the set of constraints turns out to be strictly stronger than the one needed for LTL-X. Indeed, CTL*-X is more expressive than LTL-X is, so that branching points due to nondeterministic choices should be preserved in the reduced graph. Of course, this also means that reduction for LTL-X can produce smaller state graphs, and thus be more efficient in space and time. This is compensated by the fact that some branching time logics such as CTL-X have model checking algorithms that are linear rather than exponential in the size of the checked property.

The proof of the correctness of our algorithm is novel in that it is rather different from the one used for LTL-X reductions [20]: instead of using traces [16], i.e., equivalence classes of sequences, we show stuttering equivalence between the full and the reduced state graph [1]. This equivalence was proved in [1] to be a necessary and sufficient condition for ensuring that the two stuttering equivalent structures satisfy the same CTL*-X formulas.

CTL*-X is the most expressive of the logics we discuss and, consequently, the same result holds for logics that are included in it, namely ACTL*-X, ACTL-X, and CTL-X.

Experimental results show that even with the additional constraint on selecting subsets of the enabled operations, the reduction is still substantial. We demonstrate the reduction on various algorithms and protocols and compare it to the reduction obtained for LTL-X. The simplicity of the reduction algorithm, and the small overhead in time and memory it incurs, suggests that one can obtain significant improvement for state-based model checking, by using the suggested reduction algorithm, with a relatively small investment. We also investigate using our algorithm as part of a branching bisimulation checker. Experiments indicate that it is more efficient to use our reduction strategy to generate a state graph to be checked than it is to generate and check the full state graph.

2 Basic Notions

Syntax of CTL*-X

Let $PV$ be a finite set of propositions. The set of state formulas and the set of path formulas are defined inductively:

S1. every member of $PV$ is a state formula.

S2. if $\phi$ and $\psi$ are state formulas, then so are $\neg \phi$, $\phi \land \psi$, and $\phi \rightarrow \psi$. 


S3. if \( \varphi \) is a path formula, then \( A \varphi \) is a state formula.

P1. any state formula \( \varphi \) is also a path formula.

P2. if \( \varphi, \psi \) are path formulas, then so are \( \varphi \land \psi \) and \( \neg \varphi \).

P3. if \( \varphi, \psi \) are path formulas, then so is \( \mathcal{U}(\varphi, \psi) \).

The modal operator \( A \) has the intuitive meaning: "for all paths". \( \mathcal{U} \) denotes the standard Until.

\( \text{CTL}^*-\text{X} \) consists of the set of all state formulae.

The following abbreviations will be used:

\[ E\varphi \overset{\text{def}}{=} \neg A \neg \varphi, \quad F\varphi \overset{\text{def}}{=} \mathcal{U}(\text{true}, \varphi), \quad G\varphi \overset{\text{def}}{=} \neg F \neg \varphi. \]

Sublogics of \( \text{CTL}^*-\text{X} \)

\( \text{CTL}\text{-X} \). The state modalities \( E \) and \( A \) and the path modalities \( U, F \) and \( G \) may only appear paired, i.e., in the combinations \( EU, EF, EG, AU, AF \) and \( AG \).

\( \text{ACTL}^*-\text{X} \). The modality \( E \) is prohibited, and negation can be applied only to subformulas that do not contain modalities.

\( \text{ACTL}\text{-X} \). The sublogic of \( \text{CTL}\text{-X} \) in which the modality \( E \) is prohibited, and negation can be applied only to subformulas that do not contain modalities.

\( \text{LTL}\text{-X} \). Restriction to formulas of the form \( A\varphi \), where \( \varphi \) does not contain \( A \) and \( E \). We usually write \( \varphi \) instead of \( A\varphi \) if confusion is unlikely.

Semantics of \( \text{CTL}^*-\text{X} \)

Let \( T \) be a set of labels. A model for \( \text{CTL}^*-\text{X} \) is a pair \( (F, V) \), where \( F = (W, (\overset{a}{\rightarrow})_{a \in T}, w^0) \) is a directed, rooted, edge-labeled graph with node set \( W \) and initial node \( w^0 \), while \( V \) is a valuation function that assigns to each node \( w \) the set of propositions \( V(w) \subseteq PV \) that are 'true' in \( w \). The edge relation is assumed to be total: i.e. \( \forall u \exists v. a \overset{a}{\rightarrow} v \). The labels on the edges in the definition of the graph are only used in the sequel for the benefit of the description of the suggested algorithm, but are ignored by the interpretation of the temporal logics.

Let \( M = ((W, (\overset{a}{\rightarrow})_{a \in T}, w^0), V) \) be such a model and let \( \pi = (v_0, v_1, \ldots) \) be an infinite path starting at \( v_0 \in W \) such that \( v_i \overset{a_i}{\rightarrow} v_{i+1} \) for every \( i \). Let \( \pi_i \) denote the suffix \( (v_i, v_{i+1}, \ldots) \) of \( \pi \).

Satisfaction of a formula \( \varphi \) in a state \( v \) of \( M \). (written \( M, v \models \varphi \) or just \( v \models \varphi \)), is defined inductively as follows:

S1. \( v \models q \) iff \( q \in V(v) \), for \( q \in PV \).

S2. \( v \models \neg \varphi \) iff \( v \not\models \varphi \). \( v \models \varphi \land \psi \) iff \( v \models \varphi \) and \( v \models \psi \).

S3. \( v \models A\varphi \) iff \( \pi \models \varphi \) for every path \( \pi \) starting at \( v \).

P1. \( \pi \models \varphi \) iff \( v_0 \models \varphi \) for any state formula \( \varphi \).

P2. \( \pi \models \neg \varphi \) iff \( v \not\models \varphi \). \( \pi \models \varphi \land \psi \) iff \( v \models \varphi \) and \( v \models \psi \).

P3. \( \pi \models \mathcal{U}(\varphi, \psi) \) iff there is an \( i \geq 0 \) such that \( \pi_i \models \psi \) and \( \pi_j \models \varphi \) for all \( 0 \leq j < i \).

Observe that satisfaction over our models coincides with the standard definition as provided in, e.g., [1].
Programs, State Graphs and Independence

For purposes of state graph generation and model checking, the specific syntactic structure of programs is not important. Instead, a finite-state program $P$ is viewed as a 4-tuple $(Q, T, q_0, I)$ where $Q$ is a finite set of states giving, e.g., the values of the variables and the contents of the message queues, $T$ is a finite set of operations such as assignments and send and receive actions to and from the message queues, $q_0 \in Q$ is the initial state, and $I$ is a so-called independence relation on the program’s actions that will be discussed below (see Definition 2.1). The enabling condition $en(q) \subseteq T$ is the set of operations that can be executed from a state $q$. Each operation $a \in T$ is identified with a partial function $a: Q \rightarrow Q$ (its denotation) that needs to be defined at least for each $q$ such that $a \in en(q)$.

We assume that $en(q) \neq \emptyset$ for any $q$.

Ignoring the independence relation for a moment, programs can directly be represented as state graphs, namely by taking $Q$ as nodes and $T$ for edges. The aim for reduction prompts to also consider subgraphs; a homomorphism is used in making the connection to the program.

A state graph, $G_P$, for a program $P$ is a directed, rooted and edge labeled graph $(S, (\stackrel{a}{\rightarrow})_{a \in T}, s^0)$, with nodes $S$ and initial node $s^0$ for which there is a homomorphism $st: S \rightarrow Q$ that maps nodes to program states such that (1) $st(s^0) = q_0$ and (2) if $s \stackrel{a}{\rightarrow} t$ then $a \in en(st(s))$ and $st(t) = a(st(s))$.

Hence, a state graph explicitly represents computations of the program as paths through the graph. In fact, because actions are functional (i.e., their denotations are), computations, i.e., paths in a state graph that start in the initial state, are uniquely determined by the sequence of actions that occur along them. Observe that a state graph need not contain all states of the corresponding program. If it does contain all reachable states and, moreover, if $st$ is an isomorphism with respect to the set of reachable states, we call it the full state graph. We shall assume state graphs not to contain unreachable states.

A program $P$ together with a valuation function $V: Q \rightarrow 2^{P^T}$ defines models $(G_P, \hat{V})$ for any state graph $G_P$ for $P$ and function $\hat{V}: S \rightarrow 2^{P^T}$ defined by $\hat{V}(s) = V(st(s))$. In the sequel, we shall not distinguish between $V$ and $\hat{V}$.

Partial order reduction exploits concurrency in programs and the fact that truth of specifications is often insensitive to the order in which so-called independent actions from different concurrent components occur in computations. Such independent actions can be, e.g., assignments to variables that are local to different components and send-actions in different components that affect separate message queues. The information as to which actions are independent can be given in an abstract way as follows:

Definition 2.1 An independence relation is an irreflexive and symmetric relation $I \subseteq T \times T$ such that for each pair of operations $(a, b) \in I$ (called independent operations), for each $q \in Q$,

- if $a \in en(q)$ (i.e., $a$ is enabled from $q$), then $b \in en(a(q))$ iff $b \in en(q)$.
- if $a, b \in en(q)$ then $a(b(q)) = b(a(q))$.

If $(a, b) \not\in I$, then $a$ and $b$ are called dependent.

3 Stuttering Equivalence

The correctness of the reduction method will be based on the notion of stuttering equivalence.

Let $M = ((W, (\stackrel{a}{\rightarrow})_{a \in T}, w^0), V)$ and $M' = ((W', (\stackrel{a'}{\rightarrow})_{a' \in T}, w'^0), V')$ be two finite models.

Definition 3.1 ([1]) A relation $\sim \subseteq W \times W'$ is a stuttering equivalence between $M$ and $M'$ if the following conditions hold:

1. $w^0 \sim w'^0$. 


2. If \( w \sim w' \), then \( V(w) = V'(w') \) and for every path \( \pi \) of \( M \) that starts at \( w \), there is a path \( \pi' \) in \( M' \) that starts at \( w' \), a partition \( B_1, B_2 \ldots \) of \( \pi \), and a partition \( B'_1, B'_2 \ldots \) of \( \pi' \) such that for all \( j \geq 0, B_j \) and \( B'_j \) are nonempty and finite, and every state in \( B_j \) is related by \( \sim \) to every state in \( B'_j \) and

3. the same condition as (2) interchanging \( \pi \) and \( M' \) with \( \pi' \) and \( M' \).

In [1] stuttering equivalence is defined using approximants \( \sim^n \). Because our models are finite, it is easy to see that the two definitions are equivalent.

Theorem 3.2 ([1]) Let \( \varphi \) be a \( \text{CTL}^+ - X \) formula with the set of atomic propositions \( PV \). Let \( M \) and \( M' \) be two models with the same set of atomic propositions \( PV \) and let the relation \( \sim \) be a stuttering equivalence between \( M \) and \( M' \). If \( w \sim w' \), then \( M, w \models \varphi \) iff \( M', w' \models \varphi \).

The reverse of this theorem, essentially stating that no coarser equivalence preserves the truth value of \( \text{CTL}^+ - X \) formulas, also holds [1]. This means that as far as the equivalence is concerned, stuttering equivalence enables maximal reduction.

4 The Algorithm

The reduction algorithm is based upon a modified depth-first-search algorithm. It generates a reduced state graph \( G' \) for the checked program \( P \) such that the correctness of any checked property \( \varphi \) under \( G' \) is the same as under the full state graph \( G \) of \( P \). This is guaranteed by ensuring that the model corresponding to \( G \) is stuttering equivalent with the model corresponding to \( G' \) (see Theorem 3.2).

The idea of the reduction is that from each state in the reduced state graph the set of enabled operations is examined, and only a subset of it is used to generate successors. This contrasts with the construction of the full state graph, where all of the enabled operations are expanded. The subset of the operations \( E(q) \) taken from a state \( q \) satisfies restrictions \( C0, \ldots, C3 \) below, in order to preserve stuttering equivalence between the full and the reduced model.

![Figure 1: Possibilities of reduction by not expanding a.](image)

To explain the restrictions imposed on the set \( E(q) \) let’s assume first that the full model \( M \) does not contain loops except for self loops. Definition 3.1 describes the cases in which the model \( M' \) resulting after removing a transition from \( M \) is stuttering equivalent with \( M \). In Figure 1 we have indicated the two situations in which the \( a \)-labeled transition need not be expanded from state \( q \) in \( M \). This is the case when the states \( q'' \) and \( q' \) are stuttering equivalent (denoted \( q'' \sim q' \)) or when the states \( q \) and \( q' \) are stuttering equivalent (\( q \sim q' \)) in the full model \( M \).

If \( q'' \sim q' \) or \( q \sim q' \), then for each path \( \pi \) with the prefix \( qq'' \), there is a path \( \pi' \) with the prefix \( qq' \) such that two partitions of \( \pi \) and \( \pi' \) satisfying Condition (2) of Definition 3.1 exist. Notice that it is because of the absence of non-trivial self-loops in \( M \) that the path \( \pi' \) cannot contain the transition.
\[ q \xrightarrow{a} q' \]. If it could, \( M \) would not need to be stuttering equivalent with \( M' \) since \( q \xrightarrow{a} q' \) is not present in \( M' \). Therefore, the full model \( M \) remains stuttering equivalent with the model \( M' \) obtained after removing the transition \( q \xrightarrow{a} q'' \) from \( M \).

As for ensuring \( q'' \sim q' \), we do not know of any efficiently checkable condition that would imply this. Indeed, the general problem is PSPACE-hard in the number of program operations, as it depends on the subgraphs of nodes reachable from \( q'' \) and \( q' \). Therefore, we concentrate on the second case: ensuring that \( q \sim q' \).

First observe that by repeatedly applying the argument above, it follows that if \( q \sim q' \) then it suffices to only have the transition \( q \xrightarrow{b} q' \) from \( q \) in \( M' \), i.e., the subtree of other transitions (indicated by the triangle in the figure) can be ignored. Hence, to reduce most effectively, we shall require the algorithm to use singleton sets whenever reduction is possible; thus we obtain

\[ C0 \] \( E(q) \) contains either all operations enabled in state \( q \), or exactly one of these; i.e., \( E(q) \) is a singleton.

The next condition will make sure that the execution of \( b \) does not change any propositional variable used in \( \varphi \) assigned to \( q \) and \( q' \), which is a necessary condition for \( q \sim q' \).

To present it, the following definition is needed:

**Definition 4.1** An operation \( a \in T \) is visible if it can change the truth value of some proposition in \( \varphi \); i.e., if \( V(a(q)) \neq V(\alpha(q)) \) for some state \( q \) such that \( a \in \text{en}(q) \). Let \( \text{Vis} \) be the set of visible operations in \( T \).

A good practical approximation for calculating \( \text{Vis} \) is as follows: Consider each program operation that can change one of the propositions as visible. (This approximation might not calculate the minimal set of visible operations, therefore allowing less reduction. However, it is safe in the sense that it would not miss a visible operation.)

In keeping with extant literature [21], this condition is called \( C2 \):

\[ C2 \] If \( E(q) \neq \text{en}(q) \), the operation in \( E(q) \) is not visible.

In formulating the subsequent conditions we use the fact that we have already imposed conditions \( C0 \) and \( C2 \).

The general problem of showing that \( q \sim q' \) still is PSPACE-hard in the number of program operations. So, we aim for a stronger condition: for every path \( \pi \) starting in \( q \) there is a path \( \pi' \) starting in \( q' \) that is the same up to invisible actions. Now, consider \( \pi \). As long as the actions along \( \pi \) are independent of \( b \) there is no problem in constructing \( \pi' \) because independent actions commute, so that these \( \pi \)-actions can still occur, in the required order, after the (invisible) \( b \)-action. Note that, like before, absence of non-trivial loops matters here. Dependent actions do cause a problem because there is no way to ensure that such actions can still occur without exploring all paths starting in \( q' \). So, we disallow this situation by stipulating that such actions can only occur after the \( b \)-action has occurred.

\[ C1 \] No operation \( a \in T \setminus E(q) \) that is dependent on the operation in \( E(q) \) can be executed in \( P \) before the operation from \( E(q) \) is executed.

Now, consider the first action \( c \) along \( \pi \) that depends on \( b \) and let \( \bar{q} \) be the state on \( \pi \) from which \( c \) is taken. Since the \( b \)-action must have occurred along \( \pi \) before reaching state \( \bar{q} \), commutativity of independent actions implies that the constructed prefix of \( \pi' \) ends in state \( \bar{q} \) from which \( c \) and, indeed, the whole sequence of subsequent actions along \( \pi \) can be taken.

Condition \( C1 \) occurs in many variations in LTL-X preserving reduction methods [12, 20, 23, 6].
The final restriction [20] is needed in case \( M \) does have non-trivial loops and restricts omitting operations along such loops. Consider the figure on the right, essentially extending Figure 1 with a non-trivial loop. Choosing \( E(q) = \{ b \} \), \( E(q') = \{ b' \} \), which satisfy \( C_0 \), \( C_1 \) and \( C_2 \) (when only \( a \) is visible), yields an incorrect reduction, where \( a \) is absent. \( C_3 \) prohibits completely omitting such operations occurring along a nontrivial loop. In the algorithm it is detected, (as closing a cycle on the search stack) when a node is part of a loop. It is at this node that the operation is added.

\( C_3 \) If \( E(q) \neq \operatorname{en}(q) \) then the operation in \( E(q) \) when applied to the current state \( q \), does not close a cycle on the search stack (i.e., we don't allow that an open node with value \( a(q) \) exists on the search stack).

Note that in the example \( C_3 \) only requires to add one of the two \( a' \), namely the one from \( q \) to \( q'' \). This in turn necessitates that \( q'' \) together with the corresponding \( b \) and \( b' \) are added to the reduced model. So only the \( a \) operation from \( q \) to \( q'' \) can be removed.

The conditions \( C_1, C_2, C_3 \) are sufficient to guarantee that the reduced state graph will preserve any checked linear temporal logic property \( \varphi \) [21]. The condition \( C_0 \) is newly introduced for branching temporal logics.

\[
Vis = \{ b, c \} \quad \text{Vis} = \{ d \}
\]

![Figure 2: Correct CTL*-X reductions](image)

Examples of reductions that use a subset satisfying the conditions \( C_0, \ldots, C_3 \) are shown in Figure 2. The dotted and solid transitions together constitute the full model. The reduced model consists of only the solid transitions.

The reduced state graph generation program is given in Figure 3. Each node \( s \) represents some program state \( st(s) \). The construction starts with a node whose state is the initial state \( q_0 \) of the program. The main program consists of a depth-first search (lines 4-17). Each new state is marked by the flag \( \text{open} \) (lines 2 and 12), and when its expansion is finished (i.e., it is removed from the search stack), it is marked by \( \text{closed} \) (line 16). For each new state, the subset of successors to be used is calculated by the procedure \( \text{ample} \) (the procedure call is in line 5, the procedure body is in lines 18-25). This procedure returns either a single (invisible) operation that satisfies conditions \( C_0, \ldots, C_3 \) (line 22) or the set of all enabled operations (line 21).

Checking that a singleton set \( \{ a \} \) satisfies condition \( C_1 \) is not detailed in the algorithm given in Figure 3. Because finding optimal ample sets is NP-complete ([20]), any implementation of \( \text{ample} \)
will use heuristics that may also depend on the specific programming language used (to define a finite state program in our sense).

Such heuristics are based on checking the type of the operation \( a \) (e.g., a local assignment, a synchronous receive operation, etc.) and some conditions on the rest of the program, and the state of the current node \( s \): according to the type of the operation, there are certain conditions whose satisfaction in the current state \( s \) guarantee that \( \{ a \} \) satisfies \( C1 \). For example, the simplest condition is that \( a \) is a local assignment, and is not within a non-deterministic choice with other operations. A slightly more complicated condition applies when \( a \) is a non-synchronous receive. Then \( C1 \) is guaranteed if there is no other receive operation from the same queue in any other process (this holds vacuously when a communication queue can be shared only by a pair of processes), and the queue is not empty in the state \( st(s) \). A more complete description of checking \( C1 \) appears in [11].

```
create_node(s, qo);
set open(s);
expand_node(s);

proc expand_node(s):
  working_set(s) := ample(st(s)); /* find set of operations to expand from s */
  while working_set(s) \= \emptyset do
    a := some operation of working_set(s);
    working_set(s) := working_set(s) \ {a};
    succ_state := a(st(s)); /* the a-successor of st(s) */
    if new(succ_state) then
      create_node(s', succ_state); /* s' has value succ_node */
      set open(s'); /* set s' to open, i.e., on the search stack */
      expand_node(s') fi; /* expand the successors of s' */
    create_edge(s, a, s');
  end while;
  set closed(s); /* close s, i.e., s is not on the search stack */
end expand_node.

proc ample(x) : element of T:
  foreach a \in T \ Vis /* for every invisible (due to C2) operation */
  if \( \{ a \} \) satisfies C1 for x and
    not exists s' with st(s') = a(x) and opw(s') /* Check C3 */
  then return(\{ a \}) fi /* singleton set, due to C0 */
end ample;
```

Figure 3: A reduced state graph expansion algorithm

Let \( M_P \) be the model corresponding to the full state graph \( G_P \) of given program \( P \) and \( M_r \) be the model corresponding to the reduced state graph \( G' \) generated by our algorithm. We have the following:

**Theorem 4.2** \( M_P \) is stuttering equivalent with \( M_r \).

The correctness of the algorithm is based on the relation found between \( \text{CTL}^*\)-X validity and a stuttering equivalence of models \([1]\). Proving the correctness of the above algorithm is based on constructing an equivalence relation \( \sim \) on the states of the full state graph \( M_P \) and showing that it
is a stuttering equivalence. By construction, the relation \( \sim \) relates only pairs of states that have the same set of atomic propositions. Moreover, for each \( v \sim w \) such that \( w \) remain in the reduced state graph \( M_r \), if \( \pi \) is a finite or infinite sequence from \( v \), then there exists a sequence \( \pi' \) from \( w \) such that

- \( \pi' \) contains only edges that remain in the reduced state graph \( M_r \), and
- \( \pi \) and \( \pi' \) can be partitioned into segments that pairwise correspond according to the conditions 2 and 3 of Definition 3.1.

Since the initial state \( w^0 \) remains in the reduced state graph, the above conditions guarantee that the reduced state graph \( M_r \) simulates the full state graph \( M_f \). The other direction of the bisimulation stems from the fact that \( \sim \) is an equivalence relation, hence reflexive. Thus, it follows from Theorem 3.2 that the reduction preserves the validity of the checked formula.

The construction of the relation \( \sim \) and the detailed proof are delegated, due to lack of space, to an appendix.

### 4.1 Complexity of the Algorithm

The time complexity of the algorithm is \( O(n_r \cdot C + m_r) \), where \( n_r \) is the number of states in the generated state graph, \( m_r \) is the number of edges and \( C \) is the complexity of computing an ample set. This is obvious as the algorithm is a modified depth-first search through the state graph. Computing ample sets can be done in constant time along the lines of the earlier explanation. As to the amount of space, this is clearly linear in the number of states and edges. Hence we obtain an \( O(n_r + m_r) \) space and time complexity for the algorithm.

### 5 Experimental Results

#### 5.1 Reduced State Space for Various Algorithms and Protocols

The algorithm described in this paper was implemented by Gerard Holzmann in SPIN [9] and run on several examples. The table in figure 4 below contains the number of states and edges, memory used in bytes, and time in seconds of generating full state graphs and reduced state graphs for both LTL-X and CTL-X. The reduction for CTL-X contains an additional restriction, namely \( C_0 \), on selecting the subset of successors. This restricts the subsets of successors to be either the full set of the enabled operations, or a singleton set. In order to make the comparison unbiased towards any particular checked property, all operations were considered invisible during the tests. The set of properties that can be checked without making any program operation visible includes properties such as deadlock and termination.

All measurements were made on a Sparc-10 workstation with 128Mbyte of RAM. The runtimes are the sum of system-time and user-time. The algorithms checked are as follows: leader is a leader election algorithm for an unidirectional ring [3], sorting is a pipeline distributed sorting algorithm, urp is AT&T universal receiver protocol, dltp is a data transfer protocol, snoopy is a cache coherence protocol, pftp is a file transfer protocol [9] and lpc is a model of a telephone switch. For the first two examples, leader and sorting, the reduction with and without the additional restriction \( C_0 \) are the same. Both give (when repeated with different numbers of processes) an exponential reduction of the state graph. For urp and dltp, the reduction in space and time is very similar with and without the additional restriction. For snoopy, the CTL-X reduction generates a state graph that is about 25% bigger in space, and takes about 50% more time. For pftp, the reduction is about twice better in space and time for the LTL-X reduction, and for lpc, it is about 2.5 times better in space and more than three times better in time.

9
<table>
<thead>
<tr>
<th>Algorithm</th>
<th>Reduction</th>
<th>States</th>
<th>Transitions</th>
<th>Memory</th>
<th>Time</th>
</tr>
</thead>
<tbody>
<tr>
<td>leader</td>
<td>Non</td>
<td>382,061</td>
<td>1,847,294</td>
<td>88,029,016</td>
<td>97.1</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>94</td>
<td>94</td>
<td>1,112,896</td>
<td>(&lt;0.1)</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>94</td>
<td>94</td>
<td>1,116,304</td>
<td>(&lt;0.1)</td>
</tr>
<tr>
<td>sorting</td>
<td>Non</td>
<td>659,683</td>
<td>3,454,989</td>
<td>113,629,016</td>
<td>145.5</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>182</td>
<td>182</td>
<td>1,120,088</td>
<td>(&lt;0.1)</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>182</td>
<td>182</td>
<td>1,124,496</td>
<td>0.1</td>
</tr>
<tr>
<td>urp</td>
<td>Non</td>
<td>4,329</td>
<td>16,563</td>
<td>1,943,416</td>
<td>0.9</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>1,149</td>
<td>2,160</td>
<td>1,542,008</td>
<td>0.2</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>1,231</td>
<td>2,388</td>
<td>1,598,304</td>
<td>0.3</td>
</tr>
<tr>
<td>dtp</td>
<td>Non</td>
<td>251,109</td>
<td>648,467</td>
<td>36,369,976</td>
<td>32.2</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>16,159</td>
<td>17,603</td>
<td>4,290,104</td>
<td>1.5</td>
</tr>
<tr>
<td></td>
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<td>17,100</td>
<td>18,320</td>
<td>4,540,224</td>
<td>1.6</td>
</tr>
<tr>
<td>snoopy</td>
<td>Non</td>
<td>191,498</td>
<td>938,719</td>
<td>24,446,520</td>
<td>48.2</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>34,541</td>
<td>65,468</td>
<td>6,080,056</td>
<td>5.1</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>46,329</td>
<td>99,733</td>
<td>7,620,416</td>
<td>7.5</td>
</tr>
<tr>
<td>pftp</td>
<td>Non</td>
<td>1,067,787</td>
<td>4,545,778</td>
<td>170,410,248</td>
<td>625.5</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>136,208</td>
<td>257,019</td>
<td>23,565,880</td>
<td>30.1</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>259,252</td>
<td>597,819</td>
<td>43,120,448</td>
<td>59.5</td>
</tr>
<tr>
<td>tpc</td>
<td>Non</td>
<td>3,912,286</td>
<td>11,762,426</td>
<td>251,827,768</td>
<td>7202.9</td>
</tr>
<tr>
<td></td>
<td>LTL-X</td>
<td>391,534</td>
<td>461,534</td>
<td>26,863,160</td>
<td>31.0</td>
</tr>
<tr>
<td></td>
<td>CTL-X</td>
<td>977,454</td>
<td>1,727,034</td>
<td>64,472,896</td>
<td>98.1</td>
</tr>
</tbody>
</table>

Figure 4: CTL-X versus LTL-X reductions

These results demonstrate that the inclusion of the reduction algorithm is beneficial for all the above examples. A substantial reduction can be achieved with relatively small cost, as the implementation of the reduction algorithm is simple and incurs only very small overhead (for further implementation details, refer to [11], where an efficient LTL-X implementation is described). Even in the cases where the reduction is not very big (in comparison to other reductions, as for the leader algorithm), such as in the tpc algorithm, where the gain in space is a factor of four, one can obtain a considerable benefit: since the algorithm is complicated enough to consume a large amount of memory, even the fourfold memory reduction could reduce the execution time from over two hours to about a minute and a half (avoiding needless memory swaps).

5.2 Verifying Branching Bisimulation

The reduction method described in Section 4 can be further exploited in the context of process algebra. It can be used to verify whether two states of a program are branching bisimilar [15, 19].

Let \( \text{Invis} \) denote the set of invisible operations, i.e., \( \text{Invis} = T \setminus \text{Vis} \). As we identify each operation from \( \text{Invis} \) with a silent step, the definition of branching bisimulation can be formulated as follows:

**Definition 5.1** A relation \( \subseteq \subseteq W \times W' \) is a branching simulation between \( M \) and \( M' \) if it satisfies the following conditions:

(a) \( w^0 \subseteq w'^0 \) and

(b) if \( w \subseteq w' \), then if \( w \xrightarrow{a} v \), then either \( a \in \text{Invis} \) and \( v \subseteq w' \) or there are states \( w' = w'_0, w'_1, \ldots, w'_{n+1} = v' \) such that for each \( 0 \leq i < n \), \( w'_i \xrightarrow{a'_i} w'_{i+1}, a_i \in \text{Invis}, w \subseteq w'_{i+1}, \) and
\[ w'_n \xrightarrow{a'} v', v \subseteq v'. \]

A relation \( \sim \) is a branching bisimulation ([18]) if both \( \sim \) and \( \sim^T \) (the transpose of \( \sim \)) are branching simulations.

\( M \) and \( M' \) are branching bisimilar if there is a branching bisimulation relation between \( M \) and \( M' \).

We have the following:

**Theorem 5.2** Given a program \( P \), \( M_p \) is branching bisimilar with \( M_r \).

The proof is given in the Appendix.

Verifying whether or not two states are branching bisimilar is done by constructing the minimal branching bisimilar state graph of the program. The most efficient algorithm for this was published by Vaandrager and Groote in [8] and has time and space complexity of \( O(n^2 + nm) \) and \( O(n + m) \), respectively, where \( n \) and \( m \) are the number of states and edges in the full state graph. For determinate programs, [22] presents a more efficient algorithm with time and space complexity \( O(n \log n + m) \) and \( O(n + m) \), respectively. In [8] it was conjectured that the same time complexity suffices for minimizing arbitrary state graphs.

By Theorem 5.2, our algorithm generates a state graph that is branching bisimilar to the full state graph. Moreover it has time and space complexity \( O(n_r + m_r) \), where \( n_r \) and \( m_r \) are the numbers of generated states and edges. This raises the possibility of using our algorithm as a preprocessing phase to constructing a minimal branching bisimilar state graph, thus allowing the minimization algorithm to run in time \( O(n^2 + n_r + m_r) \) using \( O(n_r + m_r) \) space. The algorithms [8, 22] can be applied to a reduced state graph, instead of the full one.

The 'benchmark' example used in literature, and by us, is Milner's scheduler as described in [17]. This is a simple token ring consisting of \( k \) cyclic processes \( C_i \), which, on having received the token, communicate with some system and then concurrently wait for acknowledgment and wait to pass on the token. Process \( C_i \) is described by the CCS equation \( C_i = \bar{t}_i \cdot c_i \cdot (\bar{a}_i \mid t_{\text{mod}+1}) \cdot C_i \). The complete scheduler on \( k \) processes is described by \( \text{Sch}_k = (t_1 \cdot \text{nil} \mid C_1) \cdots (C_k \mid t_k \cdot \text{nil}) \), where the first process starts \( C_1 \). The operator \( \cdot \) denotes a concurrent composition, and \( \cdot \) means sequencing. Letters correspond to operations. Two letters, where one is barred, e.g., \( e \) and \( \bar{e} \), may synchronize, thus producing an invisible action. The operator \( \cdot \) is the restriction operator which, in this case, forces the \( t_i, \bar{t}_i \) synchronizations to occur. \text{nil} is the idle process that does nothing.

In Figure 5 we have collected some results for various sizes \( k \) of the token ring. The measurements were done on a Sparc1+ workstation with 16MB of memory. The actual generation of both the reduced and the full state graphs are achieved by a script written in PERL, an interpreted language with heavy use of pattern matching, and the absolute times should be interpreted accordingly. A C implementation can be expected to run at least an order of magnitude faster. The number of states and edges in the full state graph is given in the 2nd and 3rd columns. We consider both the case that only the communication actions \( (c_i) \) are visible and the case that both \( c_i \) and the acknowledgment actions \( (\bar{a}_i) \) remain visible. For both cases we give the sizes of the state graphs as generated by our algorithm and the minimal state graphs (as given by an implementation of the Vaandrager/Groote algorithm, part of the PSF [13] toolkit). The line column gives the time in seconds that our algorithm needs to generate the state graph, where only \( c_i \) is visible. We see that in this case not only that the resulted reduced state graphs is small but the time to generate them is small as well. This should be contrasted with the figures in the 2nd to last column that give the times it takes to generate the full state graph. The time for the actual minimalization of the reduced state graph is negligible for these sizes of state graphs. In other words, one gains considerably here by generating the state graph using our reduction algorithm.
**Table 1: Verifying Branching Bisimulation**

<table>
<thead>
<tr>
<th>Milner's scheduler</th>
<th>Only c_i visible</th>
<th>Both c_i and a_i visible</th>
</tr>
</thead>
<tbody>
<tr>
<td>k</td>
<td>states(s)</td>
<td>edges(e)</td>
</tr>
<tr>
<td>4</td>
<td>97</td>
<td>241</td>
</tr>
<tr>
<td>5</td>
<td>241</td>
<td>721</td>
</tr>
<tr>
<td>6</td>
<td>577</td>
<td>2017</td>
</tr>
<tr>
<td>7</td>
<td>1,345</td>
<td>5,377</td>
</tr>
<tr>
<td>8</td>
<td>3,073</td>
<td>13,825</td>
</tr>
<tr>
<td>9</td>
<td>6,913</td>
<td>34,561</td>
</tr>
<tr>
<td>10</td>
<td>15,361</td>
<td>84,481</td>
</tr>
<tr>
<td>11</td>
<td>33,793</td>
<td>202,733</td>
</tr>
<tr>
<td>50</td>
<td>151</td>
<td>131</td>
</tr>
<tr>
<td>100</td>
<td>301</td>
<td>301</td>
</tr>
</tbody>
</table>

*Using a PERL script for state graph generation.

Figure 5: Verifying branching bisimulation

A second experiment shows that even in the case when our algorithm does not substantially reduce the state graph, the overhead of doing the reduction is negligible. In this case, both c_i and a_i are made visible and the reduction of the number of edges is only between 11% (k=4) and 26% (k=8). Here, more than a half of the operations are visible, which defies most of the reduction. This is fortunately untypical. Furthermore, one can see that the minimal state graph also grows exponentially, producing a minimal state graph that is only about 50% smaller than the full state graph. The 3rd to last and 2nd to last columns, marked as "full" and "PO", show the time it takes to generate the full state graph and reduced state graph, respectively. One can see that even though in this case the reduction is small, the overhead that our algorithm incurs is minimal when compared to generating the full state graph; in fact, the algorithm still runs a little faster.

Although minimizing a state graph w.r.t. branching bisimulation is a global process, certain equivalence preserving transformations can be done locally during state graph generation. For instance, states that have precisely one outgoing transition can be removed if that transition is invisible\(^1\). The column labeled 'PO & τ-removal' shows the result of augmenting our partial order algorithm with invisible-step removal. There is now a reduction in the number of states as well as a more substantial reduction in the edges. Interestingly, the resulting state graphs are in fact the minimal branching bisimilar ones. The last column shows that there is no time penalty. In fact, the running times are almost the same, which is not surprising because the algorithm has to visit the same number of nodes as before. Note however that the minimization algorithm will run in time and space proportional to the size of the reduced graph. Hence, invisible-step removal is advantageous for the minimization phase.

6 Conclusions

We have presented an algorithm for generating reduced state graphs to be used for model-checking branching temporal properties. The usual DFS expansion algorithm was modified so that only subsets of the successors from each state are expanded. This allows reducing the number of states and edges,

\(^1\)More formally: the rewrite rule \(x \cdot \tau \rightarrow x\) preserves branching bisimulation.
and thus allows reducing the space and time used for this construction and for model checking. The branching time logics include the temporal logic $CTL^*\times$, which is more expressive than the linear time logic $LTL\times$. They also include the logic $CTL\times$ which has a model-checking algorithm that is linear in the size of the checked property [2]. These advantages in either expressiveness or efficiency can now be combined with the ability to reduce the state graph using partial order methods.

On the other hand, we have shown that, in general, the reduction of the state graph for preserving branching properties is more restricted than the one for $LTL\times$: an additional restriction was added, limiting the subset of successors taken from each state to be either the full set of successors or a singleton set.

Experimental results show that the suggested algorithm results in a substantial reduction in both space and time over the traditional full state graph exploration. Also, the algorithm proved to be the preferred way to generate state graphs to verify branching bisimulation.

Acknowledgments The authors would like to thank Gerard Holzmann for implementing the additional constraint into the partial order version of SPIN, and helping with the experiments. Also, hats off for Larry Wall's Perl, which proved to be an excellent rapid prototyping tool as well.

References

A Appendix: Correctness of the Algorithm

Let $M_P = ((W, \rightarrow, w^0, L), V)$ be a model, where $(W, \rightarrow, w^0, L)$ is the full state graph of $P$ and let $M_r = ((W_r, \rightarrow, w^0, L_r), V_r)$ be a model with a reduced state graph as generated by the algorithm, where $w^0 \in W_r \subseteq W$, $\rightarrow_r \subseteq \rightarrow$, $L_r = L | \rightarrow_r$ (reads $L$ restricted to the states that participate in the reduced transition relation $\rightarrow_r$) and $V_r = V | W_r$. Because of Theorem 3.2, we just have to prove that $M_P$ is stuttering bisimilar with $M_r$.

To do so we assign to each state $w \in W$ in $M_P$ the set of operations $E(w) \subseteq T$ such that

- if $w \in W_r$, then $E(w)$ is equal to the set of operations expanded by the algorithm from $w$; i.e., $E(w)$ is equal either to $en(w)$ or to one invisible operation from $en(w)$ satisfying the conditions C1–C3.

- if $w \notin W_r$, then $E(w) = en(w)$; i.e., $E(w)$ is equal to the set of enabled operations at $w$.

In the sequel, let $M' = ((W', \rightarrow', w^0', L'), V')$ be the model obtained from $M_P$ by removing all the transitions $w \xrightarrow{a} w'$ such that $a \notin E(w)$. Hence, $w^0' = w^0$. One can easily notice that the reachable part of $M'$ is equal to $M_P$. So, to prove that $M_P$ is stuttering bisimilar with $M_r$ we just have to show that $M'$ is stuttering bisimilar with $M_P$.

We now discuss how such a stuttering equivalence is obtained. The following notation will be used, where $\sim$ is an equivalence to be fixed in the sequel:

- $w \xrightarrow{a} w'$ denotes $w \xrightarrow{a} w'$ and $w \sim w'$.

- $w \xrightarrow{\sim} E w'$ denotes $w \xrightarrow{a} w'$ and $a \in E(w)$.

- $w \xrightarrow{\sim} E w'$ denotes $w \xrightarrow{a} w'$ and $a \in E(w)$.

- when convenient, we will sometimes omit labels of transitions writing $w \xrightarrow{\sim} E w', w \Rightarrow_E w'$ etc.

As $M'$ is a submodel of $M_P$, we shall simplify notation by constructing a stuttering equivalence on $M_P$. We start out by building a relation $\sim$ in $M_P$ which, as shown in Lemma A.7, only relates states with identical valuations. Furthermore, as shown in Lemmas A.11 and A.13, $\sim$ satisfies the following properties:

P1) if $v \sim w$, $v \in W' \subseteq W$ and $v \xrightarrow{r} v'$ and $\sim (v \sim v')$, then there is a non-empty sequence of transitions $v = v_0 \Rightarrow_E w_1 \Rightarrow_E \cdots \Rightarrow_E w_k \xrightarrow{r'} v'$ in $M_P$ such that $r' \sim w'$.

P2) if $v \sim w$ and there is an infinite stutter path $v = v_0 \Rightarrow_E w_1 \Rightarrow_E \cdots$, then there is an infinite stutter path $w = w_0 \Rightarrow_E w_1 \Rightarrow_E \cdots$.

The following theorem shows that $\sim$ is a stuttering equivalence consisting only of transitions which are not removed from $M_P$.

Theorem A.1 The relation $\sim$ is a stuttering equivalence on $M_P$ such that if $v \sim w$, $w \in W' \subseteq W$ and $\pi$ is a path starting at $v$, then a corresponding path $\pi'$ (as in Definition 3.1) can be taken, starting with $w$, and consists of $\sim$ transitions only.

Proof. By definition of $\sim$, $w^0 \sim w^0$. Because $\sim$ is symmetrical, it suffices to establish Clause 2 of Definition 3.1. Let $v \sim v'$. Then $V(v) = V(v')$ holds by definition. Next, let $\pi$ be a path $v = v_0 \xrightarrow{r_1} v_1 \xrightarrow{r_2} \cdots$ starting at $v$. Partition $\pi$ into blocks $B_0, B_1, \ldots$, where every $B_j$ is a maximal subsequence of $\pi$ of $\sim$-related states. Every $B_j$ is finite, except for possibly the last block. If $B_0$ is finite with last state $v_k$, then $v_k \sim r'$ and $v_k \not\sim v_{k+1}$. By Lemma A.11. Property P1 holds so that there
is a non-empty sequence of transitions $v' = v'_0 \Rightarrow E \cdots \Rightarrow E v'_{k+1}$ s.t. $v_{k+1} \sim v'_{k+1}$. We take $B'_0 = v'_0 \cdots v'_k$ and have $v_i \sim v'_i$ for every $v_i \in B_0$ and $v'_i \in B'_0$. Because $v_{k+1} \sim v'_{k+1}$, this sets us up to construct $B'_1, B'_2, \ldots$ as long as the corresponding $B_j$'s are finite. Finally, let some $B_n$ be infinite with starting state $v_r$ and let $B'_{n-1}$ have $v'_{n-1}$ as last state, so that $v_r \sim v'_r$.

By Lemma A.13, $\sim$ satisfies Property P2, so that there is an infinite path $v'_s = v'_0 \Rightarrow E v'_1 \Rightarrow E \cdots$. As all states in $B_n$ and the constructed path starting in $v'_s$ are $\sim$-related, we can partition these suffixes into finite blocks any way we want. This concludes the proof as an infinite block obviously is the last block. □

**Corollary A.2** $M_P$ is stuttering equivalent with $M_r$.

**Proof.** We have $M_P$ is stuttering equivalent with $M_r$. This is immediate from Theorem A.1 and the fact that all $\neg E$ steps are present in $M_r$. As the reachable part of $M'$ and $M_r$ are isomorphic as models, $M_r$ and $M'$ are stuttering equivalent. The Corollary follows by transitivity. □

Thus, by corollary A.2, the reduction is correct.

The proof proceeds by first defining an equivalence that satisfies weaker properties.

**Definition A.3** Let $M = ((W, \rightarrow, w_0, L), V)$ and $M' = ((W', \rightarrow', w_0', L'), V')$ be two finite models.

A relation $\subseteq W \times W'$ is a divergence blind stuttering simulation between $M$ and $M'$ if it satisfies the following conditions:

(a) $w_0 \subseteq w_0'$ and

(b) if $w \subseteq w'$, then $V(w) = V'(w')$, and if $w = w'$, then there are states $w = w_0, w_1, \ldots, w_n = w'$ such that for each $i < n$, $w_i \rightarrow' w_{i+1}$ and $w \subseteq w_i$, and $v \subseteq v'$. This is called the transfer property.

A relation $\sim$ is a divergence blind stuttering equivalence ([18]) if both $\sim$ and $\sim^T$ (the transpose of $\sim$) are divergence blind stuttering simulations.

$M$ and $M'$ are divergence blind stuttering equivalent if there is a divergence blind stuttering equivalence relation between $M$ and $M'$.

The desired divergence blind stuttering equivalence is constructed as follows:

**Definition A.4** Let $\subseteq \subseteq W \times W'$ be a relation in $M_P$, defined as $U_{b \in \omega} F_b^k(\subseteq_k)$ where

(a) $\subseteq_k = \{(u, v) \mid u = v = r \text{ or } u \in W_r, v \rightarrow b \in E(n) \neq e(n)\}$

(b) for $R \subseteq W \times W$, $F_b^0(R) = R$. $F_b^{i+1}(R) = F_b^i(R) \cup F_b(F_b^i(R))$, and

(c) $F_b(R) = \{(u, v) \mid \exists u', v'. a \in R, u' \rightarrow b v', u' \rightarrow a u, v' \rightarrow a v, a = b\}$

The next lemma gives two properties of $\subseteq$, needed in the rest of the proof.

**Lemma A.5** The relation $\subseteq$ has the following properties.

(a) If $w \subseteq z$ and $w \neq z$, then for some $b$, $w \rightarrow b z$ and $(u, b) \in I$ for all $a \in e(n) - \{b\}$.

(b) $\subseteq$ is a divergence blind stuttering simulation in $M_P$.

**Proof.** (a) We prove by induction on $k$ the property $\forall u, r, b \forall k P_b^k(u, v)$, where $P_b^k(u, v)$ is defined as:
if $(u, v) \in F_b^k(\mathbb{E})$ we have both $V(u) = V(v)$ and if $u \neq v$ then there are $u_0$ and $v_0$ such
that $u_0 \in W_r$, $u_0 \rightarrow_v b \rightarrow v_0$, $b \in E(u_0) \neq \epsilon(u_0)$ (hence, $u_0 \mathbb{E} v_0$), $(u, v) \in F_b^k(\{(u_0, v_0)\})$,
\[ u \rightarrow_b v \text{ and } \forall a \in \epsilon(u) - \{b\} \text{ } (a, b) \in I. \]

Obviously, this implies the first part of the Lemma. For the proof below, fix some action $b$ and nodes $u, v$ such
that $(u, v) \in F_b^k(\mathbb{E})$.

For $k = 0$, we have $u \mathbb{E} v$ so that $u = v$ or $u \rightarrow_b v$ for some invisible (by C3) $b$. In either case
$V(u) = V(v)$. To establish the remainder, take $u_0 = u$ and $v_0 = v$. Since we may assume $u \neq v$, then
$u \in W_r$. As $E(u) \neq \epsilon(u)$, we have $\forall a \in \epsilon(u) - E(u)$ $(a, b) \in I$ by Property C1. This suffices as
$E(u) = \{b\}$ by Property C0.

For $k > 0$, we may assume that $(u, v) \in F_b^k(\mathbb{E}) \cap F_b^{k-1}(\mathbb{E})$. If $u = v$ then $V(u) = V(v)$, so
assume otherwise. Since $(u, v) \in F_b^k(\mathbb{E}) \cap F_b^{k-1}(\mathbb{E})$, $u \neq v$. There are $u', v'$ and $a$ such that $(u', v') \in$ F_b^{k-1}(\mathbb{E}), $u' \rightarrow_b v'$, $u' \rightarrow_a u$, $v' \rightarrow_a v$ and $a \neq b$. As $u \neq v$, we must have $u' \neq v'$, by definition of a
model. Since $P_b^{k-1}(u', v')$ holds by induction, we obtain $u_0$ and $v_0$ such that $u_0 \in W_r$, $u_0 \rightarrow_b v_0$,
b \in E(u_0) \neq \epsilon(u_0)$, $(u', v') \in F_b^{k-1}(\{(u_0, v_0)\})$ and $u' \rightarrow_b v'$. Moreover, $\forall a \in \epsilon(u') - \{b\}$, $(a, b) \in I$.

Hence, $u \rightarrow_b v$ and $V(u) = V(v)$ because $b$ is invisible. It follows that $(u, v) \in F_b^k(\mathbb{E})$ and
$(a, b) \in I$ for any $a \in \epsilon(u) - \{b\}$ by Property C1 and because $\epsilon(u) \cap E(u) = \{b\}$ by Property C0.

We conclude that $P_b^k(u, v)$ holds.

(b) As for proving that $\mathbb{E}$ is a branching simulation. Clearly $u^0 \mathbb{E} u^0$. Now, set $u \mathbb{E} v$ and
\[ u \rightarrow_a v'. \] This means that $P_b^k(u', v)$ holds for some $k$ and $b$ so that $V(u) = V(v)$. If $u = v$ then
the transfer property (of Definition A.3 (b)) holds because $\forall c \mathbb{E} u$. Otherwise, for some $b, u \rightarrow_b v$ and
$\forall c \in \epsilon(u) - \{b\}$ we have $(b, c) \in I$. If $a = b$ then the transfer property holds, again, because $v \mathbb{E} v$.
Finally, if $a \neq b$ then $(a, b) \in I$ and, hence, $\forall c \mathbb{E} v'$ for some $v'$. We obtain $v' \mathbb{E} v'$ by definition of
$\mathbb{E}$ and this establishes the transfer property and completes the proof. 

**Definition A.6** The relation $\mathbb{E}$ is the symmetric and transitive closure of $\mathbb{E}$.

Next we will prove that $\mathbb{E}$ is a divergence blind stuttering equivalence.

**Lemma A.7** The relation $\mathbb{E}$ is a divergence blind stuttering equivalence in $MP$.

**Proof.** Firstly we prove that $\mathbb{E}$ is a divergence blind stuttering simulation. Clearly $w^0 \mathbb{E} w^0$. Now,
\[ u \mathbb{E} v \text{ and } u \rightarrow_a u'. \] So, $\forall c \mathbb{E} u$. This means that $P_b^k(v, u)$ holds for some $k$ so that $V(u) = V(v)$.
If $u = v$ then the transfer property (of Definition A.3 (b)) holds because $\forall c \mathbb{E} v$. Otherwise, by Lemma
A.5 (a), for some $b, v \rightarrow_b u$. Therefore, the transfer property holds, again, because $v \mathbb{E} u$ and $u' \mathbb{E} u$.

Therefore, by definition, $\mathbb{E}$ is a divergence blind stuttering equivalence and its symmetric and
transitive closure is as well. 

In order to prove that $\mathbb{E}$ satisfies the property P1, we need two auxiliary Lemmas A.8 and A.9.

The first lemma states that if a transition labeled $a$ is removed from some state, then a sequence of non-removable
stuttering transitions (with labels independent on $a$) followed by a non-removable transition labeled $a$ will start at
that state.

**Lemma A.8** Let $w \in W' \subseteq W$ and $a \in \epsilon(w)$. Then, there is a (possibly empty) sequence of
transitions $w = w_0 \rightarrow_{V'} b_1 \rightarrow_{V'} b_2 \rightarrow_{V'} \rightarrow_{V'} w_k$ and the transition $w_k \rightarrow_{V'} a \rightarrow_{V'} v'$ in $MP$ such that $w_i \mathbb{E} w_{i+1}$,
$w_i \neq w_{i+1}$, $b_i \neq a$ for $i < k$, and $b_1 \cdots b_k a \equiv_1 a b_1 \cdots b_k$. 

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Proof. By induction on the order that the algorithm 'closes' nodes, i.e., removing the node from the search stack, after its expansion is finished. Setting it to closed at line 16 in Figure 3. We claim that the above property holds for each node when it closes, and continues to hold throughout the execution. At the end, all the nodes of $M_p$ are closed, which proves the Lemma. Now, to prove the claim, distinguish between two cases:

1. $E(w) = en(w)$ (all the operations are taken). In this case, the Lemma holds trivially, with an empty sequence of b's. This case also correspond to the base of the induction.

2. $E(w) \subset en(w)$; i.e., $E(w)$ is a proper subset of enabled operations. Choose some $b \in E(w)$, which is invisible by C2. Then by Definition A.1, there is a transition $w \xrightarrow{b} w'$. Since by C3, this transition cannot close a cycle it holds that $w \neq w'$. It is a property of the DFS (depth-first-search) algorithm, that if this transition does not close a cycle, then $w'$ is closed before $w$ is closed. Thus, by the inductive hypothesis, when $w$ is closed, $w'$ already satisfies the condition of the Lemma: i.e. there is a path $w' = w_0 \xrightarrow{b_1} w_1 \xrightarrow{b_2} \ldots \xrightarrow{b_k} w_k \xrightarrow{a} w'$ such that $w_i \subseteq w_{i+1}$, $w_i \neq w_{i+1}$ and $b_{i+1} \neq a$ for $i < k$. Now, according to C1, all the operations that are executed before $a$ from $w$ are independent of $a$, which means that the path $w \xrightarrow{b} w_0 \xrightarrow{b_1} \ldots \xrightarrow{b_k} w_k \xrightarrow{a} w'$ satisfies the conditions of the Lemma.

The next lemma says that if at some state a sequence of stuttering transitions starts, then also a sequence of non-removable stuttering transitions starts such that the ending states of these sequences are connected by a sequence of stuttering transitions.

**Lemma A.9** Let $w \in W' \subseteq W$. If there is a non-empty sequence of transitions $w = w_0 \xrightarrow{a_1} w_1 \xrightarrow{a_2} \ldots \xrightarrow{a_n} w_n$, then there is a non-empty sequence of non-removable transitions $w = w_0 \xrightarrow{E} w_1 \xrightarrow{E} \ldots \xrightarrow{E} w_n$ such that $w_i \subseteq w_{i+1}$ and $w_i \neq w_{i+1}$ for $i < n$.

**Proof.** By induction on $n$. Consider $n = 1$. If $a_1 \notin E(w_0)$, then we are done. Assume that $a_1 \notin E(w_0)$. By Lemma A.8, there is a sequence of transitions $w = w_0 \xrightarrow{a_1} w_1 \xrightarrow{b_1} w_2 \xrightarrow{b_2} \ldots \xrightarrow{b_k} w_k$ and the transition $w_k \xrightarrow{a_1} w'$ such that $w_i \subseteq w_{i+1}$, $w_i \neq w_{i+1}$ for $i < k$, and $b_1 \ldots b_k a_1 = a_1 b_1 \ldots b_k$.

Since $a_1$ is independent of all $b_i$, there are states $w'_i$ such that $w_1 = w_0 \xrightarrow{b_1} w_1 \xrightarrow{b_2} w_2 \xrightarrow{b_3} \ldots \xrightarrow{b_k} w_k = w'$ and $w'_i \xrightarrow{a_1} w'_i$, for $i < k$. Because of $w'_i \subseteq w'_{i+1}$ for $i < k$, it follows from the definition of $E$ that $w_i \subseteq w_{i+1}$ for $i < k$.

Notice that, by assumption, $w_0 \sim w_1$ and by definition of $\sim$, $w_1 \sim w'$. Thus, by transitivity of $\sim$, $w_0 \sim w'$. Next, by definition of $\sim$, $w_0 \sim w'_i$. It follows from the definition of $\sim$ that $w'_i \sim w'$. If any transition labeled by $b_i$ is a self-loop, then we remove the state $w_{i-1}$ from the sequence $w'_i = w_0 \xrightarrow{b_1} w_1 \xrightarrow{b_2} \ldots \xrightarrow{b_k} w_k = w'$, which completes the proof for $n = 1$.

Inductive step. Consider a sequence $w = w_0 \xrightarrow{a_1} w_1 \xrightarrow{d_{k+1}} w_{k+1}$. By the inductive hypothesis, there is a non-empty sequence of non-removable transitions $w = w_0 \xrightarrow{E} w_1 \xrightarrow{E} \ldots \xrightarrow{E} w'_k$ and there is a (possibly empty) sequence of transitions $w_k = w_0 \xrightarrow{d_1} w_1 \xrightarrow{d_2} \ldots \xrightarrow{d_m} w_m = w_k$ such that $w_k \subseteq w_{k+1}$ and $w_k \neq w_{k+1}$ for $i < m$.

There are now two possibilities: either $a_{k+1} \neq d_i$ for all $i \leq m$ or there is some $d_j = a_{k+1}$. If $a_{k+1} \neq d_i$ for all $i \leq m$, then from Lemma A.5(a) we have $(a_{k+1}, d_i) \in I$ for all $i \leq m$ and therefore
we can complete the "rectangle" such that \( w_k \xrightarrow{a_{k+1}} w_{k+1} \) and there is a sequence of transitions \( w_k = w_0 \xrightarrow{d_1} w_1 \xrightarrow{d_m} w_{k+1} \). By definition of \( \mathcal{E}_c, w_{k+1} \subseteq w_{k+1}^+ \) for \( i < m \). If any of the transitions labeled by \( d_i \) is a self-loop, then we remove the state \( w_k \) from the sequence. For \( a_{k+1} \) from \( w_i \) we apply the construction shown for the case \( n = 1 \).

Assume now that \( j \) is the smallest index such that \( d_j = a_{k+1} \). Then, we can build the "rectangle" until \( w_{j-1} \) such that \( w_{j-1} \xrightarrow{\alpha_{j-1}+1} w_{j-1} \) and there is a sequence of transitions \( w_{k+1} = w_0 \xrightarrow{d_1} w_1 \xrightarrow{d_m} w_{j-1} \). By definition of \( \mathcal{E}_c, w_{k+1} \subseteq w_{j-1}^+ \) for all \( i < j - 1 \). If any of the transitions labeled by \( d_i \) is a self-loop, then we remove the state \( w_{j-1} \) from the sequence. Since \( d_j = a_{k+1} \) and there are no two different transitions with the same label, \( w_j = w_{j-1} \) and we take the existing sequence from \( w_j \). This completes the inductive step.

\[ \square \]

**Corollary A.10** Let \( w \in \mathcal{W} \). If there is an infinite sequence of transitions \( w = w_0 \xrightarrow{a_1} w_1 \xrightarrow{a_2} \cdots \), then there is an infinite sequence of non-removable transitions \( w = w_0 \xrightarrow{E} w_1 \xrightarrow{E} \cdots \xrightarrow{E} w_k \xrightarrow{E} w' \) in \( M_\mathcal{P} \) such that \( w' \sim w' \).

**Proof.** A consequence of the induction step in the proof of Lemma A.9 and the definition of \( \Rightarrow \). \( \square \)

Now, we are ready to prove that \( \sim \) satisfies \( P1 \).

**Lemma A.11** The relation \( \sim \) satisfies the property \( P1 \), i.e., if \( v \sim w, w \in \mathcal{W}' \subseteq \mathcal{W} \) and \( v \mapsto v' \) and \( \neg(\nu \sim v') \), then there is a non-empty sequence of transitions \( w = w_0 \xrightarrow{E} w_1 \xrightarrow{E} \cdots \xrightarrow{E} w_k \xrightarrow{E} w' \) in \( M_\mathcal{P} \) such that \( v' \sim w' \).

**Proof.** Let \( v \sim w, v \mapsto v' \) and \( \neg(\nu \sim v') \). From Lemma A.7 the relation \( \sim \) is a divergence blind stuttering equivalence. Then, by definition of divergence blind stuttering equivalence and transitivity of \( \sim \), there is a non-empty sequence of transitions \( w = u_0 \xrightarrow{E} u_1 \xrightarrow{E} \cdots \xrightarrow{E} u_n \xrightarrow{E} u' \) such that for all \( i \leq n, v \sim u_i \) and \( v' \sim u' \). It follows from transitivity of \( \sim \) that \( \neg u_n \sim u' \). Let \( b \) be the label such that \( u_n \xrightarrow{a} u' \). Then, by Lemma A.9, there is a non-empty sequence of non-removable transitions \( w = w_0 \xrightarrow{E} w_1 \xrightarrow{E} \cdots \xrightarrow{E} w_k \xrightarrow{E} w' \) and there is a (possibly empty) sequence of transitions \( u_n = w_0 \xrightarrow{d_1} w_1 \xrightarrow{d_m} w_m = w' \) such that \( w_n \subseteq w_{i+1}^+ \) and \( w_i \neq w_{i+1}^+ \) for \( i < m \). We show now that \( a \neq d_i \) for all \( i \leq m \). If \( a = d_1 \), then \( u' = w_1 \). Since \( a_n \sim w_1, u_n \sim u' \). Contradiction with the assumption that \( \neg (u_n \sim u') \). If \( a \neq d_1 \), then by Lemma A.5(a), \( a \in E(w_1^+) \). Let \( w_n \xrightarrow{a} w_{n+1}^+ \). Then, by definition of \( \mathcal{E}_w, w_n \subseteq w_{n+1}^+ \). By definition of \( \sim, \neg (w_n \sim w_{n+1}^+) \). Therefore, the same argument can be repeated to show that \( a \neq d_2 \) and then that \( a \neq d_i \) for each \( i > 2 \).

Therefore, by Lemma A.5(a), \( a \) is independent with all \( d_i \). So \( a \in E(w_1^+) \). Let \( w_n \xrightarrow{a} w_{n+1}^+ \) for \( i < m \) and denote \( u_{n+1}^+ \) by \( u \). Then, by definition of \( \mathcal{E}_c, w_{i+1} \subseteq w_{i+1}^+ \) for \( i < m \). Notice that \( u' = w_{n+1}^+ \) and \( u = w_{n+1}^+ \). Therefore, by definition of \( \sim, u' \sim u \).

If \( a \in E(w_1^+) \), then we are done. If not, then by Lemma A.8, there is a sequence of transitions \( w_i = v_0 \xrightarrow{b_1} v_1 \xrightarrow{b_2} \cdots \xrightarrow{b_k} v_k \) and the transition \( r_k \xrightarrow{a} w'' \) such that \( v_i \subseteq v_{i+1}^+, v_i \neq w_{i+1}^+, b_{i+1} \neq a \) for \( i < k \). and \( b_1 \ldots b_n \equiv b_1 \ldots b_k \). As before, using properties of \( \mathcal{E}_c \) and \( \sim \), we can show that \( u \sim w'' \). Therefore, by transitivity of \( \sim, u' \sim w'' \) and \( u' \sim w'' \), which ends the proof. \( \square \)

In order to prove that the relation \( \sim \) satisfies also the property \( P2 \), the following auxiliary lemma is used to show that if at some state an infinite sequence of stuttering transitions starts, then also an infinite sequence of stuttering transitions starts at each state related by \( \sim \).
Lemma A.12 Let \( w \sim v \). If there is an infinite sequence \( w \Rightarrow w_1 \Rightarrow w_2 \Rightarrow \cdots \Rightarrow w_i \Rightarrow \cdots \), then there is an infinite sequence \( v \Rightarrow v_1 \Rightarrow v_2 \Rightarrow \cdots \Rightarrow v_i \Rightarrow \cdots \).

Proof. Let \( \sim^1_c = \subseteq_c \cup \subseteq^T_c \) and \( \sim^{k+1}_c = \sim_k^c \cap \subseteq_c \). The proof is by induction on \( \text{length}(w \sim v) \), where \( \text{length}(w \sim v) = \min\{k > 0 | w \sim^k_c v\} \).

For the base case \( w \sim_c v \), if \( w = v \), then we are done. So, assume otherwise. Then, according to Lemma A.5(a), either \( v \sim w \) or \( w \sim v \).

In case \( v \sim w \), this fact together with \( w \sim v \) gives \( v \Rightarrow w \). Concatenation this edge to \( w \Rightarrow w_1 \Rightarrow w_2 \Rightarrow \cdots \Rightarrow w_i \Rightarrow \cdots \) provides the desired sequence.

To facilitate discussion of the second case where \( w \Rightarrow v \), we make some transition labels explicit. Let \( w \xrightarrow{a} v \) and \( w \xrightarrow{b_1} w_1 \xrightarrow{b_2} w_2 \xrightarrow{b_3} \cdots \xrightarrow{b_i} w_i \xrightarrow{b_{i+1}} \cdots \). Also assume \( w \) and \( v \) to have index 0 wherever applicable.

We build a sequence of transitions \( v_{i-1} \xrightarrow{c_i} v_i \) starting at \( v \), using the transitions \( w_i \xrightarrow{b_i} w \); from the sequence starting at \( w \).

As long as \( a \neq b_i \), proceed as follows: start from \( r \). If \( a \neq b_1 \) then, by Lemma A.5(a), \( (a, b_1) \in I \) and hence \( b_1 \in s\text{en}(v) \). So, for some \( v_1 \), \( v \xrightarrow{b_1} r_1 \). Note that if \( r_1 = r \) this is a self-loop and therefore, by Definition A.4(a), also \( v \sim v \) and thus \( v \Rightarrow v \); hence the self-loop yields the desired sequence and we are done in this case. If \( v_1 \neq v \), because \( (a, b_1) \in I \), also \( w_1 \xrightarrow{a} v_1 \). By the Definition A.4(c), then \( w_1 \sim v_1 \). Hence, by transitivity of \( \sim \), \( v \sim r_1 \). This establishes \( v \Rightarrow r_1 \).

As mentioned, \( v_1 \xrightarrow{a} r_1 \), thus making the situation at \( v_1 \) with respect to \( w_1 \) exactly the same as at \( v \) with respect to \( w \). This shows that, as long as \( a \neq b_i \), the above construction can be applied iteratively.

Now when \( a = b_j \) is encountered, this means that \( v_j = w_j \). But that means that from \( v_{j-1} \) on one can take as transitions \( v_{k-1} \xrightarrow{c_j} v_k \). \( k \geq j \), just the transitions \( w_{k-1} \xrightarrow{b_i} w_k \), yielding the desired sequence.

For the induction step let \( w \sim^{k-1}_c z \sim_c r \) for some \( z \). With respect to \( w \sim^{k-1}_c z \), the induction hypothesis provides a sequence \( z \xrightarrow{d_1} z_1 \xrightarrow{d_2} z_2 \xrightarrow{d_1} \cdots \xrightarrow{d_i} z_i \xrightarrow{d_{i+1}} \cdots \). In precisely the same manner as in the base case, a sequence \( v \Rightarrow v_1 \Rightarrow v_2 \Rightarrow \cdots \Rightarrow v_i \Rightarrow \cdots \) can now be constructed. This concludes the proof.

Lemma A.13 The relation \( \sim \) satisfies the property \( P2 \), i.e., if \( v \sim w \) and there is an infinite stutter path \( v = v_0 \Rightarrow v_1 \Rightarrow v_2 \Rightarrow \cdots \), then there is an infinite stutter path \( w = w_0 \Rightarrow w_1 \Rightarrow w_2 \Rightarrow \cdots \).

Proof. Let \( v \sim w \) and \( v \Rightarrow v_1 \Rightarrow v_2 \Rightarrow \cdots \). Then there is an infinite stutter sequence starting at \( v \). By Lemma A.12, there is an infinite sequence \( w \Rightarrow w_1 \Rightarrow w_2 \Rightarrow \cdots \) starting at \( w \). By applying Corollary A.11, we obtain an infinite stutter sequence \( w = w_0' \Rightarrow w_1' \Rightarrow w_2' \Rightarrow \cdots \).

Finally, we prove that the algorithm produces a reduced state graph that is branching bisimilar to the full state graph. i.e., Theorem 5.2 from Section 5.2:

Proof. From the proofs of Lemmas A.5(b), A.7, and A.11 one immediately concludes that \( \subseteq \) is a branching bisimulation between \( M_P \) and \( M_F \).
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