

Synthesis of Concurrent Systems with Many Similar Processes

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Methods for synthesizing concurrent programs from temporal logic specifications based on the use of a decision procedure for testing temporal satisfiability have been proposed by Emerson and Clarke and by Manna and Wolper. An important advantage of these synthesis methods is that they obviate the need to manually compose a program and manually construct a proof of its correctness. One only has to formulate a precise problem specification; the synthesis method then mechanically constructs a correct solution. A serious drawback of these methods in practice, however, is that they suffer from the state explosion problem. To synthesize a concurrent system consisting of K sequential processes, each having N states in its local transition diagram, requires construction of the global product-machine having about N^K global states in general. This exponential growth in K makes it infeasible to synthesize systems composed of more than 2 or 3 processes. In this article, we show how to synthesize concurrent systems consisting of many (i.e., a finite but arbitrarily large number K of) similar sequential processes. Our approach avoids construction of the global product-machine for K processes; instead, it constructs a two-process product-machine for a single pair of generic sequential processes. The method is uniform in K , providing a simple template that can be instantiated for each process to yield a solution for any fixed K . The method is also illustrated on synchronization problems from the literature.

Categories and Subject Descriptors: C.2.4 [**Computer-Communication Networks**]: Distributed Systems; D.1.2 [**Programming Techniques**]: Automatic Programming; D.1.3 [**Programming Techniques**]: Concurrent Programming; D.2.4 [**Software Engineering**]: Program Verification; F.3.1 [**Logics and Meanings of Programs**]: Specifying and Verifying and Reasoning about Programs—*mechanical verification*; I.2.2 [**Artificial Intelligence**]: Automatic Programming—*program synthesis*

A preliminary version of some of these results was presented at the ACM Symposium on Principles of Programming Languages, Austin, Texas, 1989, under the title “Synthesis of Concurrent Systems with Many Similar Sequential Processes.” P.C. Attie was sponsored by the Air Force Office of Scientific Research, Air Force Materiel Command, USAF, under grant number F49620-96-1-0221. E.A. Emerson was supported in part by NSF Grant CCR-9415496 and by SRC Contract 97-DP-388. The U.S. Government is authorized to reproduce and distribute reprints for Governmental purposes notwithstanding any copyright notation thereon. The views and conclusions contained herein are those of the authors and should not be interpreted as necessarily representing the official policies or endorsements, either expressed or implied, of the Air Force Office of Scientific Research or the U.S. Government.

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General Terms: Theory, Verification

Additional Key Words and Phrases: Concurrent programs, program synthesis, temporal logic

1. INTRODUCTION

We exhibit a method of synthesizing a system of K *similar sequential processes* executing in parallel from a temporal logic specification, where K is an arbitrarily large natural number. The method can be automated. It may be seen as an extension of the synthesis method of Emerson and Clarke [1982] and the related method of Manna and Wolper [1984]. The basic idea underlying both these earlier approaches to synthesis was to exploit the finite-model property for the temporal logic serving as a specification language: if a formula is satisfiable then it is satisfiable in a finite model. This makes it possible to develop a decision procedure for testing satisfiability of formulas in the specification logic. Given a formula, the decision procedure determines whether there exists a model of the formula. If such a model does exist, the procedure constructs a model which is a finite, labeled, directed graph that may be viewed as the global state transition diagram of a concurrent system meeting the specification. From this global state transition diagram, the *synchronization skeletons* of the individual sequential processes may be obtained by projecting out onto the coordinates corresponding to each of the individual sequential processes.

In principle, these earlier model-theoretic methods of program synthesis would have permitted synthesis of concurrent systems consisting of many (i.e., a finite but arbitrarily large number K of) similar sequential processes running in parallel. In practice, however, the methods were not feasible for systems with many processes because of the state explosion problem: the global state transition diagram is essentially the automata-theoretic product of the sequential processes (these processes being viewed as state-machines), and is therefore of size exponential in K , in general. Thus, even if the individual sequential processes had only two local states each, the global state transition diagram would contain on the order of 2^K global states. More recent synthesis methods [Anuchitanukul and Manna 1994; Pnueli and Rosner 1989a; Pnueli and Rosner 1989b] suffer from the same intractability problem.

In this article, we propose a method (which we refer to as MP-synthesis, i.e., many-process synthesis, in the sequel) to overcome the synthesis state explosion problem. We show how to reduce the problem of synthesizing a concurrent system of K similar sequential processes, each with N local states—which would naively require building the global state transition diagram of size on the order of N^K —to the problem of constructing the product of small numbers of sequential processes, and in particular, the product of a pair of sequential processes (which we call a *pair-system*), thereby avoiding the exponential complexity in K .

One other highly desirable feature of MP-synthesis is that it is uniform in K . Applied to a “generic” pair of representative processes i, j (since all processes are similar, any pair is representative, so we can select an arbitrary pair of *process indices* i, j , which we use as names for our representative pair) MP-synthesis yields a template for process i that can be instantiated for each K to get a synchronization

skeleton for process i in the K -process solution. The method is uniform so that the $(K + 1)$ -process solution is analogous to the K -process solution. Furthermore, our method can generate systems under arbitrary *process interconnection* schemes, e.g., fully connected, ring, star. In our model of parallel computation, two processes are interconnected if and only if either (1) one process can inspect the local state of the other process or (2) both processes read and/or write a common variable, or both.

The method requires the pair-systems to satisfy certain technical assumptions, which are given in the sequel. Thus it is not completely general. Nevertheless, it is applicable in many interesting cases. In particular, we illustrate MP-synthesis by applying it to the K -process mutual exclusion problem, the K -process dining philosophers problem, and the K -process drinking philosophers problem [Chandy and Misra 1984; Chandy and Misra 1988] for arbitrarily large finite K . This is a significant improvement over the previous literature. For example, Wolper [1982] considers the dining philosophers problem, but only for three philosophers; even five philosophers made the synthesis problem computationally intractable. Similarly, the solutions synthesized in Emerson and Clarke [1982] and Mamma and Wolper [1984] for the mutual exclusion problem were only for two processes. Consideration of just three processes made the problem infeasible for hand computation, while an actual implementation automating the Emerson and Clarke [1982] synthesis method [Inaba 1984] was unable to handle more than four processes. Thus, we are able to overcome the severe limitations previously imposed by the state explosion problem on the applicability of automatic temporal logic synthesis methods.

A crucial aspect of our method is its soundness: what correctness properties of the pair-systems are preserved by the K -process systems? We demonstrate, using the technical assumptions mentioned above, that (1) a propositional invariant of any pair-system is also an invariant of the K -process system and (2) a temporal leads-to property (expressed over a single process) satisfied by a pair-system is also satisfied by the K -process system. (A temporal leads-to property is of the following form: if condition 1 is true now, then condition 2 will eventually be true.) Part (1) means that safety properties which can be expressed pairwise are preserved. For example, mutual exclusion of K processes from a critical section can be expressed pairwise by stating that no two processes are in the critical section at the same time. Part (2) means that typical progress properties of a *single process* are preserved. For example, absence of starvation in the mutual exclusion problem can be expressed as “if a process requests entry to the critical section, then it eventually gains entry to the critical section.”

We do not consider other correctness properties, e.g., liveness properties expressed over a pair of processes. Such properties arise, for example, when expressing data dependencies in distributed or heterogeneous databases [Attie et al. 1996; Elmagarmid 1992; Klein 1991], e.g., “if transaction 1 commits then transaction 2 must also commit.” There is no reason in principle that MP-synthesis would not preserve such properties; however, different technical assumptions than the ones we make here would be required. We defer this topic for future work.

The rest of this article is organized as follows. Section 2 defines the model of computation, and Section 3 presents the specification language. Section 4 presents our syntax for concurrent programs and other notation. Section 5 describes MP-synthesis, and Section 6 gives a proof of its soundness. Section 7 presents a way of

implementing the programs we synthesize in terms of interprocess communication and synchronization primitives that are available in practice, namely asynchronous message passing and broadcast. Section 8 discusses some important aspects of the synthesis method: its complexity and scope of applicability. Finally, Section 9 discusses related work by others; Section 10 discusses possible extensions of the results presented here; and Section 11 presents our conclusions. Appendices B and C present those proofs that are omitted from the main text for reasons of space.

2. MODEL OF PARALLEL COMPUTATION

We consider concurrent programs of the form $P = P_1 \parallel \dots \parallel P_K$ that consist of a finite number of fixed sequential processes P_1, \dots, P_K running in parallel. With every process P_i , we associate a single, unique index, namely i . We then say that two processes are *similar* if and only if one can be obtained from the other by swapping their indices (this is formalized in Section 6.2). Intuitively, this corresponds to concurrent algorithms where a single “generic” indexed piece of code gives the code body for all processes.

We observe that for most actual concurrent programs the portions of each process responsible for interprocess synchronization can be cleanly separated from the sequential applications-oriented computations performed by the process. This suggests that we focus our attention on *synchronization skeletons* which are abstractions of actual concurrent programs where detail irrelevant to synchronization is suppressed.

We may view the synchronization skeleton of an individual process P_i as a state-machine where each state represents a region of code intended to perform some sequential computation and where each arc represents a conditional transition (between different regions of sequential code) used to enforce synchronization constraints. For example, there may be a node labeled C_i representing the critical section of process P_i . While in C_i , the process P_i may simply increment a single variable, or it may perform an extensive series of updates on a large database. In general, the internal structure and intended application of the regions of sequential code in an actual concurrent program are unspecified in the synchronization skeleton. By virtue of the abstraction to synchronization skeletons, we thus eliminate all steps of the sequential computation from consideration.

Formally, the synchronization skeleton of each process P_i is a directed graph where each node is labeled by a unique name (s_i) which represents a *local state* of P_i , and where each arc is labeled with a synchronization command $B \rightarrow A$ consisting of an enabling condition (i.e., guard) B and corresponding action A to be performed (i.e., a guarded command [Dijkstra 1976]). Self-loops, where there is an arc from a node to itself, are disallowed. Also, each node must have at least one outgoing arc, i.e., a skeleton contains no “dead ends.” A *global state* is a tuple of the form $(s_1, \dots, s_K, v_1, \dots, v_m)$ where each node s_i is the current local state of P_i and where v_1, \dots, v_m is a list giving the current values of the shared variables x_1, \dots, x_m (we assume that these are ordered in some fixed way, so that v_1, \dots, v_m specifies a unique value for each shared variable). A guard B is a predicate on states, and an action A is a parallel assignment statement that updates the values of the shared variables. If the guard B is omitted from a command, it is interpreted as *true*, and we simply write the command as A . If the action A is omitted, the shared variables

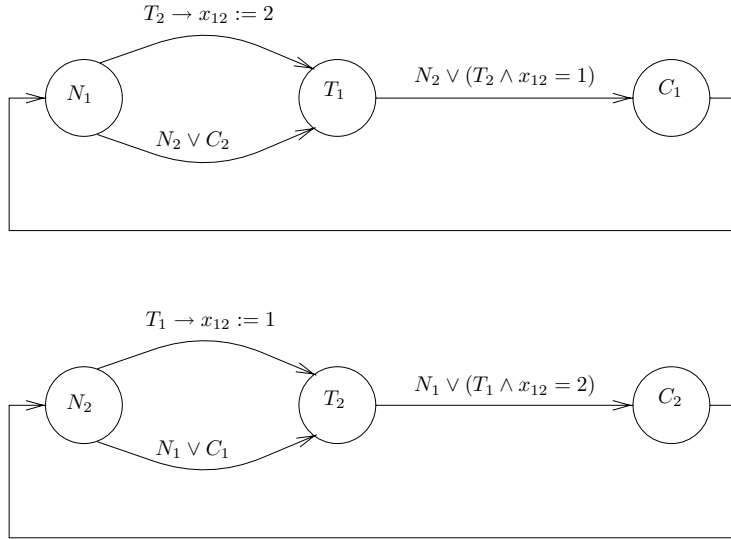


Fig. 1. Two-process mutual exclusion example.

are unaltered, and we write the command as B . For example, Figure 1 displays the synchronization skeletons (taken from Emerson and Clarke [1982]) synthesized by the synthesis method of Emerson and Clarke for the two-process mutual exclusion problem.

We model parallelism in the usual way by the nondeterministic interleaving of the “atomic” transitions of the individual synchronization skeletons of the processes P_i . Hence, at each step of the computation, some process with an “enabled” arc is nondeterministically selected to be executed next. Assume that the current state is $s = (s_1, \dots, s_i, \dots, s_K, v_1, \dots, v_m)$ and that process P_i contains an arc from node s_i to s'_i labeled by the command $B \rightarrow A$. If B is true in the current state then a permissible next state is $(s_1, \dots, s'_i, \dots, s_K, v'_1, \dots, v'_m)$ where v'_1, \dots, v'_m is the list of updated values for the shared variables resulting from action A . The arc from node s_i to s'_i is said to be *enabled* in state s . An arc that is not enabled is *disabled*, or *blocked*. A *computation path* is any sequence of states where each successive pair of states is related by the above next-state relation.

3. THE SPECIFICATION LANGUAGE MPCTL*

Our specification language is an extension of the temporal logic CTL* (Computation Tree Logic* [Emerson 1990]). CTL* is a propositional branching-time temporal logic. The basic modalities of CTL* consist of a path quantifier, either A (for all paths) or E (for some path) followed by a linear-time formula, which is built up from atomic propositions, the boolean operators \wedge, \vee, \neg , and the linear-time modalities G (always), F (sometime), X_j (strong nexttime), Y_j (weak nexttime) and U (until). CTL* formulae are built up from atomic propositions, the boolean operators \wedge, \vee, \neg , and the basic modalities. We remark that since synchronization

skeletons are in general finite state, the propositional version of temporal logic can be used to specify their properties.

3.1 CTL*

We have the following syntax for CTL*. We inductively define a class of state formulae (true or false of states) using rules (S1)–(S3) below and a class of path formulae (true or false of paths) using rules (P1)–(P3) below:

- (S1) Each atomic proposition p is a state formula;
- (S2) if f, g are state formulae, then so are $f \wedge g$, $\neg f$ (where \wedge, \neg indicate conjunction and negation, respectively); and
- (S3) if f is a path formula, then Ef and Af are state formulae.
- (P1) Each state formula is also a path formula;
- (P2) if f, g are path formulae, then so are $f \wedge g$, $\neg f$; and
- (P3) if f, g are path formulae, then so are $X_j f$, fUg .

Ef is a formula which intuitively means that there is some maximal path for which f holds, and Af is a formula which intuitively means that f holds of every maximal path. $X_j f$ (process indexed strong nexttime) is a formula which intuitively means that the immediate successor state along the maximal path under consideration is reached by executing one step of process P_j , and formula f holds in the immediate successor state.¹ fUg is a formula which intuitively means that there is some state along the maximal path under consideration where g holds, and f holds at every state along this path up to, but not necessarily including, that state.

CTL* consists of the set of state formulae generated by the above rules. The linear-time temporal logic PTL [Manna and Wolper 1984] consists of the set of path formulae generated by rules (S1) and (P1)–(P3). The logic CTL [Emerson 1990; Emerson and Clarke 1982] is obtained by replacing rules (P1)–(P3) by (P0):

- (P0) If f, g are state formulae, then $X_j f$, fUg are path formulae.

The set of state formulae generated by rules (S1)–(S3) and (P0) forms CTL.

Formally, we define the semantics of CTL* formulae with respect to a (K -process) structure $M = (S, R_{i_1}, \dots, R_{i_K})$ consisting of

- S , a countable set of states. Each state is a mapping from the set of atomic propositions into $\{true, false\}$ and
- $R_i \subseteq S \times S$, a binary relation on S giving the transitions of sequential process i .

Let $R = R_{i_1} \cup \dots \cup R_{i_K}$. A *path* is a sequence of states (s_1, s_2, \dots) such that $\forall i, (s_i, s_{i+1}) \in R$, and a *fullpath* is a maximal path. A fullpath (s_1, s_2, \dots) is infinite unless for some s_k there is no s_{k+1} such that $(s_k, s_{k+1}) \in R$. We use the convention (1) that $\pi = (s_1, s_2, \dots)$ denotes a fullpath and (2) that π^i denotes the suffix $(s_i, s_{i+1}, s_{i+2}, \dots)$ of π , provided $i \leq |\pi|$, where $|\pi|$, the length of π , is ω when π is infinite and k when π is finite and of the form (s_1, \dots, s_k) ; otherwise

¹Note that we follow the presentation in Emerson and Clarke [1982], which introduced the process indexed nexttime operator, rather than the more conventional presentation of Emerson [1990], where the nexttime operator is unindexed.

π^i is undefined. We also use the usual notation to indicate truth in a structure: $M, s_1 \models f$ (respectively $M, \pi \models f$) means that f is true in structure M at state s_1 (respectively of fullpath π). In addition, we use $M, S \models f$ to mean $\forall s \in S. (M, s \models f)$, where S is a set of states. We define \models inductively:

- (S1) $M, s_1 \models p$ iff $s_1(p) = \text{true}$
- (S2) $M, s_1 \models f \wedge g$ iff $M, s_1 \models f$ and $M, s_1 \models g$
 $M, s_1 \models \neg f$ iff $\text{not}(M, s_1 \models f)$
- (S3) $M, s_1 \models Ef$ iff there exists a fullpath $\pi = (s_1, s_2, \dots)$ in M such that
 $M, \pi \models f$
 $M, s_1 \models Af$ iff for every fullpath $\pi = (s_1, s_2, \dots)$ in M : $M, \pi \models f$
- (P1) $M, \pi \models f$ iff $M, s_1 \models f$
- (P2) $M, \pi \models f \wedge g$ iff $M, \pi \models f$ and $M, \pi \models g$
 $M, \pi \models \neg f$ iff $\text{not}(M, \pi \models f)$
- (P3) $M, \pi \models X_j f$ iff π^2 is defined and $(s_1, s_2) \in R_j$ and $M, \pi^2 \models f$
 $M, \pi \models fUg$ iff there exists $i \in [1 : |\pi|]$ such that $M, \pi^i \models g$ and for
 all $j \in [1 : (i - 1)]$: $M, \pi^j \models f$

We write $\models f$ to indicate that f is *valid*, i.e., true at all states in all structures. When the structure M does not affect the truth of the formula, (e.g., $M, s_1 \models p$ depends only on s_1 and p), it may be omitted (e.g., $M, s_1 \models p$ is written as $s_1 \models p$).

We introduce the abbreviations $f \vee g$ for $\neg(\neg f \wedge \neg g)$, $f \Rightarrow g$ for $\neg f \vee g$, and $f \equiv g$ for $(f \Rightarrow g) \wedge (g \Rightarrow f)$, which denote logical disjunction, implication, and equivalence, respectively. We also introduce a number of additional modalities as abbreviations: $Y_j f$ for $\neg X_j \neg f$, Ff for $[trueUf]$, Gf for $\neg F\neg f$, $fU_w g$ for $[fUg] \vee Gf$, $\overset{\infty}{F}f$ for GFf , and $\overset{\infty}{G}f$ for FGf . Y_j is the “process indexed weak nexttime” modality. $Y_j f$ is a formula which intuitively means that if the immediate successor state along the maximal path exists, and is reached by executing one step of process P_j , then formula f holds in the immediate successor state. F is the “eventually” modality. Ff is a formula which intuitively means that there is some state along the maximal path where f holds. G is the “always” modality. Gf is a formula which intuitively means that f holds at every state along the maximal path. U_w is the “weak until” (or “unless”) modality. $fU_w g$ is a formula which intuitively means that either g holds at some state along the maximal path, and f holds at every state along the maximal path up to (but not necessarily including) that state; or g never holds along the maximal path, and f holds at all states of the maximal path. Finally, $\overset{\infty}{F}$ and $\overset{\infty}{G}$ are the infinitary modalities “infinitely often” and “almost everywhere,” respectively. $\overset{\infty}{F}f$ and $\overset{\infty}{G}f$ are formulae which intuitively mean respectively that f holds in infinitely many states along the maximal path and that f holds in all but a finite number of states along the maximal path. Particularly useful modalities are AFf , which means that for every maximal path there exists a state on the path where f holds, and AGf , which means that f holds at every state along every path. Note that the unindexed nexttime operators can be viewed as abbreviations:

$$Xf \equiv (X_{i_1}f \vee \dots \vee X_{i_K}f)$$

$$Yf \equiv (Y_{i_1}f \wedge \dots \wedge Y_{i_K}f)$$

3.2 Sublogics of CTL*

In the sequel, we define various sublogics of CTL* and PTL. We denote a sublogic by \mathcal{L} followed by a list (enclosed in parentheses) of all the

- sets of atomic propositions,
- boolean operators, and
- temporal modalities

that can be used in constructing formulae of the sublogic. For example, $\mathcal{L}(\{Q_1, R_1\}, \{Q_2, R_2\}, \neg, \wedge, U)$ denotes the sublogic of PTL consisting of all those formulae which contain only the atomic propositions Q_1, Q_2, R_1, R_2 and which contain no next-time (X, Y) modalities, while $\mathcal{L}(\wedge, \vee, AG(q_i \Rightarrow AFR_i), AG(q_i \Rightarrow AXr_i))$ is the sublogic of CTL* consisting of positive boolean combinations of formulae of the form $AG(q_i \Rightarrow AFR_i)$ or $AG(q_i \Rightarrow AXr_i)$.

3.3 The Interconnection Relation

As stated in the introduction, we present a method of synthesizing a K -process system from a pair-system. We also stated that our method can deal with arbitrary process interconnection schemes. The desired process interconnection scheme is given by the *interconnection relation* I , which is, therefore, part of the problem specification. The domain of I is the set $\{i_1, \dots, i_K\}$ of indices of the processes in the K -process system being specified. Thus, $I \subseteq \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\}$, and $i I j$ iff processes i and j are interconnected. In the sequel, we say that i and j are *neighbors* when $i I j$. Since our definition of interconnection (Section 1) is symmetric, we stipulate that I is symmetric. We also stipulate that I is irreflexive, since a process cannot be interconnected to itself. Furthermore, we assume that I is a total relation: $\bigwedge i \in \{i_1, \dots, i_K\}. (\bigvee j. (i I j))$. Thus every process is interconnected to at least one other process.

Since there are many possible interconnection schemes for a given number K of processes, we henceforth use the more precise term *I-system* in place of the (up till now used) term “ K -process system.” We also introduce the following abbreviations. $dom(I)$ denotes the domain of I (i.e., the set $\{i_1, \dots, i_K\}$); $I(i)$ denotes the set $\{j \mid i I j\}$; and $\hat{I}(i)$ denotes the set $\{i\} \cup \{j \mid i I j\}$.

3.4 MPCTL*

An MPCTL* (Many-Process CTL*) formula consists of a spatial modality followed by a CTL* state formula over a “uniformly” indexed family $\mathcal{AP} = \{\mathcal{AP}_{i_1}, \dots, \mathcal{AP}_{i_K}\}$ of atomic propositions. The propositions in \mathcal{AP}_i are the same as those in \mathcal{AP}_j except for their subscripts; for example, we could have $\mathcal{AP}_1 = \{Q_1, R_1\}$ and $\mathcal{AP}_2 = \{Q_2, R_2\}$. Q_1 corresponds to Q_2 , R_1 to R_2 , etc. A spatial modality is of the form \bigwedge_i , or \bigwedge_{ij} . \bigwedge_i quantifies the process index i which ranges over $\{i_1, \dots, i_K\}$. \bigwedge_{ij} quantifies the process indices i, j which range over the elements of $\{i_1, \dots, i_K\}$ which are related by I . If the spatial modality is \bigwedge_i , then only atomic propositions in \mathcal{AP}_i are allowed in the CTL* formula, and if the spatial modality is \bigwedge_{ij} , then only atomic propositions in $\mathcal{AP}_i \cup \mathcal{AP}_j$ are allowed in the CTL* formula. In other words, every process index within an MPCTL* formula must be “bound” by the spatial modality.

Thus mutual exclusion over all pairs of processes that are related by I is expressed by $\bigwedge_{ij} AG(\neg(C_i \wedge C_j))$, (no two processes related by I are in their critical regions simultaneously), while absence of starvation for each process is expressed by $\bigwedge_i AG(T_i \Rightarrow AFC_i)$ (every process which enters its trying region will eventually enter its critical region). The semantics of these assertions over a structure depends on the interconnection relation I associated with the structure. For example, the semantics of the mutual exclusion specification over a structure is given by viewing the spatial modalities as abbreviations and expanding them in the obvious way: $AG(\neg(C_1 \wedge C_2))$ for a two-process structure, with $I = \{\{1, 2\}\}$, and $AG(\neg(C_1 \wedge C_2)) \wedge AG(\neg(C_1 \wedge C_3)) \wedge AG(\neg(C_2 \wedge C_3))$ for a three-process structure, with $I = \{\{1, 2\}, \{1, 3\}, \{2, 3\}\}$, and so forth. Similarly, the semantics of the absence of starvation assertion is given by $AG(T_1 \Rightarrow AFC_1) \wedge AG(T_2 \Rightarrow AFC_2)$ over a two-process structure with $I = \{\{1, 2\}\}$, and by $AG(T_1 \Rightarrow AFC_1) \wedge AG(T_2 \Rightarrow AFC_2) \wedge AG(T_3 \Rightarrow AFC_3)$ over a three-process structure with $I = \{\{1, 2\}, \{1, 3\}, \{2, 3\}\}$.

In general, an MPCTL* formula q is interpreted over a K -process structure $M = (S, R_{i_1}, \dots, R_{i_K})$ where S and R_{i_1}, \dots, R_{i_K} are as before. We say that K is the *arity* of the structure M . The definition of truth in structure M at state s of formula q is given by $M, s \models q$ iff $M, s \models q'$ where q' is the CTL* formula obtained from q by viewing q as an abbreviation and expanding it as shown above, i.e.,

$$-M, s \models \bigwedge_i f_i \text{ iff } M, s \models \bigwedge i \in \{i_1, \dots, i_K\}. (f_i)$$

$$-M, s \models \bigwedge_{ij} f_{ij} \text{ iff } M, s \models \bigwedge (i, j) \in I. (f_{ij})$$

where f_i is a CTL* formula containing atomic propositions from \mathcal{AP}_i only, and f_{ij} is a CTL* formula containing atomic propositions from $\mathcal{AP}_i \cup \mathcal{AP}_j$ only. Note that we use $\bigwedge v \in R. (f)$, $\bigvee v \in R. (f)$, to denote $f(v_1) \wedge \dots \wedge f(v_n)$, $f(v_1) \vee \dots \vee f(v_n)$ respectively, where $R = \{v_1, \dots, v_n\}$. Sometimes we will also use $\bigwedge_{v \in R} f$, $\bigvee_{v \in R} f$. We also use “ $\bigwedge (i, j) \in I$ ” to denote “ $\bigwedge i, j \in \text{dom}(I)$ such that $i I j$.” In MPCTL*, $\bigwedge_{ij} f_{ij}$ is permitted, and is equivalent to $\bigwedge_i f_i$. MPCTL* is similar to the Indexed CTL* of Clarke et al. [1986], but allows quantification over pairs of processes (via the quantifier \bigwedge_{ij}) as well as over single processes. Furthermore, if the CTL* subformula of the MPCTL* formula is restricted to be in CTL [Emerson 1990; Emerson and Clarke 1982], then the resulting logic is called MPCTL.

An example of an MPCTL* specification is given below for the mutual exclusion problem. In this specification, process P_i is always in exactly one of three states N_i , T_i , C_i , which represent, respectively, the Noncritical, Trying, and Critical code regions. A process in its Noncritical state does not require the use of critical (shared) resources, and it performs computations that can proceed in parallel with computations by the other processes. A process requiring a critical resource moves into the Trying state. From there, it enters the Critical state as soon as possible, provided that all constraints on the use of the critical resource are met. In the Critical state, P_i has access to the critical resource. This specification is given by the following set of MPCTL* formulae (which are implicitly conjoined):

- (1) initial state (every process is initially in its noncritical region):

$$\bigwedge_i N_i$$

- (2) it is always the case that any move P_i makes from its noncritical region is into its trying region, and such a move is always possible:

$$\mathbf{\bigwedge}_i AG(N_i \Rightarrow (AY_i T_i \wedge EX_i T_i))$$

- (3) it is always the case that any move P_i makes from its trying region is into its critical region:

$$\mathbf{\bigwedge}_i AG(T_i \Rightarrow (AY_i C_i))$$

- (4) it is always the case that any move P_i makes from its critical region is into its noncritical region, and such a move is always possible:

$$\mathbf{\bigwedge}_i AG(C_i \Rightarrow (AY_i N_i \wedge EX_i N_i))$$

- (5) P_i is always in exactly one of the states N_i , T_i , or C_i :

$$\mathbf{\bigwedge}_i AG(N_i \equiv \neg(T_i \vee C_i))$$

$$\mathbf{\bigwedge}_i AG(T_i \equiv \neg(N_i \vee C_i))$$

$$\mathbf{\bigwedge}_i AG(C_i \equiv \neg(N_i \vee T_i))$$

- (6) P_i does not starve:

$$\mathbf{\bigwedge}_i AG(T_i \Rightarrow AFC_i)$$

- (7) a transition by one process cannot cause a transition by another:

$$\mathbf{\bigwedge}_{ij} AG((N_i \Rightarrow AY_j N_i) \wedge (N_j \Rightarrow AY_i N_j))$$

$$\mathbf{\bigwedge}_{ij} AG((T_i \Rightarrow AY_j T_i) \wedge (T_j \Rightarrow AY_i T_j))$$

$$\mathbf{\bigwedge}_{ij} AG((C_i \Rightarrow AY_j C_i) \wedge (C_j \Rightarrow AY_i C_j))$$

- (8) no two processes access critical resources together:

$$\mathbf{\bigwedge}_{ij} AG(\neg(C_i \wedge C_j))$$

If $I = \text{dom}(I) \times \text{dom}(I) - \{(i, i) \mid i \in \text{dom}(I)\}$, i.e., all pairs of processes are I -related, then the specification is for the K -process mutual exclusion problem, where at most one process is in its critical section at any time. On the other hand, by using an appropriate I , we can also specify the following problems in MPCTL*:

Five dining philosophers arranged in a ring:

$$I = \{\{1, 2\}, \{2, 3\}, \{3, 4\}, \{4, 5\}, \{5, 1\}\}.$$

One thousand dining philosophers arranged in a ring:

$$I = \{\{1, 2\}, \{2, 3\}, \dots, \{999, 1000\}, \{1000, 1\}\}.$$

Readers-Writers (with P_1, P_2 as readers, P_3, P_4 as writers, and no writer priority):

$$I = \{\{1, 3\}, \{1, 4\}, \{2, 3\}, \{2, 4\}, \{3, 4\}\}.$$

These examples generalize the mutual exclusion example: in dining philosophers, only neighboring (I -related) philosophers are mutually excluded from simultaneous access to their critical sections, and in readers-writers, only reader/writer and writer/writer pairs are mutually excluded.

4. NOTATION AND SYNTAX

In the sequel, unless otherwise indicated (e.g., by explicit quantification) i and j will denote arbitrary elements of $\text{dom}(I)$ such that $i \neq j$. Consider the pair-system consisting of processes i and j , which execute concurrently and interact by reading each other's local state and by reading/writing shared variables (when we need to specify the component processes, we call this pair-system the ij -system). We call (the synchronization skeletons for)² processes i, j in this pair-system P_i^j, P_j^i , respectively. In the sequel, we call a process of a pair-system a *pair-process*. As elaborated in Section 2, (a synchronization skeleton for) a process is a directed graph, where the nodes are local states of the process, and the arcs between nodes are labeled with synchronization commands, whose execution is required for the arc to be “traversed” by the process. These synchronization commands test the local state of other processes, and they test/modify shared variables. Let \mathcal{SH}_{ij} be the set of all the variables shared by P_i^j and P_j^i . We restrict P_i^j, P_j^i to the following syntax. Each shared variable x_{ij} assumes a value in $\{i, j\}$, i.e., all shared variables are binary. Shared variables may optionally be superscripted, which takes the place of the usual “subscript” notation, since we use subscripts here to denote the relevant pair-system. The guards of the synchronization commands that label arcs in P_i^j are built up from the standard propositional logic connectives, the atomic propositions in \mathcal{AP}_j ,³ and expressions of the form $x_{ij} = i$ or of the form $x_{ij} = j$, where x_{ij} is in \mathcal{SH}_{ij} . The actions of the synchronization commands that label arcs in P_i^j are multiple assignments of the form $x_{ij}, y_{ij}, z_{ij}, \dots := h, k, \ell, \dots$ where the left-hand side is a list of variables in \mathcal{SH}_{ij} , and the right-hand side is an equally long list of constants, all of whom have value either i or j . Each node of P_i^j assigns either *true* or *false* to each atomic proposition in \mathcal{AP}_i (i.e., is a mapping of \mathcal{AP}_i into $\{\text{true}, \text{false}\}$). We refer to the nodes of P_i^j as *i-states* (i.e., local states of P_i^j). For example, in Figure 1, the lower N_1 -to- T_1 transition has a label with guard $N_2 \vee C_2$ and assignment *skip*, and the upper N_1 -to- T_1 transition has a label with guard T_2 and assignment $x_{12} := 2$. The T_1 -to- C_1 transition has a label whose guard is $N_2 \vee (T_2 \wedge x_{12} = 1)$ and whose assignment is *skip*. The node N_1 maps N_1 to *true* and both T_1 and C_1 to *false*. Likewise, the node T_1 maps T_1 to *true* and both N_1 and C_1 to *false*, and the node C_1 maps C_1 to *true* and both N_1 and T_1 to *false*. In the sequel, we use S_i to denote the set of i -states in P_i^j , and we use $(s_i, B \rightarrow A, t_i)$ to denote an arc from i -state s_i to i -state t_i with label $B \rightarrow A$.

Since we adopt an interleaving model of parallel computation, we need to define the global state of the pair-system $P_i^j \parallel P_j^i$. A global state of this pair-system is a tuple $(s_i, s_j, v_{ij}^1, \dots, v_{ij}^m)$ where s_i, s_j are the current i -state, j -state, respectively, and $v_{ij}^1, \dots, v_{ij}^m$ give the values of all the variables in \mathcal{SH}_{ij} . We refer to global states of $P_i^j \parallel P_j^i$ as *ij-states*. When viewed as a mapping of $\mathcal{AP}_i \cup \mathcal{AP}_j$ into

²In our framework, there is no technical distinction between a process and a synchronization skeleton; a process is a synchronization skeleton. We therefore, for the most part, use the term “process” from now on instead of “skeleton,” but will sometimes use the term skeleton when we wish to emphasize the syntactic aspects of a process.

³We remind the reader that the atomic propositions in \mathcal{AP}_j can only be written by P_j^i , but may be read by P_i^j ; see Section 3.4.

$\{true, false\}$ (as required by Section 3.1 to evaluate CTL* formulae), an ij -state inherits the assignments defined by its component i - and j -states: $s_{ij}(Q_i) = s_i(Q_i)$, $s_{ij}(Q_j) = s_j(Q_j)$, where $s_{ij} = (s_i, s_j, v_{ij}^1, \dots, v_{ij}^m)$, and Q_i, Q_j are arbitrary atomic propositions in $\mathcal{AP}_i, \mathcal{AP}_j$, respectively. In addition to the skeletons for P_i^j, P_j^i , a complete description of the pair-system $P_i^j \| P_j^i$ requires that the initial states be given. We therefore include a set S_{ij}^0 of ij -states which gives the permissible initial states, and our complete notation for the pair-system consisting of processes i and j is $(S_{ij}^0, P_i^j \| P_j^i)$.

We now turn to I -systems. For a given interconnection relation I , the I -system consists of processes i_1, \dots, i_K executing concurrently, where $\{i_1, \dots, i_K\}$ is the domain of I . We call the synchronization skeleton for process i ($i \in \{i_1, \dots, i_K\}$) in the I -system P_i^I , and we call a process of the I -system an I -process. As in the case for pair-systems, the nodes of such processes are i -states, i.e., mappings of \mathcal{AP}_i into $\{true, false\}$. A global state of the I -system is a tuple $(s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, where s_i , ($i \in \{i_1, \dots, i_K\}$) is the current local state of P_i^I , and v^1, \dots, v^n give the values of all the shared variables of the I -system (once again we assume some fixed ordering of these variables, so that the values assigned to them are uniquely determined by the list v^1, \dots, v^n). We refer to global states of the I -system as **I -states**. When viewed as a mapping of $\mathcal{AP}_{i_1} \cup \dots \cup \mathcal{AP}_{i_K}$ into $\{true, false\}$ (as required by Section 3.1 to evaluate CTL* formulae), an I -state inherits the assignments defined by its component i -states ($i \in \{i_1, \dots, i_K\}$): $s_{ij}(Q_i) = s_i(Q_i)$, where $s = (s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, and Q_i is an arbitrary atomic proposition in \mathcal{AP}_i ($i \in \{i_1, \dots, i_K\}$). We usually use s_I, t_I, u_I to denote I -states. When the interconnection relation I is clear from context, it may be omitted, e.g., s_I, P_i^I are rendered as s, P_i , respectively. We use S_I^0 to denote the set of permissible initial I -states and \mathcal{SH} to denote the set of shared variables of the I -system. The I -system overall is denoted as $(S_I^0, P_{i_1}^I \| \dots \| P_{i_K}^I)$. The next section shows how $(S_I^0, P_{i_1}^I \| \dots \| P_{i_K}^I)$ is (uniquely) determined from $(S_{ij}^0, P_i^j \| P_j^i)$ and I .

Let s_i be an i -state. We define a state-to-formula operator $\{s_i\}$ that takes an i -state s_i as an argument and returns a propositional formula that characterizes s_i in that $s_i \models \{s_i\}$, and $s'_i \not\models \{s_i\}$ for all i -states s'_i such that $s'_i \neq s_i$: $\{s_i\} = (\bigwedge_{s_i(Q_i)=true} Q_i) \wedge (\bigwedge_{s_i(Q_i)=false} \neg Q_i)$, where Q_i ranges over the members of \mathcal{AP}_i .

5. DESCRIPTION OF THE METHOD

5.1 Synthesis of Many-Process Systems from Pair-Systems

As stated earlier, our aim is to synthesize an I -system without explicitly generating its global state transition diagram (and thereby incurring a time and space complexity exponential in $|dom(I)|$). We achieve this by breaking the problem down into two steps:

- (1) Synthesize a pair-system (from the given temporal logic specification)
- (2) Generalize the pair-system to an I -system

Since our focus in this article is on exploiting the similarity of processes, we shall not address the first step explicitly. We note in passing that any method for the derivation of concurrent programs can be used to generate the pair-system. In particular, if the problem specification can be written in MPCTL, then we can extract

a two-process specification (say for processes i and j) by expanding the MPCTL specification (as shown in Section 3) for $I = \{\{i, j\}\}$. We then apply the synthesis method of Emerson and Clarke [1982] to the two-process specification, resulting in the pair-system. For example, the specification for K -process mutual exclusion given in Section 3.4 is actually in MPCTL, and so the pair-system for mutual exclusion (Figure 1) can be generated automatically from this specification. Since a pair-system has only $O(N^2)$ states (where N is the size of each sequential process), the problem of deriving a pair-system from a specification is considerably easier than that of deriving an I -system from the specification. Hence, the contribution of this article, namely the second step above, is to reduce the more difficult problem (deriving the I -system) to the easier problem (deriving the pair-system). We proceed as follows.

Consider the pair-system $(S_{ij}^0, P_i^j \| P_j^i)$. We take this pair-system and generalize it in a natural way to an I -system. We show that our generalization preserves a large class of correctness properties. Roughly the idea is as follows. Consider first the generalization to three pairwise interconnected processes i, j, k , i.e., $I = \{\{i, j\}, \{j, k\}, \{k, i\}\}$. With respect to process i , the proper interaction (i.e., the interaction required to satisfy the specification) between process i and process j is captured by the synchronization commands that label the arcs of P_i^j . Likewise, the proper interaction between process i and process k is captured by the arc labels of P_i^k . Therefore, in the three-process system consisting of processes i, j, k executing concurrently, (and where process i is interconnected to both process j and process k), the proper interaction for process i with processes j and k is captured as follows: when process i traverses an arc, the synchronization command which labels that arc in P_i^j is executed “simultaneously” with the synchronization command which labels the corresponding arc in P_i^k . For example, taking as our specification the mutual exclusion problem, if P_i executes the mutual exclusion protocol with respect to both P_j and P_k , then, when P_i enters its critical section, both P_j and P_k must be outside their own critical sections.

Based on the above reasoning, we determine that the synchronization skeleton for process i in the aforementioned three-process system (call it P_i^{jk}) has the same basic graph structure as P_i^j and P_i^k , and a arc label in P_i^{jk} is a “composition” of the labels of the corresponding arcs in P_i^j and P_i^k . In addition, the initial states S_{ijk}^0 of the three-process system are exactly those states that “project” onto initial states of all three pair-systems $((S_{ij}^0, P_i^j \| P_j^i), (S_{ik}^0, P_i^k \| P_k^i), \text{ and } (S_{jk}^0, P_j^k \| P_k^j))$. More formally, $s_{ijk}^0 \in S_{ijk}^0$ if and only if $s_{ijk}^0 = (s_i, s_j, s_k, v_{ij}^1, \dots, v_{ij}^m, v_{ik}^1, \dots, v_{ik}^m, v_{jk}^1, \dots, v_{jk}^m)$, where $(s_i, s_j, v_{ij}^1, \dots, v_{ij}^m) \in S_{ij}^0$, $(s_i, s_k, v_{ik}^1, \dots, v_{ik}^m) \in S_{ik}^0$, and $(s_j, s_k, v_{jk}^1, \dots, v_{jk}^m) \in S_{jk}^0$.

Generalizing this to the case of an arbitrary interconnection relation I , we see that the skeleton for process i in the I -system, P_i^I , has the same basic graph structure as P_i^j , and a transition label in P_i^I is a “composition” of the labels of the corresponding transitions in $P_i^{j_1}, \dots, P_i^{j_n}$, where $\{j_1, \dots, j_n\} = I(i)$, i.e., processes j_1, \dots, j_n are all the I -neighbors of process i . Likewise the set S_I^0 of initial states of the I -system is exactly those states all of whose “projections” onto all the pairs in I give initial states of the corresponding pair-system.

This discussion leads to the following definition of the synthesis method, which shows how an I -process P_i^I of the I -system $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ is derived from the pair-process P_i^j of the pair-system $(S_{ij}^0, P_i^j \parallel P_j^i)$ and how the initial state set S_I^0 of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ is derived from the initial state set S_{ij}^0 of $(S_{ij}^0, P_i^j \parallel P_j^i)$.

Definition 5.1.1 (MP-Synthesis). An I -process P_i^I is derived from the pair-process P_i^j as follows:

P_i^I contains an arc from s_i to t_i with label $\bigwedge_{j \in I(i)} B_i^j \rightarrow //_{j \in I(i)} A_i^j$
iff
for every j in $I(i)$: P_i^j contains an arc from s_i to t_i with label $B_i^j \rightarrow A_i^j$.

The *initial state set* S_I^0 of the I -system is derived from the initial state set S_{ij}^0 of the pair-system as follows:

$$S_I^0 = \{(s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n) \mid \bigwedge (i, j) \in I. (s_i, s_j, v_{ij}^1, \dots, v_{ij}^m) \in S_{ij}^0\}$$

where $v_{ij}^1, \dots, v_{ij}^m$ are those values from v^1, \dots, v^n that denote values of variables in \mathcal{SH}_{ij} .

Here \bigwedge is logical conjunction, and $//$ is simultaneous parallel assignment, i.e., $//_{j \in I(i)} A_i^j$ is a single parallel assignment. Its execution is achieved by simultaneously executing every A_i^j (for fixed i , and as j varies over $I(i)$). From Definition 5.1.1, we see that the shared variables of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ are shared *pairwise*, i.e., \mathcal{SH} is partitioned into the sets \mathcal{SH}_{ij} for all I -related pairs i, j : $\mathcal{SH} = \cup (i, j) \in I. \mathcal{SH}_{ij}$ and $\bigwedge i, j, k \in \text{dom}(I). (\mathcal{SH}_{ij} \cap \mathcal{SH}_{ik} = \emptyset)$.

The MP-synthesis definition constructs the I -system as a “conjunctive composition” of the pair-systems for every pair in I . This composition operates as follows:

- For an I -process P_i^I to traverse an arc from s_i to t_i , every ij -system $(S_{ij}^0, P_i^j \parallel P_j^i)$ (for every j in $I(i)$) must be in an ij -state s_{ij} whose P_i^j -component is s_i , and each such s_{ij} must have an outgoing transition to some ij -state t_{ij} whose P_i^j -component is t_i
- When the P_i^I -transition from s_i to t_i occurs, the current ij -states of all the ij -systems are *simultaneously* changed from s_{ij} to t_{ij}

This characterization of transitions in the I -system as compositions of transitions in all the relevant pair-systems is formalized in Lemma 6.4.1. The above definition of MP-synthesis is, in effect, a *syntactic transformation* that can be carried out in linear time and space (in both $(S_{ij}^0, P_i^j \parallel P_j^i)$ and I). In particular, we avoid explicitly constructing the global state transition diagram of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$, which is of size exponential in $K = |\text{dom}(I)|$. Figure 2 gives an example application of MP-synthesis: the synthesis of a solution to the three-process mutual exclusion problem from the solution to the two-process mutual exclusion problem given in Figure 1. One of the transitions of the three-process mutual exclusion system is $[N_1 T_2 C_3] \xrightarrow{1} [T_1 x_{12} = 2 T_2 C_3]$ (it arises from the execution of the arc labeled $T_2 \wedge (N_3 \vee C_3) \rightarrow x_{12} := 2$ leading from 1-state N_1 to 1-state T_1). This transition can be regarded as a composition of the transitions $[N_1 T_2] \xrightarrow{1} [T_1 x_{12} = 2 T_2]$ and $[N_1 C_3] \xrightarrow{1} [T_1 C_3]$. The first is a transition of the two-process mutual exclusion

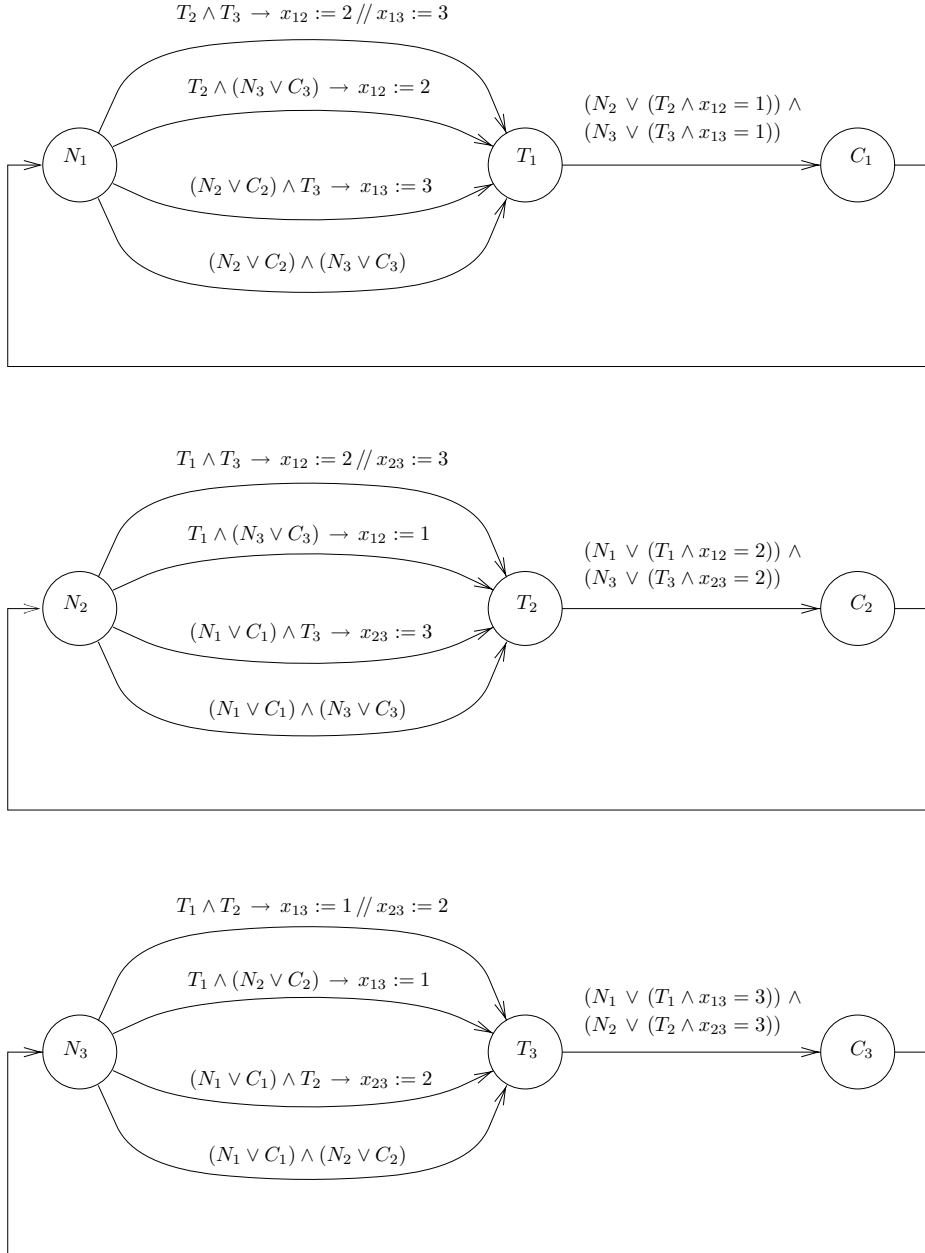


Fig. 2. Three-process mutual exclusion example.

system for processes 1 and 2, and the second is a transition of the two-process mutual exclusion system for processes 1 and 3.

5.1.1 Subsystems of the I -System. Let $J \subseteq I$, i.e., J is an interconnection relation which is a subrelation of I . Thus, J represents a subsystem (called the J -system) of the I -system in that it generally contains a subset of I 's processes, and these processes are related by a subset of the pair-constraints in I . The J -system is, in effect, the composition of the pair-systems for all the pairs in J . A global state of the J -system, called a J -state, is a tuple $(s_{j_1}, \dots, s_{j_L}, v_j^1, \dots, v_j^m)$, where $\{j_1, \dots, j_L\}$ is the domain of J , and v_j^1, \dots, v_j^m is a list of values (in some fixed ordering) for the variables in $\bigcup_{(i,j) \in J} \mathcal{SH}_{ij}$.

5.1.2 State and Path Projection. Due to the compositional nature of the I -system outlined above, it will be very convenient in the sequel to be able to refer to the components of a given I -state or ij -state. For this purpose, we define the **state projection operator** \uparrow . This operator has several variants. First of all, we define projection onto a single process from both I -states and ij -states: if $s = (s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, then $s \uparrow i = s_i$, and if $s_{ij} = (s_i, s_j, v_{ij}^1, \dots, v_{ij}^m)$, then $s_{ij} \uparrow i = s_i$. This gives the i -state corresponding to the I -state s , ij -state s_{ij} , respectively. Next we define projection of an I -state onto a pair-system: if $s = (s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, then $s \uparrow ij = (s_i, s_j, v_{ij}^1, \dots, v_{ij}^m)$, where $v_{ij}^1, \dots, v_{ij}^m$ are those values from v^1, \dots, v^n that denote values of variables in \mathcal{SH}_{ij} . This gives the ij -state corresponding to the I -state s , and is well defined only when $i I j$. We also define projection onto the shared variables in \mathcal{SH}_{ij} from both ij -states and I -states: if $s_{ij} = (s_i, s_j, v_{ij}^1, \dots, v_{ij}^m)$, then $s_{ij} \uparrow \mathcal{SH}_{ij} = (v_{ij}^1, \dots, v_{ij}^m)$, and if $s = (s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, then $s \uparrow \mathcal{SH}_{ij} = (v_{ij}^1, \dots, v_{ij}^m)$, where $v_{ij}^1, \dots, v_{ij}^m$ are those values from v^1, \dots, v^n that denote values of variables in \mathcal{SH}_{ij} . Finally, we define projection of an I -state onto a J -system. If $s = (s_{i_1}, \dots, s_{i_K}, v^1, \dots, v^n)$, then $s \uparrow J = (s_{j_1}, \dots, s_{j_L}, v_j^1, \dots, v_j^m)$, where $\{j_1, \dots, j_L\}$ is the domain of J , and v_j^1, \dots, v_j^m are those values from v^1, \dots, v^n that denote values of variables in $\bigcup_{(i,j) \in J} \mathcal{SH}_{ij}$. This gives the J -state (defined analogously to an I -state) corresponding to the I -state s and is well defined only when $J \subseteq I$.

To define projection for paths, we first extend the definition of path (and fullpath) to include the index of the process making the transition, e.g., each transition is labeled by an index denoting this process. For example, a path in M_I would be represented as $s^1 \xrightarrow{d_1} s^2 \dots s^n \xrightarrow{d_n} s^{n+1} \xrightarrow{d_{n+1}} s^{n+2} \dots$, where $\bigwedge m \geq 1. (d_m \in \text{dom}(I))$.

Let π be an arbitrary path in M_I . For any J such that $J \subseteq I$, define a J -block (cf. Clarke et al. [1986] and Browne et al. [1988]) of π to be a maximal subsequence of π that starts and ends in a state and does not contain a transition by any P_i such that $i \in \text{dom}(J)$. Thus we can consider π to be a sequence of J -blocks with successive J -blocks linked by a single P_i -transition such that $i \in \text{dom}(J)$ (note that a J -block can consist of a single state). It also follows that $s \uparrow J = t \uparrow J$ for any pair of states s, t in the same J -block. This is because a transition that is not by some P_i such that $i \in \text{dom}(J)$ cannot affect any atomic proposition in $\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$, nor can it change the value of a variable in $\bigcup_{(i,j) \in J} \mathcal{SH}_{ij}$; and a J -block contains no such P_i transition. Thus, if B is a J -block, we define $B \uparrow J$ to be $s \uparrow J$ for some state s in B . We now give the formal definition of path projection. We use the same notation \uparrow as for state projection. Let B^n denote the n th J -block of π .

Definition 5.1.2.1 (Path Projection). Let π be $B^1 \xrightarrow{d_1} \dots B^n \xrightarrow{d_n} B^{n+1} \dots$ where

B^m is a J -block for all $m \geq 1$. Then the *Path Projection Operator* $\uparrow J$ is given by

$$\pi \uparrow J = B^1 \uparrow J \xrightarrow{d_1} \dots B^n \uparrow J \xrightarrow{d_n} B^{n+1} \uparrow J \dots$$

Thus there is a one-to-one correspondence between J -blocks of π and states of $\pi \uparrow J$, with the n th J -block of π corresponding to the n th state of $\pi \uparrow J$ (note that path projection is well defined when π is finite). For example, consider the three-process mutual exclusion system given in Figure 2. This is synthesized from the two-process mutual exclusion system given in Figure 1 using $I = \{\{1, 2\}, \{2, 3\}, \{3, 1\}\}$. Consider the following path π of this system (for notational clarity, we omit the values of the shared variables):

$$\begin{aligned} &[N_1 N_2 N_3] \xrightarrow{1} [T_1 N_2 N_3] \xrightarrow{3} [T_1 N_2 T_3] \xrightarrow{1} [C_1 N_2 T_3] \xrightarrow{1} \\ &[N_1 N_2 T_3] \xrightarrow{3} [N_1 N_2 C_3] \xrightarrow{3} [N_1 N_2 N_3] \end{aligned}$$

Let $J = \{\{1, 2\}\}$. Then the J -blocks of π are $[N_1 N_2 N_3]$, $[T_1 N_2 N_3] \xrightarrow{3} [T_1 N_2 T_3]$, $[C_1 N_2 T_3]$, and $[N_1 N_2 T_3] \xrightarrow{3} [N_1 N_2 C_3] \xrightarrow{3} [N_1 N_2 N_3]$. Note that all the states within a block agree on all propositions in $\mathcal{AP}_1 \cup \mathcal{AP}_2$. The projected path $\pi \uparrow J$ can be constructed in two steps. First, eliminate all state components not in J from each state:

$$[N_1 N_2] \xrightarrow{1} [T_1 N_2] \xrightarrow{3} [T_1 N_2] \xrightarrow{1} [C_1 N_2] \xrightarrow{1} [N_1 N_2] \xrightarrow{3} [N_1 N_2] \xrightarrow{3} [N_1 N_2]$$

Then collapse all identical states into a single state, eliminating the transitions in between:

$$\pi \uparrow J = [N_1 N_2] \xrightarrow{1} [T_1 N_2] \xrightarrow{1} [C_1 N_2] \xrightarrow{1} [N_1 N_2]$$

5.2 Semantics of the Pair-Systems

The semantics of $(S_{ij}^0, P_i^j \| P_j^i)$ is given by the global state transition diagram M_{ij} generated by its execution. For example, Figure 3 gives the global state transition diagram generated by the execution of the processes in Figure 1 (with the initial state set being $\{[N_1 x_{12} = 1 N_2], [N_1 x_{12} = 2 N_2]\}$). We call the global state transition diagram of a pair-system a *pair-structure*.

Definition 5.2.1 (Pair-Structure). Let $i I j$. The semantics of $(S_{ij}^0, P_i^j \| P_j^i)$ is given by the pair-structure $M_{ij} = (S_{ij}^0, S_{ij}, R_{ij})$ where

- (1) S_{ij} is a set of ij -states,
- (2) $S_{ij}^0 \subseteq S_{ij}$ gives the initial states of $(S_{ij}^0, P_i^j \| P_j^i)$, and
- (3) $R_{ij} \subseteq S_{ij} \times \{i, j\} \times S_{ij}$ is a transition relation giving transitions of $(S_{ij}^0, P_i^j \| P_j^i)$.

A transition (s_{ij}, h, t_{ij}) by $P_h^{\bar{h}}$ is in R_{ij} if and only if all of the following hold:

- (a) $h \in \{i, j\}$,
- (b) s_{ij} and t_{ij} are ij -states, and
- (c) there exists an arc $(s_{ij} \uparrow h, B_h^{\bar{h}} \rightarrow A_h^{\bar{h}}, t_{ij} \uparrow h)$ in $P_h^{\bar{h}}$ such that
 - (i) $s_{ij}(B_h^{\bar{h}}) = \text{true}$,
 - (ii) $\langle s_{ij} \uparrow \mathcal{SH}_{ij} \rangle A_h^{\bar{h}} < t_{ij} \uparrow \mathcal{SH}_{ij} \rangle$, and
 - (iii) $s_{ij} \uparrow \bar{h} = t_{ij} \uparrow \bar{h}$.

Here $\bar{h} = i$ if $h = j$ and $\bar{h} = j$ if $h = i$.

In a transition (s_{ij}, h, t_{ij}) , we say that s_{ij} is the *start* state and that t_{ij} is the *finish* state. The transition (s_{ij}, h, t_{ij}) is called a $P_h^{\bar{h}}$ -transition. In the sequel, we use $s_{ij} \xrightarrow{h} t_{ij}$ as an alternative notation for the transition (s_{ij}, h, t_{ij}) . $\langle s_{ij} \uparrow \mathcal{SH}_{ij} \rangle$ $A \langle t_{ij} \uparrow \mathcal{SH}_{ij} \rangle$ is Hoare triple notation [Hoare 1969] for total correctness, which in this case means that execution of A always terminates,⁴ and, when the shared variables in \mathcal{SH}_{ij} have the values assigned by s_{ij} , leaves these variables with the values assigned by t_{ij} . $s_{ij}(B_h^{\bar{h}}) = \text{true}$ states that the value of guard $B_h^{\bar{h}}$ in state s_{ij} is *true*.⁵

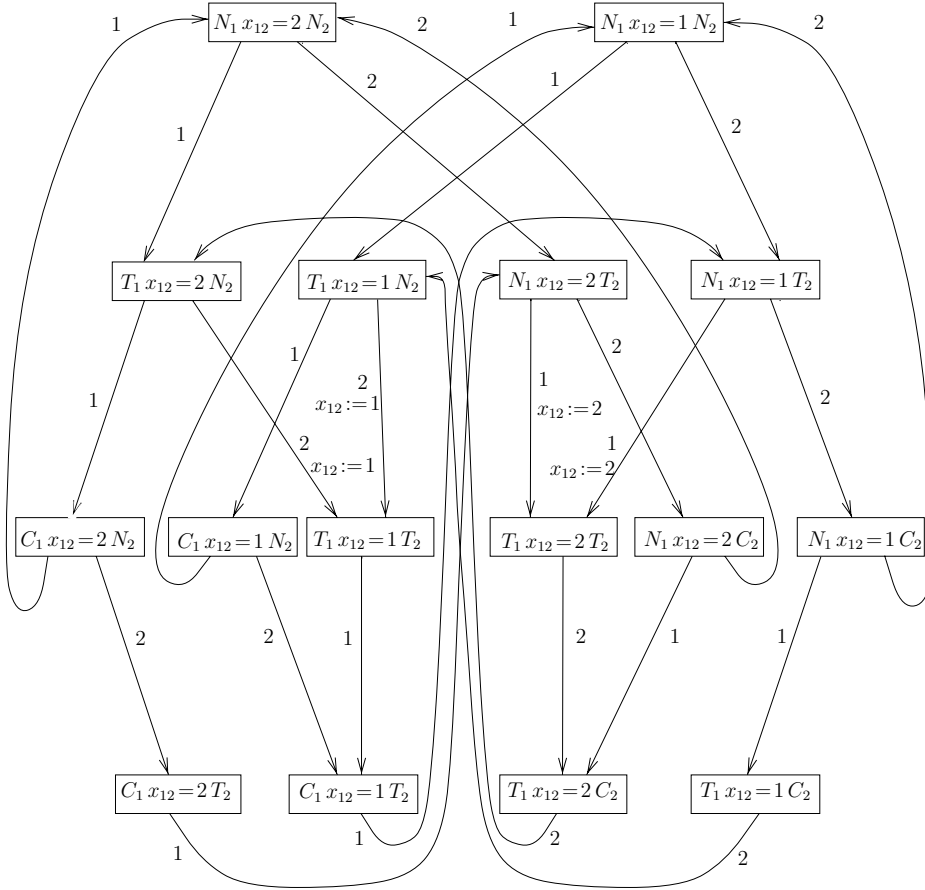
As M_{ij} gives the semantics of $(S_{ij}^0, P_i^j \| P_j^i)$, and we express correctness properties using MPCTL*, we consider that $(S_{ij}^0, P_i^j \| P_j^i)$ possesses a correctness property expressed by an MPCTL* formula $\bigwedge_{k\ell} f_{k\ell}$ if and only if $M_{ij}, S_{ij}^0 \models \bigwedge_{k\ell} f_{k\ell}$ (note that this is equivalent to $M_{ij}, S_{ij}^0 \models f_{ij}$ by MPCTL* semantics). We also remark that, as a consequence of the previous definition, M_{ij} does not contain two or more (differently named) states that are identical when viewed as mappings over the domain $\mathcal{AP}_i \cup \mathcal{AP}_j \cup \mathcal{SH}_{ij}$.

In Figure 3, we see that the only states where the value of the shared variable affects the possible outgoing transitions are $[T_1 x_{12} = 1 T_2], [T_1 x_{12} = 2 T_2]$. Hence, we can minimize the number of global states by collapsing together all other propositionally equivalent states (two states are propositionally equivalent if and only if they assign *true* to exactly the same atomic propositions). This results in the minimized structure shown in Figure 4. State minimization is performed by taking the quotient with respect to strong bisimulation, denoted by \sim [Browne et al. 1988; Clarke et al. 1986], over structures. Essentially, two states are \sim -related if and only if (1) they are propositionally equivalent and (2) all their successor states can be paired into \sim -related pairs (duplicates allowed). Given an arbitrary structure $M_{12}(= (S_{12}^0, S_{12}, R_{12}))$, let $M'_{12}(= (S_{12}^{0'}, S'_{12}, R'_{12}))$ result from minimizing the number of states in M_{12} . Then, the states of M'_{12} are the equivalence classes of S_{12} with respect to \sim : $S'_{12} \stackrel{\text{df}}{=} \{\bar{s} \mid s \in S_{12}\}$, where $\bar{s} \stackrel{\text{df}}{=} \{t \mid s \sim t, s \in S_{12}, t \in S_{12}\}$. The initial states of M'_{12} are those equivalence classes containing some initial state of M_{12} : $S_{12}^{0'} \stackrel{\text{df}}{=} \{\bar{s} \mid s \in S_{12}^0\}$. Finally, two states of M'_{12} are connected by a P_i^j -transition $(i, j \in \{1, 2\}, i \neq j)$ if and only if they have representatives in M_{12} that are connected by a P_i^j -transition: $R_{12} \stackrel{\text{df}}{=} \{\bar{s} \xrightarrow{i} \bar{t} \mid s \xrightarrow{i} t \in R_{12}\}$.

In Figure 4, we display the least information needed to uniquely identify each representative. This necessarily includes all the atomic propositions that are true. It may or may not include shared variable values. For example, $[T_1 N_2]$ represents the equivalence class $\{[T_1 x_{12} = 1 N_2], [T_1 x_{12} = 2 N_2]\}$. The value of x_{12} is not needed here: the atomic propositions T_1 and N_2 being true suffices to identify the states of this equivalence class. By contrast, $[T_1 x_{12} = 1 T_2]$ represents the (singleton) equivalence class $\{[T_1 x_{12} = 1 T_2]\}$, and here, the value of x_{12} is needed

⁴Termination is obvious, since the right-hand side of A is a list of constants.

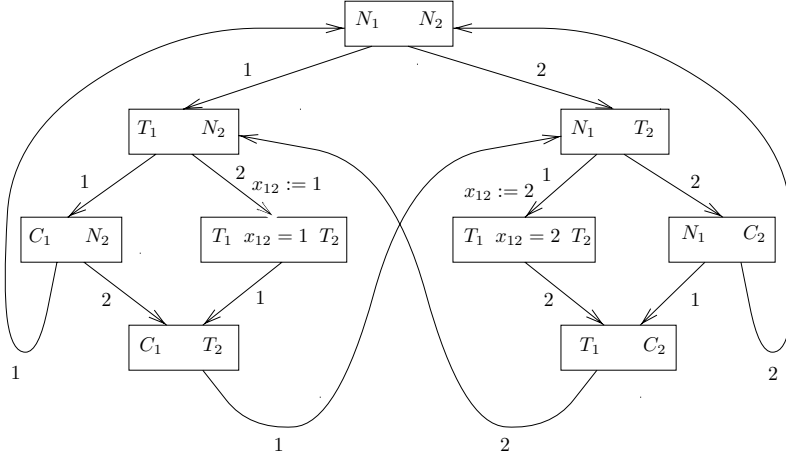
⁵ $s_{ij}(B_h^{\bar{h}})$ is defined by the usual inductive scheme: $s_{ij}(\text{"}x_{ij} = h_{ij}\text{"}) = \text{true}$ iff $s_{ij}(x_{ij}) = h_{ij}$, $s_{ij}(B_1^{\bar{h}} \wedge B_2^{\bar{h}}) = \text{true}$ iff $s_{ij}(B_1^{\bar{h}}) = \text{true}$ and $s_{ij}(B_2^{\bar{h}}) = \text{true}$, $s_{ij}(\neg B_1^{\bar{h}}) = \text{true}$ iff $s_{ij}(B_1^{\bar{h}}) = \text{false}$.



The initial state set is $\{[N_1 x_{12} = 2 N_2], [N_1 x_{12} = 1 N_2]\}$.

Fig. 3. Global state transition diagram of the two-process mutual exclusion example.

to distinguish from $[T_1 x_{12} = 2 T_2]$, which represents the (singleton) equivalence class $\{[T_1 x_{12} = 2 T_2]\}$. Note that, in Figure 3, the states $[T_1 x_{12} = 1 T_2]$ and $[T_1 x_{12} = 2 T_2]$ are \sim -inequivalent, since they have one successor state each, and these successor states (namely, $[C_1 x_{12} = 1 T_2]$ and $[T_1 x_{12} = 2 C_2]$) are propositionally inequivalent, and so cannot be \sim -related. Since the initial states of the structures of Figures 3 and 4 are \sim -related, they satisfy the same CTL* formulae, by Theorem 2 of Clarke et al. [1986]. Hence in particular, the structure of Figure 4 satisfies the CTL* specification for the two-process mutual exclusion problem. The strong bisimulation relation can be computed in time $O(|S_{12}|^2)$ as a fixpoint, as outlined in Milner [1989, Ch. 4]. Hence, the reduced form can be computed automatically in time $O(|S_{12}|^2)$.



The initial state set is $\{ [N_1 \ N_2] \}$.

Fig. 4. Reduced global state transition diagram of the two-process mutual exclusion example.

5.3 Semantics of I -Systems and J -Systems

The semantics of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ is given by the global state transition diagram M_I generated by its execution. We call the global state transition diagram of an I -system an I -structure.

Definition 5.3.1 (I -Structure). The semantics of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ is given by the I -structure $M_I = (S_I^0, S_I, R_I)$ where

- (1) S_I is a set of I -states,
- (2) $S_I^0 \subseteq S_I$ gives the initial states of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$, and
- (3) $R_I \subseteq S_I \times \{i_1, \dots, i_K\} \times S_I$ is a transition relation giving the transitions of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$. A transition (s, i, t) by P_i^I is in R_I if and only if
 - (a) $i \in \{i_1, \dots, i_K\}$,
 - (b) s and t are I -states, and
 - (c) there exists an arc $(s \uparrow i, \bigwedge_{j \in I(i)} B_i^j \rightarrow //_{j \in I(i)} A_i^j, t \uparrow i)$ in P_i^I such that all of the following hold:
 - (i) for all j in $I(i)$, $s \uparrow i j(B_i^j) = \text{true}$ and $\langle s \uparrow \mathcal{H}_{ij} > A_i^j < t \uparrow \mathcal{H}_{ij} \rangle$,
 - (ii) for all j in $\{i_1, \dots, i_K\} - \{i\}$: $s \uparrow j = t \uparrow j$, and
 - (iii) for all j, k in $\{i_1, \dots, i_K\} - \{i\}$, $j \neq k$: $s \uparrow \mathcal{H}_{jk} = t \uparrow \mathcal{H}_{jk}$.

In a transition (s, i, t) , we say that s is the *start* state, and t is the *finish* state. The transition (s, i, t) is called a P_i^I -transition. In the sequel, we use $s \xrightarrow{i} t$ as alternative notation for the transition (s, i, t) . Also, if I is set to $\{\{i, j\}\}$ in Definition 5.3.1, then the result is, as expected, the pair-structure definition (5.2.1). In other words, the two definitions are consistent. Furthermore, the semantics of a J -system, $J \subseteq I$ (see Section 5.1.1) is given by the J -structure $M_J = (S_J^0, S_J, R_J)$, which is obtained by using J for I in Definitions 5.1.1 and 5.3.1.

As M_I gives the semantics of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$, and we express correctness properties using MPCTL*, we consider that $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ possesses a correctness property expressed by an MPCTL* formula $\bigwedge_{k\ell} f_{k\ell}$ if and only if $M_I, S_I^0 \models \bigwedge_{k\ell} f_{k\ell}$ (note that this is equivalent to $M_I, S_I^0 \models \bigwedge (i, j) \in I. (f_{ij})$ by MPCTL* semantics).

5.4 A Compact Representation for Synchronization Skeletons

Suppose P_i^j (for every $j \in I(i)$) contains two arcs from i -state s_i to i -state s'_i , e.g., $a_{i,1}^j = (s_i, B_{i,1}^j \rightarrow A_{i,1}^j, s'_i)$ and $a_{i,2}^j = (s_i, B_{i,2}^j \rightarrow A_{i,2}^j, s'_i)$. Then, by the MP-synthesis definition (5.1.1), P_i^I contains $2^{|I(i)|}$ arcs from i -state s_i to i -state s'_i , one arc for each element of the cartesian product:

$$\{a_{i,1}^{j_1}, a_{i,2}^{j_1}\} \times \dots \times \{a_{i,1}^{j_n}, a_{i,2}^{j_n}\}$$

where $\{j_1, \dots, j_n\} = I(i)$. Thus, P_i^I is exponentially large in $K (= |dom(I)|)$ in the worst case (since $|I(i)| = K - 1$ when $i \neq j$ for every j in $dom(I) - \{i\}$), which defeats the purpose of MP-synthesis. We deal with this by defining a compact representation for processes in which there is at most one arc between any pair of (local) states, thereby avoiding the exponential blowup illustrated above. We provide an overview of this notation here, with full details and definitions in Appendix D.

Consider a pair-process P_i^j which has two arcs from state s_i to state s'_i , labeled with the synchronization commands $B_{i,1}^j \rightarrow A_{i,1}^j, B_{i,2}^j \rightarrow A_{i,2}^j$. In compact notation, we replace these two arcs by a single arc whose label is $B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j$. \oplus is a binary operator which takes a pair of guarded commands as arguments. Roughly, the operational semantics of $B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j$ is that if one of the guards $B_{i,1}^j, B_{i,2}^j$ evaluates to true, then the corresponding body $A_{i,1}^j, A_{i,2}^j$ respectively, can be executed. If neither $B_{i,1}^j$ nor $B_{i,2}^j$ evaluates to true, then the command “blocks,” i.e., waits until one of $B_{i,1}^j, B_{i,2}^j$ evaluates to true.⁶ We call an arc whose label has the form $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$ a (*pair*) *move*. In compact notation, a pair-process has at most one move between any pair of local states.

Now an I -process P_i^I is derived by “composing” all the moves of P_i^j (as j varies over $I(i)$) which have the same start and end states. Towards this end we define the binary composition operator \otimes on guarded commands. Roughly, the operational semantics of $B_{i,1}^j \rightarrow A_{i,1}^j \otimes B_{i,2}^j \rightarrow A_{i,2}^j$ is that if both of the guards $B_{i,1}^j, B_{i,2}^j$ evaluate to true, then the bodies $A_{i,1}^j, A_{i,2}^j$ can be executed in parallel. If at least one of $B_{i,1}^j, B_{i,2}^j$ evaluates to false, then the command “blocks,” i.e., waits until both of $B_{i,1}^j, B_{i,2}^j$ evaluate to true.

We now recast the MP-synthesis Definition (5.1.1) into compact form as follows.

Definition 5.4.1 (Compact MP-Synthesis). A *compact I-process* P_i^I is derived from the compact pair-process P_i^j as follows:

P_i^I contains a move from s_i to t_i with label $\otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$
 iff
 for every j in $I(i)$: P_i^j contains a move from s_i to t_i with label $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$.

⁶This interpretation was proposed by Dijkstra [1982].

The *initial state set* S_I^0 of the I -system is derived from the initial state S_{ij}^0 of the pair-system as follows:

$$S_I^0 = \{s \mid \bigwedge (i, j) \in I. (s \upharpoonright ij \in S_{ij}^0)\}.$$

Thus an I -process in compact notation has at most one move of the form $(s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$ between any pair of i -states s_i, t_i . The I -move $(s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$ can be regarded as a type of “conjunctive normal form” for guarded commands, being a conjunction over all j in $I(i)$ of the pair-moves $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$, each of which can be regarded as a “disjunction” of guarded commands.

A move in a pair-process is simply a special case of a move in an I -process when I is a single pair, i.e., if $I = \{\{i, k\}\}$, then $I(i)$ in $\otimes_{j \in I(i)}$ expands to the singleton $\{k\}$ giving a move in the pair-process P_i^k .

We define the following notation. We use a_i^j, a_i^I to denote moves in P_i^j, P_i^I , respectively. Consider a move a_i^I of P_i^I . By the compact MP-synthesis definition (5.4.1), a_i^I has the form $(s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$. Let $a_i^I.start, a_i^I.guard_j$ denote $s_i, \bigvee_{\ell \in [1:n]} B_{i,\ell}^j$, respectively, and let $a_i^I.guard$ denote $\bigwedge_{j \in I(i)} a_i^I.guard_j$. In the special case $I = \{\{i, j\}\}$, let a_i^j denote the move $(s_i, \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$, and let $t_k.moves$ denote the set $\{a_k^I \mid a_k^I \in P_k^I \text{ and } a_k^I.start = t_k\}$.

The semantics of compact pair-systems and compact I -systems is given by *compact pair-structures* and *compact I-structures*, respectively. The following definitions are a direct consequence of the meaning of \oplus, \otimes , and the definitions for pair-structures (Definition 5.2.1) and I -structures (Definition 5.3.1).

Definition 5.4.2 (Compact Pair-Structure). Let $i I j$. The semantics of $(S_{ij}^0, P_i^j \parallel P_j^i)$ in compact notation is given by the pair-structure $M_{ij} = (S_{ij}^0, S_{ij}, R_{ij})$ where

- (1) S_{ij} is a set of ij -states,
- (2) $S_{ij}^0 \subseteq S_{ij}$ gives the initial states of $P_i^j \parallel P_j^i$, and
- (3) $R_{ij} \subseteq S_{ij} \times \{i, j\} \times S_{ij}$ is a transition relation giving the transitions of $P_i^j \parallel P_j^i$.

A transition (s_{ij}, h, t_{ij}) by $P_h^{\bar{h}}$ is in R_{ij} if and only if all of the following hold:

- (a) $h \in \{i, j\}$,
- (b) s_{ij} and t_{ij} are ij -states, and
- (c) there exists a move $(s_{ij} \upharpoonright h, \oplus_{\ell \in [1:n]} B_{h,\ell}^{\bar{h}} \rightarrow A_{h,\ell}^{\bar{h}}, t_{ij} \upharpoonright h)$ in $P_h^{\bar{h}}$ such that there exists $m \in [1 : n]$:
 - (i) $s_{ij}(B_{h,m}^{\bar{h}}) = \text{true}$,
 - (ii) $\langle s_{ij} \upharpoonright h \mathcal{H}_{ij} \rangle A_{h,m}^{\bar{h}} \langle t_{ij} \upharpoonright h \mathcal{H}_{ij} \rangle$, and
 - (iii) $s_{ij} \upharpoonright \bar{h} = t_{ij} \upharpoonright \bar{h}$.

Here $\bar{h} = i$ if $h = j$ and $\bar{h} = j$ if $h = i$.

Definition 5.4.3 (Compact I-Structure). The semantics of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ in compact notation is given by the I -structure $M_I = (S_I^0, S_I, R_I)$ where

- (1) S_I is a set of I -states,

- (2) $S_I^0 \subseteq S_I$ gives the initial states of $P_{i_1}^I \parallel \dots \parallel P_{i_K}^I$, and
- (3) $R_I \subseteq S_I \times \text{dom}(I) \times S_I$ is a transition relation giving the transitions of $P_{i_1}^I \parallel \dots \parallel P_{i_K}^I$.
 A transition (s, i, t) by P_i^I is in R_I if and only if
- (a) $i \in \text{dom}(I)$,
 - (b) s and t are I -states, and
 - (c) there exists a move $(s \uparrow i, \otimes_{j \in I(i) \oplus \ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t \uparrow i)$ in P_i^I such that all of the following hold:
 - (i) for all j in $I(i)$, there exists $m \in [1:n]$:
 $s \uparrow ij(B_{i,m}^j) = \text{true}$ and $\langle s \uparrow \mathcal{SH}_{ij} \rangle A_{i,m}^j < t \uparrow \mathcal{SH}_{ij} \rangle$,
 - (ii) for all j in $\text{dom}(I) - \{i\}$: $s \uparrow j = t \uparrow j$, and
 - (iii) for all j, k in $\text{dom}(I) - \{i\}$, $j I k$: $s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}$.

Returning to the example given at the beginning of this section, the pair of arcs $(s_i, B_{i,1}^j \rightarrow A_{i,1}^j, s'_i)$, $(s_i, B_{i,2}^j \rightarrow A_{i,2}^j, s'_i)$ in P_i^j are replaced by the single move $(s_i, (B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j), s'_i)$, and the $2^{|I(i)|}$ arcs in P_i^I are replaced by the single move $(s_i, \otimes_{j \in I(i)} (B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j), s'_i)$. A full discussion of compact representation is given in Appendix D.

5.5 Examples

Using the compact notation, we synthesize the following programs. We start with the compact solution to the two-process mutual exclusion problem, which is shown in Figure 5 (this is simply the compact representation of the solution shown in Figure 1). From this we synthesize the process P_1 in the solution for the original dining philosophers problem (five philosophers in a circle), which is shown in Figure 6. The processes $P_2 - P_5$ can be obtained from P_1 by appropriate process index substitutions. Figure 7 gives, in compact notation, the process P_i^I in the solution for the generalized dining philosophers problem. The interconnection relation I gives the neighborhood relationship among philosophers. If I is a complete graph, then we get a solution to the K -process mutual exclusion problem.

Finally, Figure 8 gives the process P_i^I in the solution to the drinking philosophers problem [Chandy and Misra 1984; Chandy and Misra 1988]. r_i^j is a shared variable which is assigned to nondeterministically upon entry to the trying region. If it is set to *true*, then P_i^I will, upon entry into its critical section, require the “bottle” that is shared with P_j^I . If, on the other hand, r_i^j is set to *false*, then P_i^I will not require the shared bottle upon critical section entry. Thus, if r_i^j and r_j^i are both set to *true*, then P_i^I and P_j^I will contend for the bottle that they share, and so will exclude each other upon entry to the critical section this time around. If either (or both) of r_i^j, r_j^i are set to *false*, then P_i^I and P_j^I may enter their critical sections simultaneously this time around, since at least one of them will not require access to the shared bottle. We have not shown the two-process drinking philosophers solution here.

In Figures 7 and 8, the interconnection relation I is not fixed, but is a parameter to the synthesized program. A single such parameterized program thus represents an infinite family of “concrete” programs, one for each value of I . In particular, I can be “arbitrarily large.” For example, we could set $I = \{(i, j) \mid i \neq j, i, j \in \{1, \dots, 1000\}\}$. Figure 7 would then represent a mutual exclusion solution for 1000

processes. Figure 9 shows process P_1 of this program. As this program has on the order of $2^{499,500}$ global states (there are $1000 * 999/2 = 499,500$ binary pairwise shared variables), straightforward state exploration is clearly infeasible for the synthesis of this program. By using other “large” values for I , we can obtain solutions containing large numbers of processes for the generalized dining and drinking philosophers problems from Figures 7 and 8, respectively. For example, if $I = \{\{1, 2\}, \{2, 3\}, \dots, \{999, 1000\}, \{1000, 1\}\}$, then we obtain dining, drinking philosophers solutions for 1000 philosophers arranged in a circle. One very important advantage of the compact notation is that it permits this “parameterized” representation, whereas our normal notation does not.

6. SOUNDNESS OF THE METHOD

Since we synthesize I -systems from pair-systems, we formulate the soundness of the synthesis method as follows: if a pair-system has a particular correctness property, then the I -system also has that property. As stated previously, we express correctness properties *pairwise*, i.e., as MPCTL* formulae. Furthermore, we only consider a subset of MPCTL*—which nonetheless can express many safety and liveness properties that arise in practice, e.g., invariants that can be expressed pairwise, and temporal leads-to properties over individual processes. We discuss correctness properties in more detail in Section 6.1. Also, independently of our soundness results, we characterize sufficient conditions under which the I -system is deadlock-free.

Leading up to our main soundness results are many preliminary results that establish the prerequisite relationships between pair-structures and I -structures. Our results can be grouped into the following sequence:

- Similarity Results (Section 6.2)*: Formalize the notion of similarity of processes and establish that all pair-structures are similar.
- Projection Results (Section 6.3)*: Establish that a state and its projection satisfy the same propositional formulae over their common atomic propositions. Also establish that a path and its projection satisfy the same linear-time temporal logic, i.e., PTL (see Section 3.1), formulae over their common atomic propositions. These results relate states (paths) and their projections independently of the occurrence of these states (paths) in the structures of interest, namely the pair and I -structures.
- Mapping Results (Section 6.4)*: Establish that a P_i^I -transition occurs in the I -structure if and only if all of its projections onto $\{\{i, j\}\}$, for $j \in I(i)$, occur in the corresponding ij -structures. Also establish that if a state (path) occurs in the I -structure, then its projection onto J occurs in the J -structure (for any $J \subseteq I$). In other words, these results relate the occurrence of states (paths) in the I -structure to the occurrence of states (paths) in the J -structure. Together with the projection results, the mapping results provide a basis for relating the MPCTL* formulae satisfied by the I -structure to the MPCTL* formulae satisfied by the pair-structures.
- Deadlock-Freedom Results (Section 6.5)*: Establish that the I -system is free of deadlock. This is done using a technical assumption (the “wait-for-graph assumption”) which, roughly, guarantees that circular waiting chains cannot form.

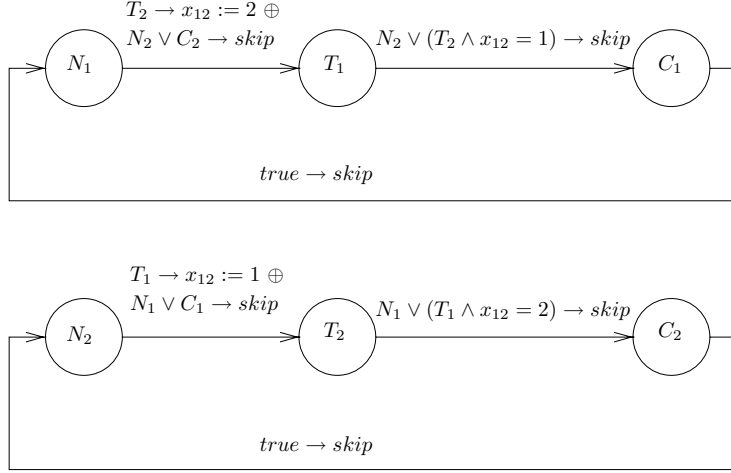


Fig. 5. Compact skeletons P_1^2 (top) and P_2^1 (bottom) for two-process mutual exclusion.

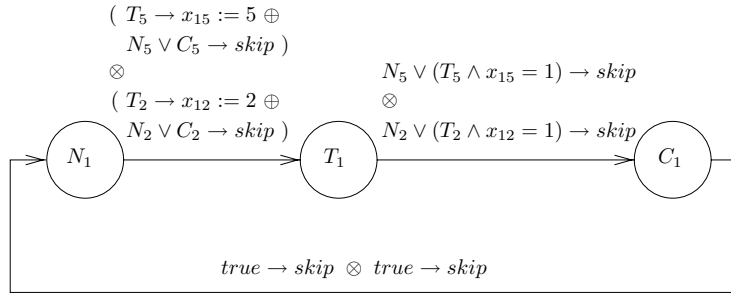


Fig. 6. Compact skeleton P_1 in the solution to the original dining philosophers problem.

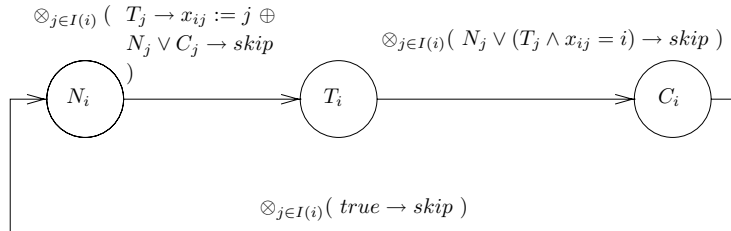
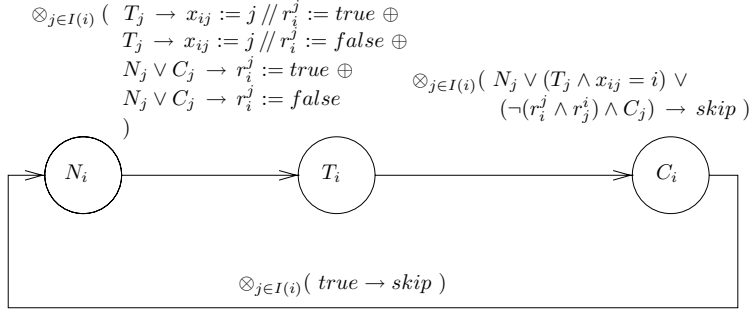
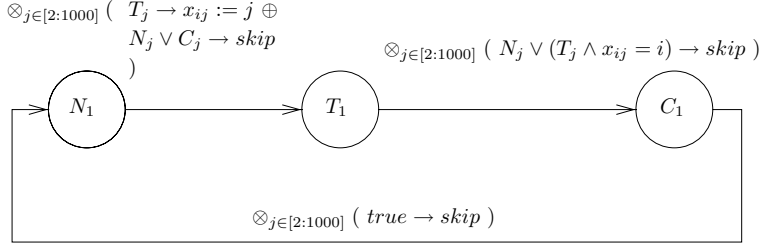


Fig. 7. Compact skeleton P_i in the solution to the generalized dining philosophers problem.

Fig. 8. Compact skeleton P_i in the drinking philosophers solution.Fig. 9. Compact skeleton P_1 in the solution to the 1000-process mutual exclusion problem.

Uses the mapping results and the compact representation. These results stand alone as characterizing conditions under which the I -system is deadlock-free. In the soundness results, deadlock-freeness is assumed.

- Liveness Results (Section 6.7)*: Establish that individual starvation of processes cannot occur in the I -system. Again this requires a technical assumption (the “liveness assumption”). This assumption prevents a process from executing forever while indirectly (i.e., via a “chain” of other processes) blocking another process that needs to progress. Uses the mapping and deadlock-freeness results and the compact representation.
- Soundness Results (Sections 6.6 and 6.8)*: Our main soundness results. We establish two results, one for completely interconnected systems (i.e., every pair of processes are I -related) and another for systems with arbitrary interconnection. Uses the projection, mapping, and liveness results.

6.1 Correctness Properties and Reachability

Correctness properties in our framework are specified as MPCTL* formulae. Since MPCTL* formulae are evaluated on particular states of a structure, a suitable notion of correctness is that the formula specifying the correctness property of interest is true in all initial states of the structure. In other words, P satisfies f if and only if $M, S^0 \models f$, where $M = (S^0, S, R)$ is the global state transition diagram

of P . Now, from CTL* semantics, we see that the truth value of an MPCTL* formula in an initial state of a structure depends only on the “reachable” portion of the structure. We therefore define an *initialized path* to be a path whose first state is an initial state, and a *reachable state* to be a state lying on some initialized path. Also define a *reachable transition* to be a transition lying along some initialized path.

In the sequel, we assume that every i -state s_i of P_i^j is reachable in M_{ij} , i.e., for every s_i , there exists a reachable ij -state s_{ij} such that $s_{ij} \uparrow i = s_i$. Any such s_i that is not reachable can be removed from P_i^j (along with all its incident arcs) without affecting the reachable portion of M_{ij} generated by the execution of $(S_{ij}^0, P_i^j \parallel P_j^i)$, and hence the CTL* formulae satisfied by M_{ij} .

We define the following notation: S_I^r is the set of reachable states of M_I , and $R_I^r = R_I \upharpoonright S_I^r$ is the restriction of the transition relation R_I to the set of reachable states of M_I . M_I^r is the structure (S_I^0, S_I^r, R_I^r) . If $I = \{\{i, j\}\}$ (i.e., the two-process case) then we use $S_{ij}^r, R_{ij}^r, M_{ij}^r$ for S_I^r, R_I^r, M_I^r , respectively. For an arbitrary $J \subseteq I$, we define S_J^r, R_J^r , and $M_J^r = (S_J^0, S_J^r, R_J^r)$ by using J for I in the previous definitions.

In this article, the correctness properties of interest are as follows, where $h_{k\ell}$ is a formula of $\mathcal{L}(\mathcal{AP}_k, \mathcal{AP}_\ell, \neg, \wedge)$, and p_k, q_k, r_k are formulae of $\mathcal{L}(\mathcal{AP}_k, \neg, \wedge)$:

- $\bigwedge_{k\ell} h_{k\ell}$: Specifies the initial states; all initial states must satisfy $h_{k\ell}$.
- $\bigwedge_{k\ell} AGh_{k\ell}$: Expresses a propositional invariant; $h_{k\ell}$ must be true in all reachable states.
- $\bigwedge_k AG(p_k \Rightarrow A[q_k Ur_k])$: Expresses a temporal leads-to properties, combined with an “unless” property; from any global state satisfying p_k , eventually a state satisfying r_k must be reached (along all outgoing fullpaths), and all intervening states must satisfy q_k .
- $\bigwedge_k AG(p_k \Rightarrow AY_k q_k)$: Expresses a “local structure” property; every local state of process k that satisfies p_k has as immediate successors in the skeleton for process k only local states that satisfy q_k .⁷
- $\bigwedge_k AG(p_k \Rightarrow EX_k q_k)$: Expresses a “local structure” property; every local state of process k that satisfies p_k has some immediate successor in the skeleton for process k which satisfies q_k , and when process k is in the former local state, it always has some enabled transition to the latter local state, no matter what the current global state is.⁸

6.2 Process Similarity

Our notion of process similarity is a syntactic one: two processes are *similar* if and only if one can be obtained from the other by swapping their indices. To formalize this, we introduce the *process index substitution operator* $\theta = \{j_1/i_1, \dots, j_m/i_m\}$ which denotes the simultaneous replacement of process indices i_1, \dots, i_m by process indices j_1, \dots, j_m , respectively. We require that i_1, \dots, i_m be pairwise distinct, and j_1, \dots, j_m be pairwise distinct, and so θ specifies a “renaming” of process indices and can be applied to all of the pair syntactic constructs defined in the article, as well as any pair model (e.g., M_{ij}). We define θ formally in Appendix C. To preserve

⁷We ignore the possibility of “dead code,” i.e., unreachable portions of the skeleton for process k .

⁸In our technical results, some further technical restrictions on the form of p_k, q_k are made.

the similarity of pair-processes under interchange of process indices, we make the convention that the shared variables x_{ij} and x_{ji} are actually the same variable, i.e., $x_{ij}\{i/j, j/i\}$, x_{ji} , and x_{ij} all denote the same variable. However, if the variable is superscripted, then index interchange results in a new variable. Thus $x_{ij}^i\{i/j, j/i\}$ and x_{ji}^j denote the same variable, which is a different variable than that denoted by x_{ij}^i . Our assumption of process similarity is now formally stated as follows:

Process Similarity Assumption. For all $i, j, i', j' \in \{i_1, \dots, i_K\}$ such that $i \neq j$, $i' \neq j'$, $i I j$, $i' I j'$,

$$P_i^j = P_{i'}^{j'}\{i/i', j/j'\}.$$

In other words, any pair-process can be derived from any other pair-process by index substitution. For special cases of the similarity relation among pair-processes, we have $P_i^j = P_j^i\{i/j, j/i\}$, $P_i^j = P_i^{j'}\{j/j'\}$, and $P_i^j = P_{i'}^j\{i/i'\}$.

The MP-synthesis definition (5.1.1) requires us to generate P_i^j for every pair (i, j) in I . Using the process similarity assumption, we only require a single pair-system in order to generate all the required pair-systems. For example, (assuming $(1, 2) \in I$), if we are given the pair-system $(S_{12}^0, P_1^2 \| P_2^1)$, then we can generate P_i^j for any pair (i, j) in I using $P_i^j = P_1^2\{i/1, j/2\}$, which is obtained from the process similarity assumption by letting $i' = 1, j' = 2$.

Analogously to our assumption that all pair-processes are similar, we assume that all pair-systems have similar sets of initial states.

Initial State Assumption. Let i, j, i', j' be arbitrary elements of $\{i_1, \dots, i_K\}$ such that $i I j$ and $i' I j'$. Then we have

$$S_{ij}^0 = S_{i'j'}^0\{i/i', j/j'\}.$$

The initial state assumption guarantees that S_I^0 (given by the MP-synthesis definition (5.1.1)) is nonempty, given that S_{ij}^0 is nonempty. Since both the initial states and the processes of all pair systems are similar, it follows that the pair-structures generated by the executions of the pair-systems are similar.

PROPOSITION 6.2.1 (PAIR-STRUCTURE SIMILARITY). *Let i, j, i', j' be arbitrary elements of $\{i_1, \dots, i_K\}$ such that $i I j$ and $i' I j'$. Then we have*

$$M_{ij} = M_{i'j'}\{i/i', j/j'\}.$$

By MPCTL* semantics (Section 3), an MPCTL* formula is not sensitive to the particular process indices in a structure. Hence, if two structures are identical up to a renaming of process indices, then they satisfy exactly the same MPCTL* formulae. Together with Proposition 6.2.1, this yields the following observation:

Observation 6.2.2. Let i, j, i', j' be arbitrary elements of $\{i_1, \dots, i_K\}$ such that $i I j$ and $i' I j'$. Then we have

$$M_{ij}, S_{ij}^0 \models \bigwedge_{k\ell} f_{k\ell} \text{ iff } M_{i'j'}, S_{i'j'}^0 \models \bigwedge_{k\ell} f_{k\ell}.$$

6.3 State and Path Projection Results

This section relates states (paths, respectively) and their projection in terms of the CTL* formulae that they agree on. No assumption is made about the occurrence of

these states (paths) in the structures of interest, namely the pair and I -structures.

Since an I -state s and its projection $s\upharpoonright J$ agree, by definition of \upharpoonright , on all propositions in $\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$, we expect these states to satisfy the same propositional logic formulae which contain atomic propositions only from $\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$.

PROPOSITION 6.3.1 (I-STATE PROJECTION). *Let $J \subseteq I$, and let s be an I -state. Then*

$$s \models f \text{ iff } s\upharpoonright J \models f$$

where f is a formula of $\mathcal{L}(\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i, \neg, \wedge)$.

Likewise, a J -state s_J and its projection $s_J\upharpoonright i$ satisfy exactly the same propositional logic formulae which contain atomic propositions only from \mathcal{AP}_i ($i \in \text{dom}(J)$).

PROPOSITION 6.3.2 (LOCAL STATE PROJECTION). *Let $i \in \text{dom}(J)$, and let s_J be a J -state. Then*

$$s_J \models f_i \text{ iff } s_J\upharpoonright i \models f_i$$

where f_i is a formula of $\mathcal{L}(\mathcal{AP}_i, \neg, \wedge)$.

We remind the reader that a J -block is a maximal subsequence of a path containing no transition by a process in the domain of J (Section 5.1.2). Since a J -block B of a path π and its corresponding state $B\upharpoonright J$ in $\pi\upharpoonright J$ agree on all atomic propositions in $\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$, the sequence of truth values assigned to these propositions along the paths π and $\pi\upharpoonright J$ is the same if we ignore repetitions, i.e., allow *stuttering* [Browne et al. 1988; Clarke et al. 1986]. Hence, these paths satisfy exactly the same linear-time temporal logic, i.e., the logic PTL (see Section 3.1), formulae provided that we disallow the next-time operator X , which allows repetition to be detected.

LEMMA 6.3.3 (PATH PROJECTION). *Let π be a path in M_I , and let $J \subseteq I$. Then*

$$\pi \models f \text{ iff } \pi\upharpoonright J \models f$$

where f is a formula of $\mathcal{L}(\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i, \neg, \wedge, U)$.

6.4 Mapping of I -Structures into J -Structures

This section relates the occurrence of transitions, states, and paths in the I -structure to the occurrence of the projected transitions, states, and paths respectively in J -structures ($J \subseteq I$), paying particular attention to the special case where J specifies a single pair (Lemma 6.4.1). Specifically, it shows that projection onto J defines a mapping of the transitions, reachable states, and paths of M_I into the transitions, reachable states, and paths of M_J .

By virtue of an arc in an I -process P_i^I being a “composition” of the corresponding arcs in the pair-processes P_i^j in all the pair-systems $(S_{ij}^0, P_i^j \| P_j^i)$ (as j ranges over $I(i)$), a transition by P_i^I can occur in the I -system if and only if the “projected” transitions by P_i^j in all the pair systems $(S_{ij}^0, P_i^j \| P_j^i)$ (as j ranges over $I(i)$) can all occur. This is formalized as the transition-mapping lemma, which provides the basis for establishing the remaining mapping results.

LEMMA 6.4.1 (TRANSITION MAPPING). *For all I -states $s, t \in S_I$ and $i \in \text{dom}(I)$,*

$s \xrightarrow{i} t \in R_I$ iff :

$$\begin{aligned} & \bigwedge j \in I(i) . (s \uparrow i j \xrightarrow{i} t \uparrow i j \in R_{ij}) \text{ and} \\ & \bigwedge j \in \{i_1, \dots, i_K\} - \hat{I}(i) . (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \{i_1, \dots, i_K\} - \{i\}, j I k . (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned}$$

We remind the reader of the example of “composition” of transitions (Section 5.1), which illustrates Lemma 6.4.1. Note that Lemma 6.4.1 relates M_I to a set of J -systems (namely, for all $J = (i, j)$ as j ranges over $I(i)$), whereas the remaining results below relate M_I to M_J for a single, arbitrary $J \subseteq I$.

Since, by Lemma 6.4.1, a transition by P_i^I entails “projected” transitions by P_i^j in all the pair systems $(S_{ij}^0, P_i^j \parallel P_j^i)$ as j ranges over $I(i)$, it follows that a transition by P_i^I entails “projected” transitions by P_i^j in all the pair-systems $(S_{ij}^0, P_i^j \parallel P_j^i)$ as j ranges over $J(i)$, where $J \subseteq I$. This latter set of transitions entails (again by Lemma 6.4.1) a corresponding transition by P_i^J in the J -system. Hence, we conclude that the projection onto J of every transition in the I -system is a transition in the J -system:

COROLLARY 6.4.2 (TRANSITION MAPPING). *Let $J \subseteq I$ and $i \in \text{dom}(J)$. If $s \xrightarrow{i} t \in R_I$, then $s \uparrow J \xrightarrow{i} t \uparrow J \in R_J$.*

By applying the transition-mapping corollary to every transition along a path π in M_I , we show that $\pi \uparrow J$ is a path in M_J :

LEMMA 6.4.3 (PATH MAPPING). *Let $J \subseteq I$. If π is a path in M_I , then $\pi \uparrow J$ is a path in M_J .*

As an example of this lemma, consider the path π given in Section 5.1.2 and its projection $\pi_{12} = \pi \uparrow \{1, 2\}$, where π is a path of the three-process mutual exclusion system given in Figure 2. We can verify by inspection that π_{12} is a path of the two-process mutual exclusion system given in Figure 1.

From Definition 5.1.1, we see that an initial I -state projects onto an initial J -state. Hence, a consequence of Lemma 6.4.3 is that initialized paths (see Section 6.1) in M_I project onto initialized paths in M_J :

COROLLARY 6.4.4 (PATH MAPPING). *Let $J \subseteq I$. If π is an initialized path in M_I then $\pi \uparrow J$ is an initialized path in M_J .*

Since every reachable state lies at the end of some initialized path, we can use the path-mapping corollary to relate reachable states in M_I to their projections in M_J :

COROLLARY 6.4.5 (STATE MAPPING). *Let $J \subseteq I$. If t is a reachable state in M_I , then $t \uparrow J$ is a reachable state in M_J .*

We relativize our notion of reachability as follows. We say that a state t is *s-reachable* in M_I if and only if there exists a path in M_I from s to t . Note that neither s nor t need be reachable states. The definition for pair-systems, is, as usual, obtained by setting I to $\{\{i, j\}\}$. Since, by Lemma 6.4.3 and the definition of path projection, the projection onto J of a path in M_I from I -state s to I -state t gives a path in M_J from $s \uparrow J$ to $t \uparrow J$, we have the following corollary:

COROLLARY 6.4.6 (RELATIVIZED STATE MAPPING). *Let $J \subseteq I$. If t is a s -reachable state in M_I , then $t \upharpoonright J$ is a $s \upharpoonright J$ -reachable state in M_J .*

6.5 Deadlock Freedom of the I -System

Our focus in this article is nonterminating concurrent systems, so we assume that $(S_{ij}^0, P_i^j \parallel P_j^i)$ is nonterminating, i.e., deadlock-free.⁹ Now it is possible that $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ is deadlock-prone; we present an example below of a deadlock-prone I -system (for $I = \{1, 2\}, \{2, 3\}, \{3, 1\}\}$) that is synthesized from a deadlock-free pair-system. Thus, even though the pair-system itself is deadlock-free, it permits a certain pattern of blocking behavior that allows deadlock to occur in the synthesized I -system. We shall see below that this blocking behavior is that a process executes a transition which results in a global state where the process both blocks some move of another process¹⁰ and has all of its own moves blocked by other processes. Hence, some technical assumption on the pair-system $(S_{ij}^0, P_i^j \parallel P_j^i)$ that is input to the MP-synthesis definition (5.1.1) is needed. To formulate this assumption, we first define a notion of “wait-for-graph” for our model of parallel computation. Our model of waiting is a restricted form of the AND-OR model [Knapp 1987], since each process has, in general, many possible moves at any time (the OR aspect), and to execute a move requires that all I -neighbors of the process allow the move to be executed (the AND aspect).¹¹ For a given I -system in a particular global state, the wait-for-graph characterizes the current blocking relations, i.e., which moves are blocked by which processes. We then characterize the occurrence of deadlocks as the occurrence within the wait-for-graph of a subgraph of a certain form that we call a “supercycle.” Essentially, a supercycle is a generalization to the AND-OR waiting model of what a cycle is in the AND model or a knot is in the OR model. Finally, we formulate an assumption on the behavior of pair-systems—the “wait-for-graph assumption”—that prevents the occurrences of supercycles, and hence of deadlocks, and formally establish that this is indeed the case.

6.5.1 An Example of a Deadlock-Prone I -System. Consider the pair-system with initial state set $\{[V_1 V_2]\}$ given in Figure 10. The variables x_{12}^2, y_{12}^2 are shared variables that are written by P_1 only and read by P_2 only. Shared variable x_{12}^2

⁹We do not distinguish between deadlock and termination in this article, and thus deadlock freedom and nontermination are expressed by the same CTL formula, namely $AGEXtrue$. It is in principle straightforward to make this distinction: for each $i \in \{i_1, \dots, i_K\}$, introduce a local proposition $term_i$ which is true if and only if process i has terminated. Deadlock-freedom is then written as $AG((term_{i_1} \vee \dots \vee term_{i_K}) \Rightarrow EXtrue)$, i.e., as long as some process has not terminated, some process can execute a transition. Nontermination, on the other hand, is expressed as $AG(\neg(term_{i_1} \vee \dots \vee term_{i_K}))$, i.e., no process terminates.

¹⁰We say an I -process P_i^I blocks a move of I -process P_j^I if and only if $i I j$ and there is an I -move a_j^I of P_j^I such that the current local state of P_j^I is the start state of a_j^I (i.e., local control is “at” a_j^I), and the conjunct of the guard of a_j^I that tests the shared state with P_i^I (i.e., $a_j^I.guard_i$) is false. In other words, P_j^I is prevented from executing a_j^I by P_i^I .

¹¹The general AND-OR model allows arbitrary combinations of “and” and “or” resource requests, whereas our model strictly alternates “and” and “or” requests, i.e., an I -process requests some move to execute, and that move in turn requests “permission” (true guards) from all the I -neighbors of the process.

controls the move of P_2 from W_2 to A_2 , which can be made only when $x_{12}^2 = 2$. Likewise, y_{12}^2 controls the move of P_2 from W_2 to B_2 . By making appropriate assignments to these variables, P_1 enables or disables the moves of P_2 out of W_2 (namely, $(W_2, x_{12}^2 = 2 \rightarrow \text{skip}, A_2)$ and $(W_2, y_{12}^2 = 2 \rightarrow \text{skip}, B_2)$) as follows. When P_1 moves from V_1 to W_1 , it either disables both these moves (if P_2 's current local state is V_2), or it non-deterministically chooses one of these moves to disable, and enables the other (if P_2 's current local state is W_2). P_2 exercises a symmetric control of the moves of P_1 out of W_1 (namely, $(W_1, x_{12}^1 = 1 \rightarrow \text{skip}, A_1)$ and $(W_1, y_{12}^1 = 1 \rightarrow \text{skip}, B_1)$) by assigning to the variables x_{12}^1, y_{12}^1 , which play a “bilaterally symmetric” role with respect to P_1 and P_2 that x_{12}^2, y_{12}^2 do. By inspection, the only potential deadlock state is when both processes are at W states and all outgoing moves are blocked, i.e., the state $[W_1 x_{12}^1 = 2 y_{12}^1 = 2 W_2 x_{12}^2 = 1 y_{12}^2 = 1]$. However we can easily verify that $(W_1 \wedge W_2) \Rightarrow (x_{12}^1 = 1 \vee y_{12}^1 = 1 \vee x_{12}^2 = 2 \vee y_{12}^2 = 2)$ is an invariant, so this state is not reachable. Hence the pair-system of Figure 10 is deadlock-free. The I -system for $I = \{\{1, 2\}, \{2, 3\}, \{3, 1\}\}$ derived from Figure 10 by the MP-synthesis definition (5.1.1) is shown in Figure 11. This system can reach a deadlocked state via the following path:

$$\begin{aligned}
& [V_1 V_2 V_3] \xrightarrow{1} \\
& [x_{13}^3 = 1 y_{13}^3 = 1 W_1 x_{12}^2 = 1 y_{12}^2 = 1 V_2 V_3] \xrightarrow{2} \\
& [x_{13}^3 = 1 y_{13}^3 = 1 W_1 x_{12}^2 = 1 y_{12}^2 = 1 x_{12}^1 = 2 y_{12}^1 = 1 W_2 x_{23}^3 = 2 y_{23}^3 = 2 V_3] \xrightarrow{3} \\
& [x_{13}^3 = 1 y_{13}^3 = 1 W_1 x_{12}^2 = 1 y_{12}^2 = 1 x_{12}^1 = 2 y_{12}^1 = 1 W_2 x_{23}^3 = 2 y_{23}^3 = 2 \\
& \quad x_{23}^2 = 3 y_{23}^2 = 2 W_3 x_{13}^1 = 1 y_{13}^1 = 3]
\end{aligned}$$

We can easily see by inspection that the last state (call it u^{dead}) of this path is a deadlock state, i.e., has no outgoing transitions: The P_1 -transition results in both moves out of W of both P_2, P_3 being disabled. The P_2 -transition results in the move $(W_1, x_{12}^1 = 1 \rightarrow \text{skip} \otimes x_{13}^1 = 1 \rightarrow \text{skip}, A_1)$ of P_1 being disabled. Finally, the P_3 transition results in the move $(W_1, y_{12}^1 = 1 \rightarrow \text{skip} \otimes y_{13}^1 = 1 \rightarrow \text{skip}, B_1)$ of P_1 being disabled. Hence, in state u^{dead} no move is enabled, and the system is deadlocked.

6.5.2 The Wait-For-Graph. We remind the reader of the following notation (see Section 5.4). A move a_i^I of P_i^I has the form $(s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$. $a_i^I.\text{start}, a_i^I.\text{guard}_j$ denote $s_i, \bigvee_{\ell \in [1:n]} B_{i,\ell}^j$, respectively; and $a_i^I.\text{guard}$ denotes $\bigwedge_{j \in I(i)} a_i^I.\text{guard}_j$. In the special case of $I = \{\{i, j\}\}$, a_i^I denotes the move $(s_i, \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$. Finally, we use $t_k.\text{moves}$ to denote the set of moves $\{a_k^I \mid a_k^I \in P_k^I \text{ and } a_k^I.\text{start} = t_k\}$.

The “wait-for-graph” $W_I(s)$ of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ at state s is a directed bipartite graph. The nodes of this graph are the I -processes P_i^I for $i \in \{i_1, \dots, i_K\}$, i.e., all the I -processes of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$, and the moves a_i^I of each P_i^I such that $s \models i = a_i^I.\text{start}$, i.e., all the moves that are candidates for execution. There is an edge in $W_I(s)$ from P_i^I to each node of the form a_i^I in $W_I(s)$. This edge indicates that a_i^I is a possibility for execution by P_i^I . Thus P_i^I is an “and” node, since P_i^I is blocked if and only if *all* of its successors (i.e., its possible moves) are blocked. Also, there is an edge in $W_I(s)$ from a_i^I to P_j^I if and only if $j \in I(i)$ and $s \models ij(a_i^I.\text{guard}_j) = \text{false}$, i.e., a_i^I is blocked by P_j^I (a_i^I can be executed in s only if $s \models ij(a_i^I.\text{guard}_j) = \text{true}$ for

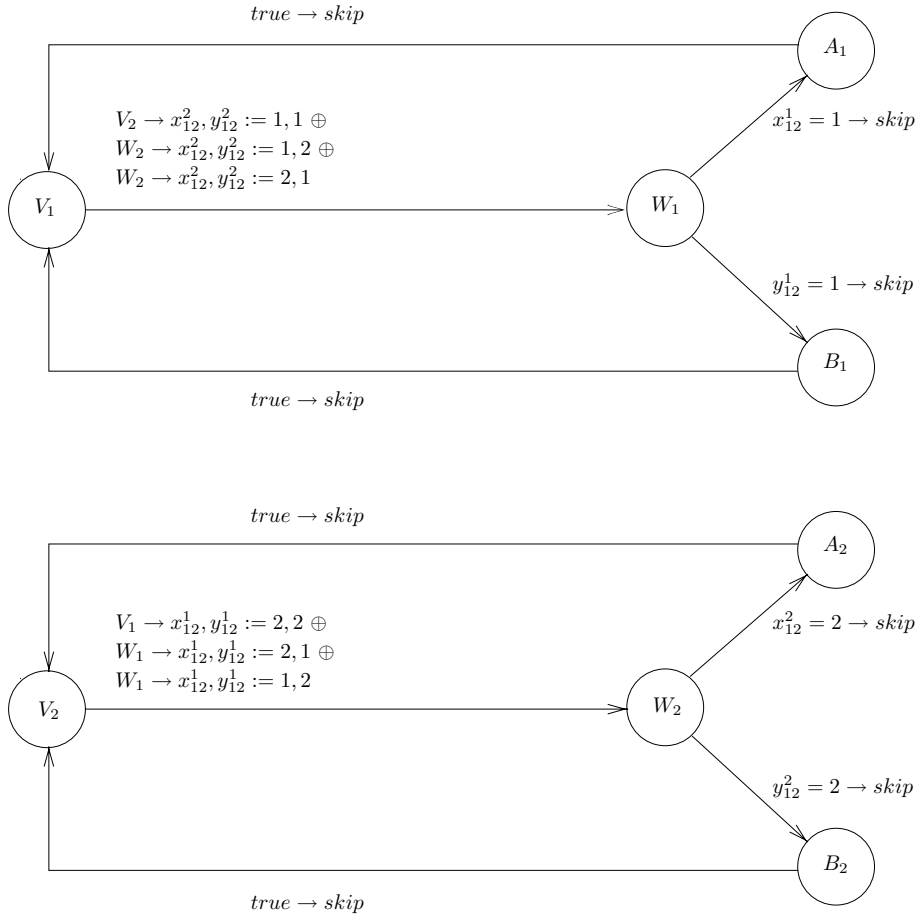


Fig. 10. Deadlock example for two processes.

all $j \in I(i)$. Thus, if there is some j in $I(i)$ such that $s \uparrow ij(a_i^I.guard_j) = false$, then a_i^I cannot be executed in state s (of M_I). Thus each move a_i^I is an “or” node, since a_i^I is blocked if and only if *some* neighbor P_j^I of P_i^I blocks a_i^I . Note, however, that we cannot say that P_i^I itself is blocked by P_j^I , since there could be another move b_i^I in P_i^I such that $s \uparrow ij(b_i^I.guard_j) = true$, i.e., b_i^I is not blocked by P_j^I (in state s), so P_i^I can progress in state s by executing b_i^I . We now give the formal definition.

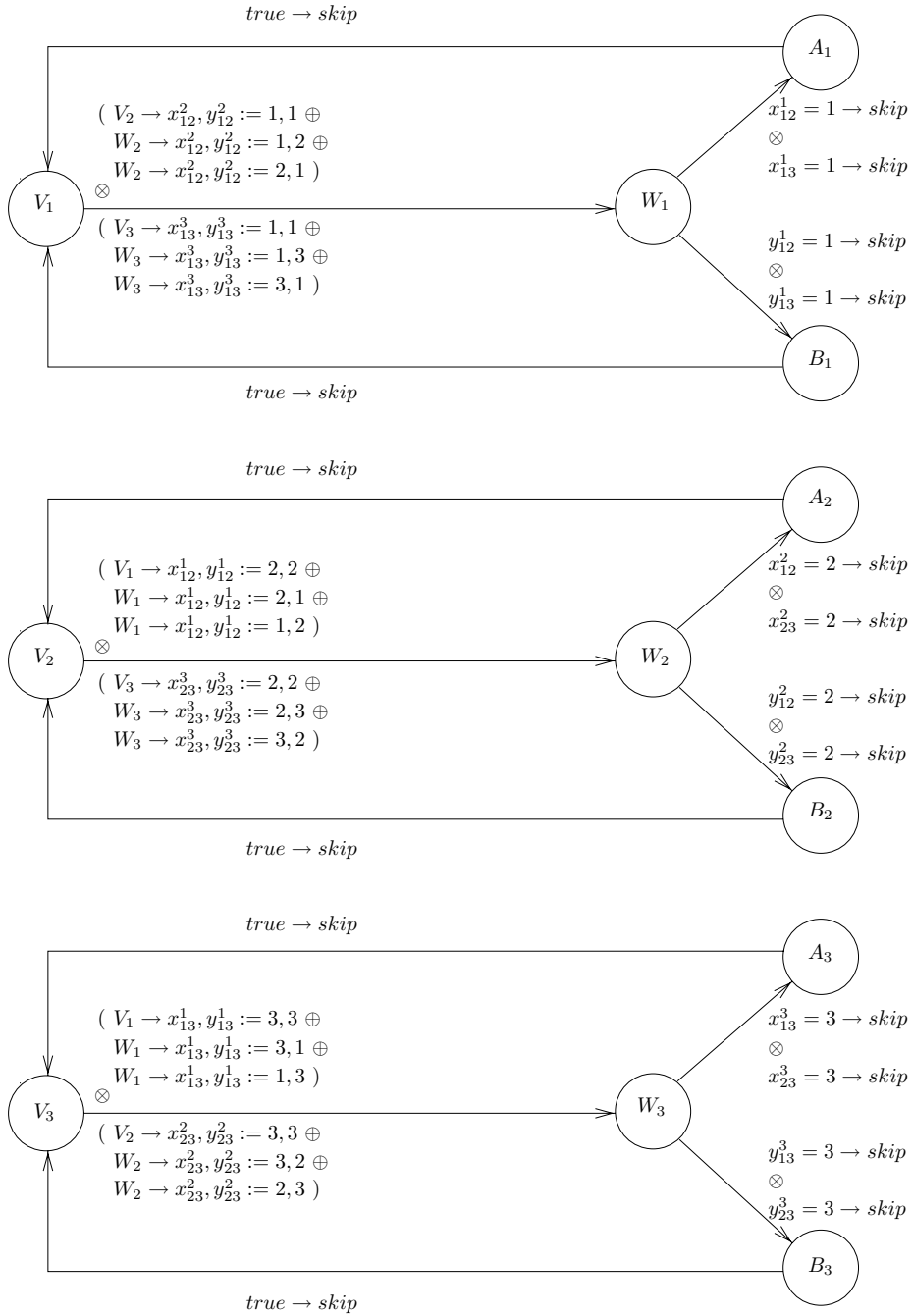


Fig. 11. Deadlock example for three processes.

Definition 6.5.2.1 (Wait-For-Graph $W_I(s)$). Let s be an arbitrary I -state. The *wait-for-graph* $W_I(s)$ of s is a directed bipartite graph, where

- (1) the nodes of $W_I(s)$ are
 - (a) the I -processes $\{P_i^I \mid i \in \text{dom}(I)\}$, and
 - (b) the moves $\{a_i^I \mid i \in \text{dom}(I) \text{ and } a_i^I \in P_i^I \text{ and } s \upharpoonright i = a_i^I.\text{start}\}$
- (2) there is an edge from P_i^I to every node of the form a_i^I in $W_I(s)$, and
- (3) there is an edge from a_i^I to P_j^I in $W_I(s)$ if and only if $i I j$ and $a_i^I \in W_I(s)$ and $s \upharpoonright j(a_i^I.\text{guard}_j) = \text{false}$.

In the sequel, we use $P_i^I \longrightarrow a_i^I \in W_I(s)$ to denote the existence of an edge from P_i^I to a_i^I in $W_I(s)$, and $a_i^I \longrightarrow P_j^I \in W_I(s)$ to denote the existence of an edge from a_i^I to P_j^I in $W_I(s)$. We also abbreviate $(P_i^I \longrightarrow a_i^I \in W(s) \text{ and } a_i^I \longrightarrow P_j^I \in W(s))$ with $P_i^I \longrightarrow a_i^I \longrightarrow P_j^I \in W(s)$, and similarly for longer “wait-chains.”

Observation 6.5.2.2 (Move Enablement). Let s be an arbitrary I -state such that $s \upharpoonright i = a_i^I.\text{start}$. If a_i^I has no outgoing edges in $W_I(s)$, then by the wait-for-graph definition (6.5.2.1), $s \upharpoonright j(a_i^I.\text{guard}_j) = \text{true}$ for all $j \in I(i)$. Hence, by the compact I -structure definition (5.4.3), a_i^I can be executed in state s .

6.5.3 Supercycles. We characterize a deadlock as the occurrence in the wait-for-graph of a graph-theoretic construct that we call a *supercycle*:

Definition 6.5.3.1 (Supercycle). SC is a supercycle in $W_I(s)$ if and only if all of the following hold:

- (1) SC is nonempty,
- (2) if $P_i^I \in SC$ then for all a_i^I such that $a_i^I \in W_I(s)$, $P_i^I \longrightarrow a_i^I \in SC$, and
- (3) if $a_i^I \in SC$ then there exists P_j^I such that $a_i^I \longrightarrow P_j^I \in W_I(s)$ and $a_i^I \longrightarrow P_j^I \in SC$.

Note that this definition implies that SC is a subgraph of $W_I(s)$.

If an I -process P_i^I is in a supercycle SC , then every potential move of P_i^I (i.e., every move out of the current local state of P_i^I) is also in SC and is blocked by some other I -process P_j that is also in SC . It follows that no I -process in SC is enabled. Furthermore this situation is stable, since every I -process P_i^I in SC is blocked by a set of I -processes (the I -processes that each block at least one potential move of P_i^I) that are all in SC . Thus, no I -process in SC can execute a transition, so the state of all these I -processes remains the same. Hence, the wait-for relationships defined by the supercycle persist forever.

Figure 12 shows the wait-for-graph $W_I(u^{\text{dead}})$ for the I -system of Figure 11 in its deadlocked state u^{dead} . And-nodes (processes) are shown as \bullet , while or-nodes (moves) are shown as \circ . $a_i, i \in \{1, 2, 3\}$ is the move of P_i from W_i to A_i , and $b_i, i \in \{1, 2, 3\}$ is the move of P_i from W_i to B_i . Since every move in $W_I(u^{\text{dead}})$ has at least one outgoing edge, i.e., is blocked by at least one process, $W_I(u^{\text{dead}})$ is also an example of a supercycle. In fact, several edges can be removed and still leave a supercycle (for example, $a_3 \longrightarrow P_1$, $b_3 \longrightarrow P_2$, $a_2 \longrightarrow P_1$ can all be removed). Thus, $W_I(u^{\text{dead}})$ contains several subgraphs that are also supercycles.

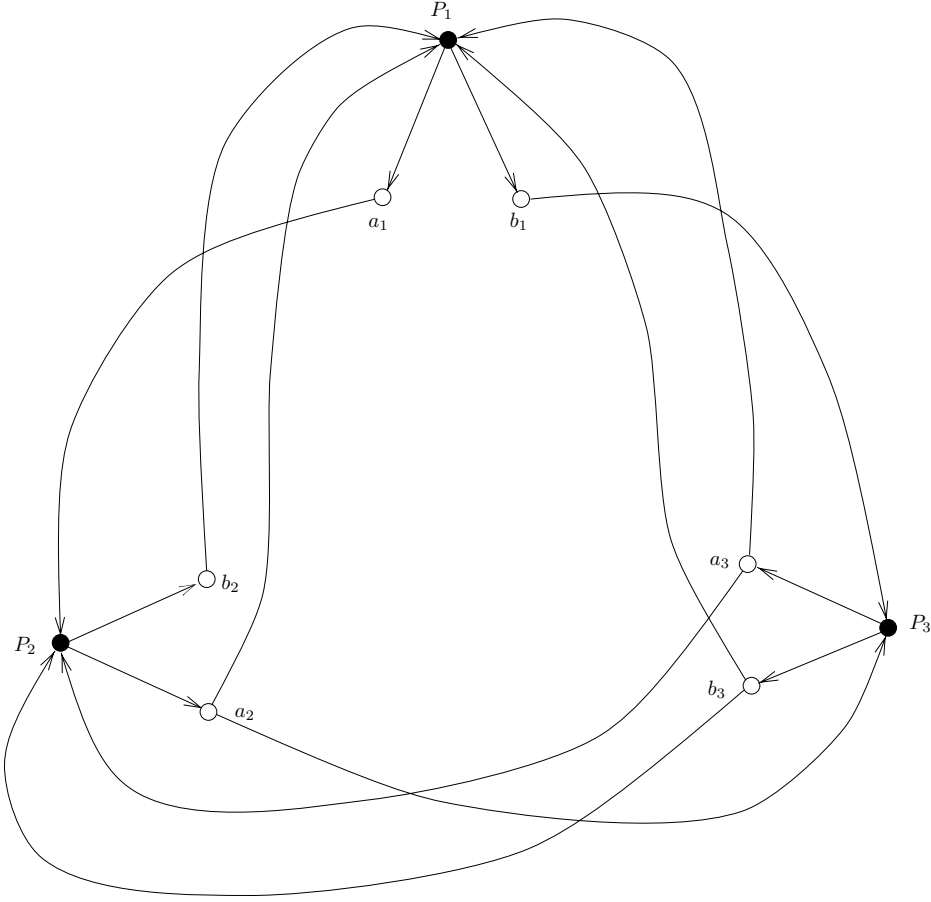


Fig. 12. Example wait-for-graph.

6.5.4 Establishing Supercycle-Freedom: The Wait-for-Graph Assumption. We shall establish deadlock freedom by showing that the wait-for-graph is always supercycle-free. We achieve this by making a technical assumption—the wait-for-graph assumption—which effectively precludes the formation of supercycles in the wait-for-graph. The essential idea is roughly as follows. The wait-for-graph assumption will allow us to conclude that in a state s resulting from a transition by P_k^I , either P_k^I has an enabled move, or P_k^I does not block any move by any other process (in neither case can P_k^I belong to a supercycle). Formally, this condition is

$$\begin{aligned}
 &\text{for every reachable } I\text{-state } t \text{ in } M_I \text{ such that } s \xrightarrow{k} t \in R_I \text{ for some} \\
 &\text{reachable } I\text{-state } s, \\
 &\quad \neg \bigvee a_j^I . (a_j^I \longrightarrow P_k^I \in W_I(t)) \quad \text{or} \\
 &\quad \bigvee a_k^I \in W_I(t) . (\bigwedge \ell \in \{i_1, \dots, i_K\} . (a_k^I \longrightarrow P_\ell^I \notin W_I(t))). \quad (**)
 \end{aligned}$$

The first disjunct says that P_k^I blocks no move by any other process, and the second

disjunct says that P_k^I has some move that is blocked by no process, i.e., some enabled move. Since our goal is to devise a completely mechanical synthesis method, the above condition must be mechanically checked. Doing so requires us, however, to generate M_I , which is, in general, exponentially large in the number of processes $K(=|dom(I)|)$. Instead, we take advantage of the state-mapping corollary (6.4.5), which relates the reachable states of M_I to the reachable states of M_J (for $J \subseteq I$), and the transition-mapping corollary (6.4.2), which relates R_I to R_J (for $J \subseteq I$), and the following proposition, which relates $W_I(s)$ to $W_J(s \uparrow J)$ (again for $J \subseteq I$), to formulate a condition over M_J (for a suitably small J) that implies condition (**). This will be the wait-for-graph assumption. If $a_i^I = (s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$, then we let a_i^J denote $(s_i, \otimes_{j \in J(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$.

PROPOSITION 6.5.4.1 (WAIT-FOR-GRAPH PROJECTION). *Let $J \subseteq I$ and $i \in J$. Furthermore, let s_I be an arbitrary I -state. Then*

- (1) $P_i^I \rightarrow a_i^I \in W_I(s_I)$ iff $P_i^J \rightarrow a_i^J \in W_J(s_I \uparrow J)$, and
- (2) $a_i^I \rightarrow P_j^I \in W_I(s_I)$ iff $a_i^J \rightarrow P_j^J \in W_J(s_I \uparrow J)$.

Now suppose that condition (**) is violated. Then, there is a transition $s \xrightarrow{k} t$ in R_I , where s is a reachable I -state and where

$$\bigvee a_k^I \in W_I(t) \cdot (a_j^I \rightarrow P_k^I \in W_I(t))$$

and

$$\bigwedge a_k^I \in W_I(t) \cdot (\bigvee \ell \in \{i_1, \dots, i_K\} \cdot (a_k^I \rightarrow P_\ell^I \in W_I(t))).$$

The first conjunct states that P_k^I has some edge incoming from some node a_j^I . The second conjunct states that for every move node of the form a_k^I in $W_I(t)$, there is some I -process P_ℓ^I that blocks a_k^I . Thus, if $n = |(t \uparrow k).moves|$, (i.e., n is the number of move nodes of the form a_k^I in $W_I(t)$), then by selecting one of the I -processes that blocks each move node of the form a_k^I , we can form a set $\{P_j^I, P_{\ell_1}^I, \dots, P_{\ell_n}^I\}$ of I -processes such that

$$a_j^I \rightarrow P_k^I \in W_I(t)$$

and

$$\bigwedge a_k^I \in W_I(t) \cdot (\bigvee \ell \in \{\ell_1, \dots, \ell_n\} \cdot (a_k^I \rightarrow P_\ell^I \in W_I(t))).$$

Now let $J = \{\{j, k\}, \{\ell_1, k\}, \dots, \{\ell_n, k\}\}$. By the state-mapping corollary (6.4.5) $s \uparrow J$ is a reachable J -state of M_J , and by the transition-mapping corollary (6.4.2), $s \uparrow J \xrightarrow{k} t \uparrow J \in R_J$. Thus, by the wait-for-graph projection proposition (6.5.4.1), there is a transition $s \uparrow J \xrightarrow{k} t \uparrow J$ in R_J , where $s \uparrow J$ is a reachable J -state and where

$$a_j^J \rightarrow P_k^J \in W_J(t \uparrow J)$$

and

$$\bigwedge a_k^J \in W_J(t \uparrow J) \cdot (\bigvee \ell \in \{\ell_1, \dots, \ell_n\} \cdot (a_k^J \rightarrow P_\ell^J \in W_J(t \uparrow J))). \quad (***)$$

It follows that if condition (***) does not hold for any transition of R_J whose initial state is reachable, then (**) cannot be violated. This leads us to the following formulation of the wait-for-graph assumption.

Definition 6.5.4.2 (Wait-For-Graph Assumption: WG). Let t_k be an arbitrary reachable local state of P_k^ℓ in $M_{k\ell}$, and let $n = |t_k.moves|$. Also let J be an

arbitrary interconnection relation of the form $\{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_n\}\}$, where $k \notin \{j, \ell_1, \dots, \ell_n\}$. Then, for every reachable J -state t_J in M_J such that $t_J \uparrow k = t_k$ and $s_J \xrightarrow{k} t_J \in R_J$ for some reachable J -state s_J , we have

$$\bigwedge a_j^J . (a_j^J \longrightarrow P_k^J \notin W_J(t_J))$$

or

$$\bigvee a_k^J \in W_J(t_J) . (\bigwedge \ell \in \{\ell_1, \dots, \ell_n\} . a_k^J \longrightarrow P_\ell^J \notin W_J(t_J)).$$

Note that, in the above discussion, ℓ_1, \dots, ℓ_n are not necessarily pairwise distinct, i.e., an I -process P_ℓ^I could block more than one move of the form a_k^I . Thus there are n cases for the form of J , as the number m of distinct members of $\{\ell_1, \dots, \ell_n\}$ varies from 1 to n . All of these cases must be considered. Likewise, the above discussion does not specify whether j is a member of $\{\ell_1, \dots, \ell_n\}$ or not. Thus both cases must be considered, which in effect generates two subcases for each of the n cases outlined above. Appendix E presents an algorithm for mechanically checking the wait-for-graph assumption. We now use the wait-for-graph assumption to establish supercycle-freedom of the wait-for-graph of any reachable state.

THEOREM 6.5.4.3 (SUPERCYCLE-FREE WAIT-FOR-GRAPH). *If the wait-for-graph assumption WG holds, and $W_I(s_I^0)$ is supercycle-free for every initial state $s_I^0 \in S_I^0$, then for every reachable state t of M_I , $W_I(t)$ is supercycle-free.*

6.5.5 Establishing Deadlock-freedom. We show that the absence of supercycles in the wait-for-graph of a state implies that there is at least one enabled move in that state.

PROPOSITION 6.5.5.1 (SUPERCYCLE). *If $W_I(s)$ is supercycle-free, then some move a_i^I has no outgoing edges in $W_I(s)$.*

PROOF. We establish the contrapositive. Since every local state of a process has at least one outgoing arc (Section 2), there exists at least one move of the form a_i^I for every $i \in \text{dom}(I)$ in $W_I(s)$. Suppose that every such move has at least one outgoing edge in $W_I(s)$. Consider the subgraph SC of $W_I(s)$ consisting of these edges together with all edges of the form $P_i^I \longrightarrow a_i^I$ in $W_I(s)$. By the wait-for-graph definition (6.5.2.1), and the supercycle definition (6.5.3.1), it is clear that SC is a supercycle in $W_I(s)$. Thus $W_I(s)$ is not supercycle-free. \square

Finally, we apply the supercycle proposition (6.5.5.1) to all of the reachable states of M_I , thereby establishing the deadlock freedom of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$.

THEOREM 6.5.5.2 (DEADLOCK FREEDOM). *If, for every reachable state s of M_I , $W_I(s)$ is supercycle-free, then $M_I, S_I^0 \models AGEX\text{true}$.*

PROOF. Let s be an arbitrary reachable state of M_I . By the antecedent, $W_I(s)$ is supercycle-free. Hence, by the supercycle proposition (6.5.5.1), some move a_i^I has no outgoing edges in $W_I(s)$. By Observation 6.5.2.2, a_i^I can be executed in state s . Since s is an arbitrary reachable state of M_I , we conclude that every reachable state of M_I has at least one enabled move a_i^I , (where, in general, a_i^I depends on s). Hence $M_I, S_I^0 \models AGEX\text{true}$. \square

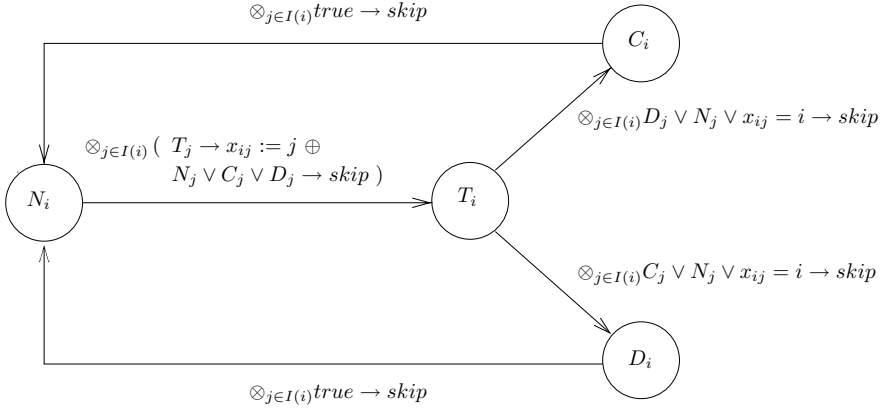


Fig. 13. Mutual exclusion example with two critical sections.

6.5.6 Examples. We now turn to some examples. Consider the solution for the generalized dining philosophers problem given in Figure 7. We verify that this solution satisfies the wait-for-graph assumption. The pair-process P_k^ℓ from which the solution in Figure 7 is derived has three local states, which, by slight abuse of terminology, we refer to as $[N_k]$, $[T_k]$, and $[C_k]$. We deal with each local state in turn.

For the move from $[N_k]$ to $[T_k]$, we observe, by inspection, that P_k^I blocks no move by any other process in the resulting I -state. For the moves from $[T_k]$ to $[C_k]$ and $[C_k]$ to $[N_k]$, we see that, in both cases, P_k^I has an enabled move in the resulting I -state.

Having checked WG for every (reachable) local state of P_k^ℓ , we conclude that the solution of Figure 7 satisfies WG (note that the above analysis takes care of both the subcases $j = \ell_1, j \neq \ell_1$). It remains to show that $W_I(s_I^0)$ is supercycle-free for every initial state s_I^0 . In an initial state, all processes are in their noncritical (“ $[N]$ ”) state, and no process blocks a move of any other process. Hence there can be no supercycles. Therefore, by the above results, we infer that every I -system corresponding to the solution in Figure 7 is deadlock-free.

Furthermore, consider the generalized example given in Figure 13 where there are now two critical sections: $[C_i]$ and $[D_i]$. Processes request entry into either of these critical sections, after which they return to their noncritical sections. When I is a “complete graph,” this example (which can clearly be generalized to an arbitrary number of critical sections) models the situation where we have some number of identical resources and where processes require exactly one of these resources. Such a resource could be a printer, for example, where a process is satisfied by acquiring any one of some number of printers. Again we can verify by simply checking all the possible moves of P_k^J that this example satisfies the wait-for-graph assumption, and therefore, all the I -systems generated from it are deadlock-free.

Looking back at the system of Figure 10, we see that it violates the wait-for-graph assumption in that it allows transitions by a process such that in the resulting

state, it blocks a move of some other process, and has all its own moves blocked by other processes. The transition by P_3 in the path displayed in Section 6.5.1 is an example of such a transition: in the resulting state u^{dead} , P_3 blocks the move $(W_1, b_{12}^1 = 1 \rightarrow skip \otimes b_{13}^1 = 1 \rightarrow skip, B_1)$ of P_1 , while both its own moves are blocked by P_2 (and P_1 , for that matter).

6.6 The Large-Model Theorem

As stated in the introduction, a crucial aspect of MP-synthesis is its soundness: what correctness properties of the pair-system are preserved by the I -system? Here, we treat the special case of complete interconnection: the process interconnection relation I is the “complete graph” on $\{i_1, \dots, i_K\}$. The general case is treated below in Section 6.8. Our soundness result for the case of complete interconnection is expressed as the large-model theorem. We show that propositional invariants expressed pairwise, temporal leads-to properties over individual processes, and local structure properties are all preserved (see Section 6.1). To facilitate the formal statement of our results, we define the sublogic “Flat CTL” (FLCTL) of CTL.

Definition 6.6.1 (Flat CTL). The logic $FLCTL_{k\ell}$ is the set of state formulae that are finite conjunctions of the following, where $h_{k\ell}$ is a formula of $\mathcal{L}(\mathcal{AP}_k, \mathcal{AP}_\ell, \neg, \wedge)$, and p_k, q_k, r_k are formulae of $\mathcal{L}(\mathcal{AP}_k, \neg, \wedge)$:

- $h_{k\ell}$
- $AGh_{k\ell}$
- $AG(p_k \Rightarrow A[q_k Ur_k])$
- $AG(p_k \Rightarrow AY_k q_k)$
- $AG(a_k \Rightarrow EX_k b_k)$, where $a_k, b_k \in \{\llbracket s_k \rrbracket \mid s_k \in S_k\}$

In terms of our notation for sublogics (Section 3.2), $FLCTL_{k\ell}$ is the sublogic $\mathcal{L}(\wedge, h_{k\ell}, AGh_{k\ell}, AG(p_k \Rightarrow A[q_k Ur_k]), AG(p_k \Rightarrow AY_k q_k), AG(a_k \Rightarrow EX_k b_k))$.

THEOREM 6.6.2 (LARGE MODEL). *Let $f_{k\ell}$ be an arbitrary formula of $FLCTL_{k\ell}$. Let $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$. Furthermore, let s be an arbitrary reachable I -state, and, for all i, j such that $i I j$, let $s_{ij} = s \upharpoonright ij$. If $M_I, S_I^0 \models AGEX true$, then*

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models \mathbf{\Lambda}_{k\ell} f_{k\ell}) \text{ implies } M_I, s \models \mathbf{\Lambda}_{k\ell} f_{k\ell}.$$

Note that the formula satisfied by the I -system depends (by $MPCTL^*$ semantics) on I itself, in that the quantification over pairs given by $\mathbf{\Lambda}_{k\ell}$ has as its range all pairs in I . Thus, we cannot directly conclude, from the large-model theorem, any correctness property that relates two processes which are not I -neighbors. Such a property could, however, be a consequence of pairwise expressible properties that relate, say, a common neighbor of the non-neighboring pair (thereby imposing some indirect interaction on the pair). The investigation of methods for deducing such indirect correctness properties is outside the scope of this article and is a topic of future research.

Our notion of a correctness property is one that holds in all the initial states of the pair-system or I -system (Section 6.1). The following corollary presents the soundness result in terms of such correctness properties.

COROLLARY 6.6.3 (LARGE MODEL). *Let $f_{k\ell}$ be an arbitrary formula of $FLCTL_{k\ell}$. Let $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$, and let (i, j) be an arbitrary pair in I . If $M_I, S_I^0 \models AGEXtrue$, then*

$$M_{ij}, S_{ij}^0 \models \bigwedge_{k\ell} f_{k\ell} \text{ implies } M_I, S_I^0 \models \bigwedge_{k\ell} f_{k\ell}.$$

6.7 Liveness Properties

We wish to generalize the large-model theorem by dropping the assumption that the process interconnection scheme is a complete graph. This enables us to deal with, for example, the dining philosophers and readers-writers problems. Unfortunately, generalizing the interconnection relation I to arbitrary schemes introduces the possibility that liveness properties may be violated. In this article, we restrict our attention to the correctness properties expressible in Flat CTL (Definition 6.6.1). The only liveness property expressible in Flat CTL is $AG(p_i \Rightarrow A[q_i Ur_i])$, where p_i, q_i, r_i are propositional formulae over the atomic propositions of \mathcal{AP}_i . Now $AG(p_i \Rightarrow A[q_i Ur_i])$ is equivalent to $AG(p_i \Rightarrow A[q_i U_w r_i]) \wedge AG(p_i \Rightarrow AFr_i)$. Since $AG(p_i \Rightarrow A[q_i U_w r_i])$ is not a liveness property, in the sequel we will use $AG(p_i \Rightarrow AFr_i)$ as our prototypical liveness property. A liveness property of this form is sometimes referred to as a “temporal leads-to” property in the literature, since its intuitive meaning is “if p_i is true in some global state, then r_i is guaranteed to eventually be true along every fullpath starting in this state.” In other words, p_i “leads to” r_i . The formula AFr_i is an *eventuality* formula. Suppose $M_{ij}, s_{ij} \models AFr_i$ for some ij -state s_{ij} . By CTL* semantics, we have the equivalence $AFr_i \equiv r_i \vee AXAFr_i$. Thus, if $s_{ij} \not\models r_i$, then we have $M_{ij}, s_{ij} \models AXAFr_i$, i.e., for every successor state t_{ij} of s_{ij} , we have $M_{ij}, t_{ij} \models AFr_i$. Thus, the eventuality AFr_i has been *propagated* to the successors of s_{ij} . This propagation continues as long as $\neg r_i$ holds. In this case we say that AFr_i is *pending* at state s_{ij} . In other words, AFr_i is a pending eventuality of s_{ij} if and only if r_i is not true in s_{ij} , but is guaranteed to eventually become true along every fullpath starting in s_{ij} . When an ij -state is reached where r_i is true, AFr_i is said to be *fulfilled*. Note that if $s_{ij} \models r_i$, then $M_{ij}, s_{ij} \models AFr_i$ follows immediately, and AFr_i is fulfilled in s_{ij} itself.

In an I -system, pending eventualities may be prevented from being fulfilled in two ways. The first way is due to the particular blocking behavior inherent in the pair-system that is input to MP-synthesis. The solution here is to make a technical assumption on the pair-system—the liveness assumption—that prevents this kind of blocking behavior. The second way is when the I -system is “dynamically partitioned” into two subsystems such that, from some point onward, no process in one subsystem ever blocks a move of any process in the other subsystem. In this case, it is possible to schedule processes only from one of the subsystems. We deal with this by requiring the use of fair scheduling, and define the particular fairness that we require—weak blocking fairness—below. We establish a progress lemma that states (roughly) that under weak-blocking fairness, and assuming the liveness-assumption, every I -process that must progress in order to fulfill a pending eventuality actually does progress. This lemma is crucial in establishing the generalized large-model theorem (Section 6.8), which in turn allows us to conclude that MP-synthesis preserves temporal leads-to properties (e.g., if a philosopher requests

entry to the critical section, it eventually gains entry).

6.7.1 Example of Nonfulfillment of Pending Eventualities Due to Blocking Behavior. Consider the I -system synthesized using the interconnection relation $I = \{\{1, 2\}, \{2, 3\}, \{3, 4\}, \{4, 5\}\}$ and the pair-structure given in Figure 14. Suppose $\mathbf{\Lambda}_k AG(D_k \Rightarrow AFE_k)$ is a conjunct of the problem specification.¹² It is easily seen by inspection that $M_{ij}, S_{ij}^0 \models \mathbf{\Lambda}_k AG(D_k \Rightarrow AFE_k)$. We show however, that $M_I, S_I^0 \not\models \mathbf{\Lambda}_k AG(D_k \Rightarrow AFE_k)$. In particular, consider the following path π in M_I , which starts in the initial state $[A_1 A_2 A_3 A_4 A_5]$:

$$\begin{aligned} & [A_1 A_2 A_3 A_4 A_5] \xrightarrow{1} [B_1 A_2 A_3 A_4 A_5] \xrightarrow{5} [B_1 A_2 A_3 A_4 B_5] \xrightarrow{5} \\ & [B_1 A_2 A_3 A_4 X_5] \xrightarrow{4} [B_1 A_2 A_3 B_4 X_5] \xrightarrow{3} [B_1 A_2 B_3 x_{34}=4 B_4 X_5] \xrightarrow{2} \\ & [B_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 X_5] \xrightarrow{1} [C_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 X_5] \xrightarrow{1} \\ & ([D_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 X_5] \xrightarrow{5} \\ & [D_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 Y_5] \xrightarrow{5})^\omega \end{aligned}$$

Let $s = [D_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 X_5]$, and $t = [D_1 x_{12}=1 B_2 x_{23}=3 B_3 x_{34}=4 B_4 Y_5]$. Clearly, $\pi \models \neg FE_1$, so $M_I, s \models \neg AFE_1$. But $s \models D_1$, so $M_I, s \models \neg(D_1 \Rightarrow AFE_1)$. Since s is a reachable state, we have $M_I, S_I^0 \not\models AG(D_1 \Rightarrow AFE_1)$. Hence, by MPCTL semantics, we finally conclude $M_I, S_I^0 \not\models \mathbf{\Lambda}_k AG(D_k \Rightarrow AFE_k)$.

6.7.2 The Liveness Assumption. The standard solution to the problem of ensuring that all pending eventualities are fulfilled is to adopt some form of fair scheduling notion [Emerson and Lei 1985; Francez 1986]. However, in this situation, no asymptotic fairness notion can help. Let ρ be the suffix $(s \xrightarrow{5} t)^\omega$ of π . By inspection, we see that only P_5 is enabled in any state along ρ . Furthermore, every state of ρ has only a single successor, namely the next state along ρ , i.e., once the first state of ρ has been entered, it is not possible to reach any state of M_I other than a state along ρ . Thus, no asymptotic fairness notion can preclude ρ as a possible computation. One would need a fairness notion which incorporated “lookahead” and prevented execution from reaching a state lying on an “unfair” path such as ρ . Since such fairness notions have not been investigated thoroughly, and in any case seem to be expensive to implement, we avoid this approach. Our only apparent alternatives are to make suitable assumptions on the structure of I and/or M_{ij} .

Since the whole point is to generalize the large-model theorem so that it applies to an arbitrary interconnection relation I , we avoid making any restrictions on I . Thus, we try to restrict the structure of M_{ij} in some way. Now the general situation which the above example illustrates is where $I = \{\{i, j_1\}, \{j_1, j_2\}, \dots, \{j_{n-1}, j_n\}, \{j_n, k\}\}$ (for arbitrary n), and there is a reachable state s of M_I such that

- (1) an eventuality of the form AFr_i is pending in s and
- (2) there is a wait-chain from P_i^I to P_k^I , and only P_k^I can be executed from some point onward.

¹²Since AFE_k is an abbreviation for $A[trueUE_k]$, this is within the fragment of CTL specifications that are preserved by our method.

We require an assumption that prevents the occurrence of such a state, i.e., all reachable states must satisfy the negation of one (or both) of the two clauses given above. We use the following:

$$M_{j_1 k}, s \uparrow j_1 k \not\models EGex_k \text{ or } \bigwedge a_i^I . (P_i^I \longrightarrow a_i^I \longrightarrow P_{j_1}^I \notin W_I(s)).$$

This is the negation of clause 2 for $I = \{\{i, j_1\}, \{j_1, k\}\}$ (where $i = k$ and $i \neq k$ are both possible, i.e., both cases must be considered). Since clause 2 implicitly universally quantifies over all I of the given form $(\{\{i, j_1\}, \{j_1, j_2\}, \dots, \{j_{n-1}, j_n\}, \{j_n, k\}\})$, it suffices to assume the negation for a subset of the possible values for I . It is imperative to use a small I to make the mechanical evaluation of our assumption tractable, since this evaluation requires M_I to be constructed, and $|M_I|$ grows exponentially with $|I|$. We formally define the required condition as the “liveness assumption.” Note that the term $P_i^I \longrightarrow a_i^I$ can be dropped without changing the meaning, since $P_i^I \longrightarrow a_i^I \notin W_I(s)$ implies $a_i^I \longrightarrow P_{j_1}^I \notin W_I(s)$.

Definition 6.7.2.1 (Liveness Assumption: LV). Let J be of the form $\{\{i, j\}, \{j, k\}\}$ where $j \notin \{i, k\}$. Then, for every reachable state s_J in M_J ,

$$M_{j k}, s_J \uparrow j k \models EGex_k \text{ implies } \bigwedge a_i^J . (a_i^J \longrightarrow P_j^J \notin W_J(s_J)).$$

Appendix E shows how the liveness assumption can be mechanically checked using CTL model-checking [Clarke et al. 1986].

6.7.3 Fairness. Notwithstanding our comments above about the futility of using asymptotic fairness to deal with the example in Figure 14, we nevertheless demonstrate that some form of fairness is essential if we are to avoid restrictions on the interconnection relation I .¹³ Suppose I is partitioned into subrelations I' and I'' such that no process in (the domain of) I' is I -related to a process in the domain of I'' , and vice-versa. Thus I' and I'' represent “uncoupled” systems. Assuming both systems are deadlock-free, it is possible, under nondeterministic scheduling, to always select processes from I' for execution, thereby causing nonfulfillment of pending eventualities of processes in I'' . Even if I is not partitioned in this way, it is possible for such uncoupled subsets of I to arise dynamically during particular executions. As a concrete example of this, consider the system generated by MP-synthesis for the dining philosophers specification given in Section 3.4, i.e., five philosophers P_1, P_2, P_3, P_4, P_5 arranged in a ring. Consider the following path π , which starts in the initial state $s^0 = [N_1 N_2 N_3 N_4 N_5]$:

$$[N_1 N_2 N_3 N_4 N_5] \xrightarrow{1} ([T_1 N_2 N_3 N_4 N_5] \xrightarrow{3} [T_1 N_2 T_3 N_4 N_5] \xrightarrow{3} [T_1 N_2 C_3 N_4 N_5] \xrightarrow{3})^\omega$$

Under purely nondeterministic scheduling π is an initialized fullpath of (the I -structure of) the solution produced by MP-synthesis for the above problem. (For ease of notation, we have omitted the values of the shared variables in π . The reader can easily ascertain, by inspecting Figure 6, that appropriate values exist.) Now π does not satisfy $G(T_1 \Rightarrow FC_1)$, and hence s^0 does not satisfy $AG(T_1 \Rightarrow AFC_1)$, which is a conjunct of the dining philosophers specification. The problem here is that P_1 is enabled, but fails to make progress (and fulfill the pending eventuality

¹³With the exception of our assumption (see Section 3.3) that every process in $dom(I)$ is I -related to at least one other process.

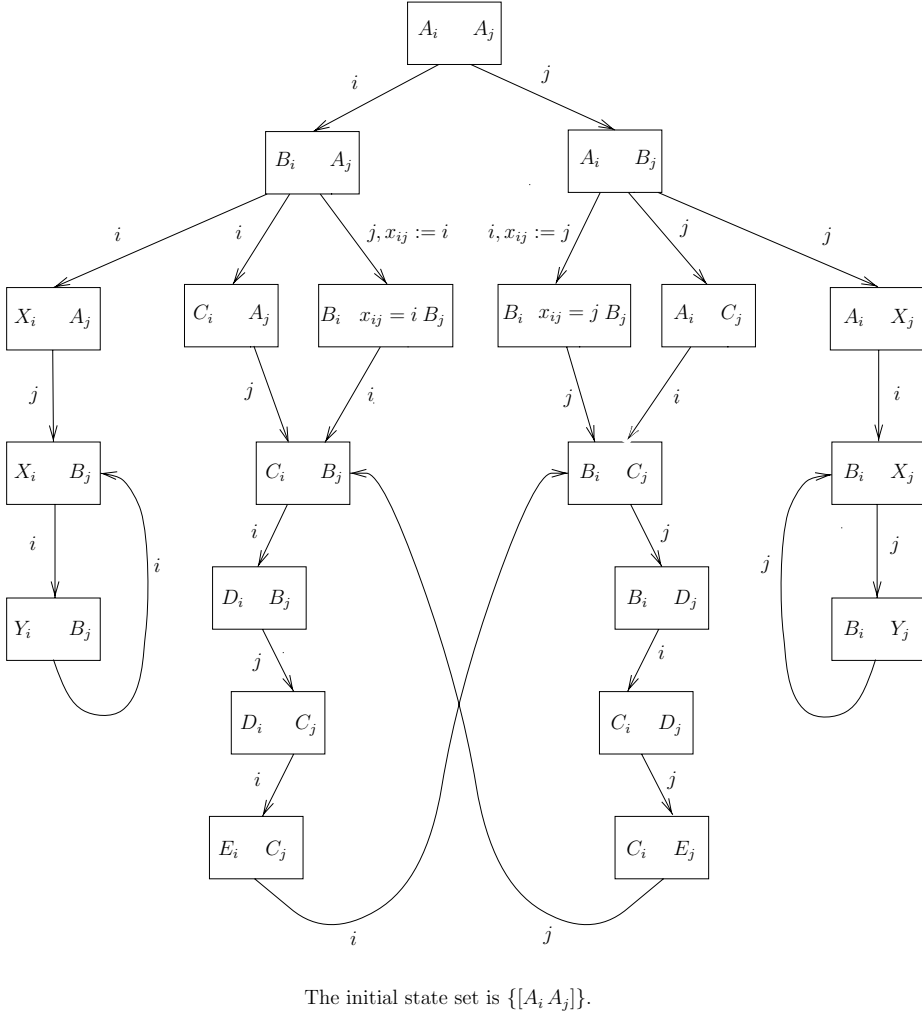


Fig. 14. Pair-structure illustrating nonfulfillment of a pending eventuality.

AFC_1) because it is not scheduled for execution. Since P_1 has become “decoupled” from P_3 (due to the other philosophers remaining permanently in their neutral (N) states), it is possible, under nondeterministic scheduling, for P_3 to be scheduled for execution continuously in spite of P_1 being continuously enabled. Note that this scenario cannot arise when I is a complete graph, because then all pairs of processes would be neighbors, so no pair of processes could ever become decoupled in the manner illustrated above.

In the dining philosophers system, fairness would eventually require that P_1 be selected for execution, and thus the path π exhibited above would no longer be a possible computation. Let en_i, ex_i mean that P_i is enabled in a state, executed along a transition, respectively (note that en_i is rendered as $EX_i true$ in CTL^*).

The three most common fairness notions are

- unconditional fairness: $\bigwedge_{i \in \text{dom}(I)} \overset{\infty}{F}ex_i$
- weak fairness: $\bigwedge_{i \in \text{dom}(I)} \overset{\infty}{G}en_i \Rightarrow \overset{\infty}{F}ex_i$
- strong fairness: $\bigwedge_{i \in \text{dom}(I)} \overset{\infty}{F}en_i \Rightarrow \overset{\infty}{F}ex_i$

Unconditional fairness is undesirable, since it requires that every process be infinitely often enabled, and this may not be the case, in general. We therefore adopt (a form of) weak fairness that is sufficient for our needs, and easier to implement than strong fairness. In the previous example, under weak fairness, P_1 is eventually executed, as desired. Unfortunately, P_2 is also eventually executed. This is not so desirable, as P_2 is in its noncritical state, and therefore weak fairness is forcing P_2 out of its noncritical state and into its trying state, i.e., it is forcing P_2 to make a request for critical section entry. In this example, we want the fairness to apply only to the trying states, where a request for critical section entry has already been made, and not to noncritical states, where it forces “spurious” requests. In addition, fairness must also apply to the critical states, in order to avoid indefinitely tying up critical resources, (which would prevent other processes from gaining access), and thereby guarantee that all eventualities are met. In the dining philosophers example, if fair scheduling is not applied to P_i in state s such that $s \uparrow i = C_i$, $s \uparrow (i+1) = T_{i+1}$, then (even though s has no pending eventuality for P_i) the pending eventuality of AFC_{i+1} for P_{i+1} might never be fulfilled, as P_i might remain forever in C_i , preventing the entry of P_{i+1} into C_{i+1} .

Generalizing this discussion to an arbitrary MPCTL* specification and the I -system derived from it, suppose that $i I j$ and $j I k$, and there is some reachable I -state s in M_I such that

- P_j^I has a pending eventuality in s ,
- in j -state $s \uparrow j$, P_j^I has two possible moves a_j^I and b_j^I (so $a_j^I.start = s \uparrow j$ and $b_j^I.start = s \uparrow j$),
- $s \uparrow i j \not\models a_j^I.guard_i$, i.e., a_j^I is blocked by P_i^I , and
- $s \uparrow j k \not\models b_j^I.guard_k$, i.e., b_j^I is blocked by P_k^I .

Clearly, if fair scheduling is not applied to P_i^I and P_k^I , then P_j^I 's pending eventuality may never be fulfilled. Note however, that $s \uparrow i j \models b_j^I.guard_i$ and $s \uparrow j k \models a_j^I.guard_k$ are both consistent with the previous example, i.e., it is not necessary for any one neighbor of P_j^I to block all the possible moves of P_j^I for P_j^I to be blocked. We must therefore apply fair scheduling to every I -process that is blocking at least one move of one of its neighbors. Furthermore we wish our fairness notion to depend only on the local state ($s \uparrow i$) of processes, and not on the shared variables (\mathcal{SH}_{ij} , for $i I j$) or the local state $s \uparrow j$ of any neighbor P_j^I . This makes the implementation of the fairness easier, since, for P_i^I , only its local state $s \uparrow i$ need be detected, rather than $\bigcup_{j \in I(i)} s \uparrow i j$ if shared variables and the local states of neighbors were included in the fairness notion. The latter involves the combined shared state over all of P_i^I 's neighbors. Detecting this combined state instantaneously is quite expensive, and we avoid the problem by formulating a fairness notion over local states only.

Hence, we must apply fair scheduling to every i -state s_i that is the P_i -component of some I -state s such that $s \upharpoonright ij \not\models a_j^I.guard_i$ for some move a_j^I of P_j^I . Note that since $a_j^I.guard_i$ depends on the shared variables as well as the current i -state, there may be other I -states t also having P_i -component s_i where $t \upharpoonright ij \models a_j^I.guard_i$. Thus, in some cases, the fair scheduling is not really necessary, i.e., our fairness is suboptimal. Making our fairness notion optimal would require instantaneous detection of the shared state with all I -neighbors, which, as mentioned above, would be quite expensive. Thus, we sacrifice some optimality in fair scheduling in return for a fairness notion that can be implemented reasonably efficiently.

Consider a move a_j^I of P_j^I . By the compact MP-synthesis definition (5.4.1), a_j^I has the form $(s_i, \otimes_{k \in I(j) \oplus \ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$. Hence, $a_j^I.guard_i = a_j^i.guard$. Hence, $s \upharpoonright ij \not\models a_j^I.guard_i$ if and only if $s \upharpoonright ij \not\models a_j^i.guard$. By the state-mapping corollary (6.4.5), we know that $s \upharpoonright ij$ is reachable in M_{ij} if s is reachable in M_I . Hence, we conclude that $s \upharpoonright ij \not\models a_j^I.guard_i$ can be evaluated by constructing M_{ij} only (and not M_I). This discussion leads us to the following definition.

Definition 6.7.3.1 (*Sometimes-Blocking State, blk_i*). An i -state s_i is *sometimes-blocking* if and only if:

$$\bigvee s_{ij}^0 \in S_{ij}^0 \cdot (M_{ij}, s_{ij}^0 \models EF(\{s_i\} \wedge (\bigvee a_j^i \in P_j^i \cdot (\{a_j^i.start\} \wedge \neg a_j^i.guard)))).$$

Note that $\bigvee a_j^i \in P_j^i \cdot (\{a_j^i.start\} \wedge \neg a_j^i.guard)$ is a purely propositional formula.

$$blk_i \stackrel{\text{df}}{=} (\bigvee \{s_i\} : s_i \text{ is sometimes-blocking}).$$

Thus, s_i is sometimes-blocking if and only if some (not necessarily all) state s_{ij} in M_{ij} blocks some move a_j^i of P_j^i and has s_i as its P_i -component. Note that it is entirely possible for M_{ij} to contain another state t_{ij} such that $t_{ij} \upharpoonright i = s_i$ and t_{ij} does not block a_j^i . Thus, being in a sometimes-blocking state does not imply that a neighbor's move is blocked: it only indicates (as the name suggests) the possibility that a neighbor's move is blocked. Also, being in a state that is not sometimes-blocking implies that no move of any neighbor is blocked. blk_i is a propositional formula (over the propositions in \mathcal{AP}_i) which is true in ij -state s_{ij} (respectively, global state s) exactly when $s_{ij} \upharpoonright i$ (respectively, $s \upharpoonright i$) is sometimes-blocking. If an i -state s_i is not sometimes-blocking, we say that it is *always-nonblocking*, and we define $nblk_i \stackrel{\text{df}}{=} \neg blk_i$.

For example, in the pair-structure (of the two-process mutual exclusion system) of Figure 4, $[N_1]$ is an always-nonblocking state: it occurs in three global states, namely $[N_1 N_2]$, $[N_1 T_2]$, and $[N_1 C_2]$, and all these states have an outgoing P_2^1 -transition. On the other hand, $[T_1]$ is a sometimes-blocking state, since the global state $[T_1 x_{12} = 1 T_2]$ has no outgoing P_2^1 -transition (but note that the global state $[T_1 x_{12} = 2 T_2]$ does have an outgoing P_2^1 -transition). $[C_1]$ is also a sometimes-blocking state, since $[C_1 T_2]$ has no outgoing P_2^1 -transition.

Observation 6.7.3.2. By Definition 6.7.3.1, if s_i is always-nonblocking, then $M_{ij}, S_{ij}^0 \models AG(\{s_i\} \Rightarrow \bigwedge a_j^i \in P_j^i \cdot (\{a_j^i.start\} \Rightarrow a_j^i.guard))$. This implies that the wait-for-graph $W_I(s)$ contains no edge whose target is P_i^I when $s \upharpoonright i$ is always-nonblocking.

Since we want fairness to apply only to sometimes-blocking states, we relativize

weak fairness to the sometimes-blocking states by defining our fairness notion Φ as a CTL* path formula as follows:

Definition 6.7.3.3 (Weak Blocking Fairness Φ).

$$\Phi \stackrel{\text{df}}{=} \bigwedge_{i \in \text{dom}(I)} \tilde{G}(\text{blk}_i \wedge \text{en}_i) \Rightarrow \tilde{F} \text{ex}_i.$$

We saw above that in the pair-structure (of the two-process mutual exclusion system) of Figure 4, T_1 and C_1 are the sometimes-blocking states. Hence, for the I -system (Figure 7) synthesized from the two-process mutual exclusion system (Figure 1), weak blocking fairness is: $\bigwedge_{i \in \text{dom}(I)} \tilde{G}((T_i \vee C_i) \wedge \text{en}_i) \Rightarrow \tilde{F} \text{ex}_i$. In other words, processes that remain enabled and in their trying or critical sections continuously must eventually be executed (note that requiring execution “eventually” or “infinitely often” makes no difference).

We denote the use of weak blocking fairness in our technical results by relativizing our notion of satisfaction of CTL* formulae (\models) so that only fullpaths that satisfy weak blocking fairness are taken into account. The resulting notion of satisfaction is denoted by \models_{Φ} , and is defined by Emerson and Lei [1985] (for the more traditional weak- and strong-fairness notions):

$$\begin{aligned} M, s \models_{\Phi} A f & \text{ iff } M, s \models A(\Phi \Rightarrow f) \\ M, s \models_{\Phi} E f & \text{ iff } M, s \models E(\Phi \wedge f) \end{aligned}$$

Effectively path quantification is only over the paths that satisfy Φ .

The following proposition is a consequence of the definition of sometimes-blocking. The essential idea is that if $M_{ij}, s_{ij} \models \neg r_i \wedge A F r_i$, then along every path leaving s_{ij} , P_i^j must eventually be executed, since the current i -state must eventually change for r_i (which depends only on the current i -state) to go from *false* to *true*. If s_i is not sometimes-blocking (i.e., always-nonblocking), then there will be at least one fullpath leaving s_{ij} that consists entirely of P_j^i transitions. This is because every P_j^i transition from s_{ij} to, say s'_{ij} , leaves the current i -state unchanged, i.e., still always-nonblocking. This implies that there is still (another) enabled P_j^i -transition from s'_{ij} . Repeating this argument, we conclude the existence of the aforementioned fullpath.

PROPOSITION 6.7.3.4 (SOMETIMES-BLOCKING). *Let s_{ij} be a reachable state of M_{ij} and $r_i \in \mathcal{L}(\mathcal{AP}_i, \neg, \wedge)$. If $M_{ij}, s_{ij} \models \neg r_i \wedge A F r_i$, then $s_{ij} \uparrow i$ is sometimes-blocking.*

The converse of this proposition — namely, “if $s_{ij} \uparrow i$ is sometimes-blocking, then an eventuality of the form $A F r_i$ is pending in s_{ij} ” — does not hold in general. For example, in the pair-structure for the mutual exclusion problem (given in Figure 4), $[C_i]$ is a sometimes-blocking i -state, but there is no eventuality of the form $A F r_i$ pending in state $[C_i T_j]$ of that structure.

6.7.4 The Progress Lemma. We can now demonstrate that, under weak blocking fairness as defined above, and given the liveness assumption, every I -process with a pending eventuality (clause (3) in the antecedent of Lemma 6.7.4.1) is eventually executed. Very roughly, the proof idea is as follows. The liveness assumption (clause (1)) guarantees that a process with a pending eventuality is not delayed forever by

another process via a wait-chain. Given that a process cannot be delayed forever, weak blocking fairness (\models_{Φ}) is sufficient to guarantee that the process is eventually executed, provided that a deadlock (clause (2)) does not occur.

LEMMA 6.7.4.1 (PROGRESS). *If*

- (1) *the liveness assumption LV holds,*
- (2) *for every reachable I-state s , $W_I(s)$ is supercycle-free, and*
- (3) *v is a reachable I-state of M_I such that $\bigwedge k \in I(\ell). (M_{k\ell}, v \upharpoonright_{k\ell} \models \neg r_\ell \wedge AFr_\ell)$ for some $\ell \in \text{dom}(I)$ and $r_\ell \in \mathcal{L}(\mathcal{AP}_\ell, \neg, \wedge)$, then*

$$M_I, v \models_{\Phi} AFe_{x_\ell}.$$

6.8 The Generalized Large-Model Theorem

We now establish soundness of MP-synthesis for the general case of arbitrary interconnection relations. It turns out that the argument differs from the large-model theorem only for temporal leads-to properties. For this case, our argument is roughly as follows. Assume that some temporal leads-to property is satisfied by the pair-systems. We use the progress lemma above to show that a process with a pending eventuality is always guaranteed to be eventually executed. Thus, until the eventuality is fulfilled, the process will be executed repeatedly. Nonfulfillment of the pending eventuality coupled with infinitely often execution, would, by path projection, imply the same phenomenon in the pair-systems. Hence, the pending eventuality is not fulfilled in the pair-systems, so the pair-systems do not satisfy the temporal leads-to property. But this contradicts our initial assumption.

THEOREM 6.8.1 (GENERALIZED LARGE MODEL). *Let $f_{k\ell}$ be an arbitrary formula of $FLCTL_{k\ell}$ (Definition 6.6.1). Let s be an arbitrary reachable I-state, and, for all i, j such that $i I j$, let $s_{ij} = s \upharpoonright_{ij}$. If the liveness assumption LV holds, and $W_I(u)$ is supercycle-free for every reachable I-state u , then*

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models \mathbf{\Lambda}_{k\ell} f_{k\ell}) \text{ implies } M_I, s \models_{\Phi} \mathbf{\Lambda}_{k\ell} f_{k\ell}.$$

As in Section 6.6, we also present the soundness result in terms of correctness properties, i.e., properties that hold in all the initial states of the pair-system or I-system.

COROLLARY 6.8.2 (GENERALIZED LARGE MODEL). *Let $f_{k\ell}$ be an arbitrary formula of $FLCTL_{k\ell}$ (Definition 6.6.1). Let (i, j) be an arbitrary pair in I . If the liveness assumption LV holds, and $W_I(u)$ is supercycle-free for every reachable I-state u , then*

$$M_{ij}, S_{ij}^0 \models \mathbf{\Lambda}_{k\ell} f_{k\ell} \text{ implies } M_I, S_I^0 \models_{\Phi} \mathbf{\Lambda}_{k\ell} f_{k\ell}.$$

PROOF. Identical to that of the large-model corollary (6.6.3), except that the generalized large-model theorem (6.8.1) is invoked instead of the large-model theorem (6.6.2). \square

We now look at our solution to the generalized dining philosophers problem given in Figure 7. We verify that this solution satisfies the liveness assumption.

Here M_{jk} is the pair-structure for the two-process mutual exclusion problem. We easily check, by inspection of Figure 4, that if $M_{jk}, s_J \uparrow jk \models EGex_k$, then $s_J \uparrow j = [N_j]$. From Figure 5 and the compact MP-synthesis definition (5.4.1), we infer, by straightforward checking, that, for $J = \{\{i, j\}, \{j, k\}\}, \bigwedge a_i^J \cdot (a_i^J \longrightarrow P_j^J \notin W_J(s_J))$ is always true when $s_J \uparrow j = [N_j]$, i.e., a process in its noncritical state does not block any move of any neighbor. Thus, we conclude that the dining philosophers solution satisfies the liveness assumption.

In Section 6.5, we showed that the solution for the generalized dining philosophers problem given in Figure 7 satisfies the wait-for-graph assumption. Now the wait-for-graph assumption depends only on the pair-system from which the K -process system is derived. It is independent of the interconnection relation I . Since the dining philosophers solution is derived from the same pair-system as the K -process mutual exclusion solution (namely, the two-process mutual exclusion system), the only difference being the particular interconnection relation used, we conclude that the dining philosophers solution satisfies the wait-for-graph assumption. We can likewise verify that $W_I(s_I^0)$ is supercycle-free for every initial state s_I^0 of the dining philosophers solution, by an argument analogous to that for the K -process mutual exclusion solution given in Section 6.5. Thus, by the supercycle-free wait-for-graph theorem (6.5.4.3), we conclude that $W_I(t)$ is supercycle-free for every reachable state t of the dining philosophers solution.

We have thus shown that the dining philosophers solution meets both assumptions of the generalized large-model corollary (6.8.2), so we conclude that the dining philosophers solution satisfies the MPCTL* specification given in Section 3.4. Since the interconnection relation I can specify any interconnection pattern, our solution solves the *generalized* dining philosophers problem put forth by Chandy and Misra [1984]. In particular, when interpreted for $I = \{\{1, 2\}, \{2, 3\}, \{3, 4\}, \{4, 5\}, \{5, 1\}\}$, our solution solves the original dining philosophers problem put forth by Dijkstra.

7. IMPLEMENTATION ON REALISTIC ARCHITECTURES

The programs that we synthesize have a large grain of atomicity in that a process must in one atomic step (1) inspect the local states and (pairwise) shared variables of all of its neighbors and (2) modify all of the shared variables. We now sketch a method of implementing our programs using low atomicity interprocess communication and synchronization primitives that are easily provided by most distributed systems, namely broadcast (as in most local area networks, which are ethernet based), and asynchronous point-to-point FIFO message passing (as in most wide-area networks).

First, we translate a synthesized program into a form which uses the *multiparty interaction* [Back and Kurki-Suonio 1984; 1988; 1989; Charlesworth 1987; Forman 1987; Francez and Forman 1996; Francez et al. 1986; Ramesh and Mehndiratta 1987] as the sole interprocess communication and synchronization primitive. A multiparty interaction is an atomic synchronous event involving a fixed set of processes (the *participants*). The participant processes must all reach a point where executing the interaction is one of the possible continuations. The interaction is then *enabled*. Only enabled interactions can be executed. During an interaction, data can be exchanged between any pair of processes [Forman 1987; Francez and Forman 1996; Ramesh and Mehndiratta 1987]. However, we do not make use of this feature; we

use interactions for synchronization only.

The translation of an I system $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ into a multiparty-interaction-based program proceeds as follows:

- For each P_i^I , $i \in \{i_1, \dots, i_K\}$, every move is replaced by a single arc labeled with a unique interaction name.
- Every pair-structure M_{ij} , $(i I j)$ is instantiated as a process, which we call a *pair-controller*. Each P_i^J -transition in M_{ij} is relabeled with the same interaction name as the corresponding move in P_i^I . The variables in \mathcal{SH}_{ij} now become local variables of the process M_{ij} .
- For every initial state $s_I^0 \in S_I^0$ we have an initial state t_I^0 of the translated program. In t_I^0 , the state of every P_i^I , $i \in \{i_1, \dots, i_K\}$, is $s_I^0[i]$, and the state of every M_{ij} , $i I j$, is $s_I^0[ij]$.

For example, the generalized dining philosophers solution is given in Figure 15, where i and j are replicated for all $(i, j) \in I$ (it is not necessary to display P_j^I as well as P_i^I ; this is done to help clarify the example). In P_i^I , the move from T_i to C_i has been replaced by a single multiparty interaction, $enter_i$, whose participants are P_i^I , and M_{ij} for all $j \in I(i)$. Since every M_{ij} such that $j \in I(i)$ must participate in the execution of $enter_i$, we see that P_i^I cannot enter the critical section without the “permission” of all the pair-controllers just as in the original solution. In fact, it is relatively straightforward to see that the original and translated programs have the same semantics, i.e., they generate isomorphic I -structures. The K -process mutual exclusion solution is a special case, which contains pair-controllers for every pair P_i^I, P_j^I of processes.

The second step of our implementation method uses one of the many distributed algorithms available [Back and Kurki-Suonio 1988; Back et al. 1985; Bagrodia 1989; Chandy and Misra 1987; Chandy and Misra 1988] for implementing multiparty interactions in terms of asynchronous message passing or ethernet broadcast. This then gives us our final implementation in terms of widely available, low atomicity primitives.

The number of pair-controllers (and therefore of processes introduced) is equal to the number of pairs in I , which, in the worst case, is $O(K^2)$. Thus, this scheme may, in some cases, produce somewhat inefficient programs, with large numbers of processes whose only purpose is to implement the pairwise constraints. This problem may be alleviated by merging some of the pair-controllers. For example, pair-controllers M_{ij} and M_{jk} may be merged into a process M_{ijk} as follows: M_{ijk} stores M_{ij} and M_{jk} as data structures, together with the “current states” of M_{ij} and M_{jk} . Using this information, M_{ijk} can determine which interactions it should participate in, and what the next states of M_{ij} and M_{jk} should be after the execution of an interaction. We emphasize that we do not generate M_{ijk} by taking the automata-theoretic product of M_{ij} and M_{jk} . Rather, M_{ijk} acts as an “interpreter” that simulates the execution of M_{ij} and M_{jk} (thus, we avoid generating in the worst case an exponentially large structure). Note that pair-controllers which are merged do not have to have common process indices, but the most efficient programs are generated when pair-controllers with common indices are merged.

Since we can merge any number of pair-controllers in this way, we can generate multiparty-interaction-based programs that range across the distributed/centralized

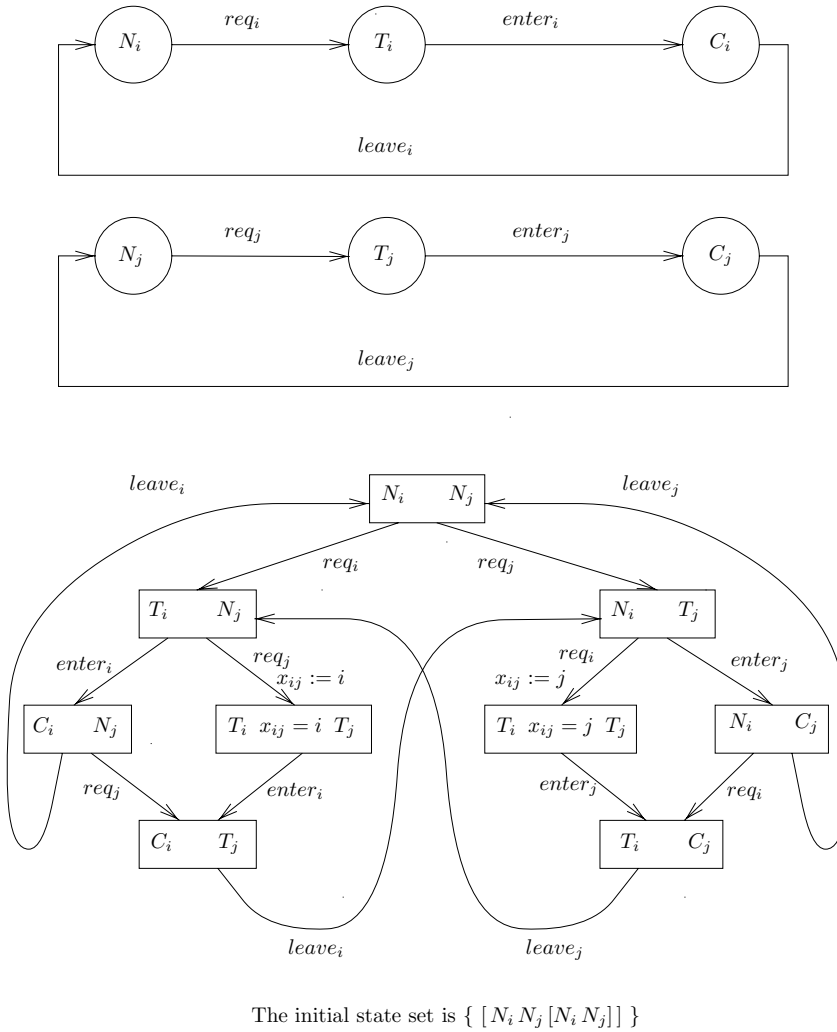


Fig. 15. Multiparty-interaction solution to the generalized dining philosophers problem.

spectrum. The more centralized programs will use fewer messages overall, but will have worse bottleneck performance, since a process that represents a large number of pair-controllers will have to handle a large number of messages, as it will interact with all of the P_i^I for every i that is an index of one of the pair-controllers represented. Likewise, more distributed programs will have better bottleneck performance, but will generate more messages overall. Thus, we have the usual tradeoffs and can make the appropriate choice based on the desired performance characteristics and on the target architecture. A detailed study of these issues is beyond the scope of this article.

In certain special cases, we can reduce the atomicity of the operations in our synthesized programs without appealing to a multiparty-interaction scheduling al-

gorithm. We outline this in the following two sections.

7.1 Reducing Read Atomicity: Guard Stability

The solutions produced by MP-synthesis contain read and write operations on vectors of shared variables. There are certain conditions under which these vector operations can be broken down into a sequence of simple operations. For example, in the MP-synthesis solution to the mutual exclusion problem, the guard on the T_i to C_i arc is $\bigwedge_{j \in I(i)} (N_j \vee (T_j \wedge x_{ij} = i))$. Each individual conjunct of this guard, i.e., $(N_j \vee (T_j \wedge x_{ij} = i))$ for some particular value of j , is *temporarily stable* [Katz 1986] in the sense that if P_i^I is “at” T_i , then once $(N_j \vee (T_j \wedge x_{ij} = i))$ holds, it will continue to hold until P_i^I enters the critical section C_i . Hence the conjuncts of the guard do not need to be checked simultaneously in one “large” vector operation. Any sequential order of checking the guard conjuncts will do.

We generalize this discussion as follows. Let $(s_i, \bigwedge_{j \in I(i)} B_i^j \rightarrow \bigvee_{j \in I(i)} A_i^j, t_i)$ be an arc of P_i^I . If $M_{ij}, S_{ij}^0 \models AG((\{s_i\} \wedge B_i^j) \Rightarrow A[(\{s_i\} \wedge B_i^j) \cup t_i])$, then the conjuncts B_i^j in $\bigwedge_{j \in I(i)} B_i^j$ can be checked sequentially instead of simultaneously, since once true, each B_i^j will remain true until the arc is executed. The condition $M_{ij}, S_{ij}^0 \models AG((\{s_i\} \wedge B_i^j) \Rightarrow A[(\{s_i\} \wedge B_i^j) \cup t_i])$ can be checked in polynomial time by the model-checking algorithm of Clarke et al. [1986].

7.2 Reducing the Atomicity of Assignment Statements

In the previous section we gave a way of reducing read atomicity. We now show how to reduce write atomicity in some cases. The solutions produced by MP-synthesis contain write operations on vectors of shared variables, e.g., $\bigvee_{j \in I(i)} A_i^j$ in P_i^I , where the A_i^j have to be $x_{ij} := i$ or $x_{ij} := j$. We now assume that P_i^I contains only assignments of the form $x_{ij} := j$; P_j^I contains only assignments of the form $x_{ij} := i$; and every move assigns to at most one shared variable. We now implement these assignments in terms of the solution to the generalized dining philosophers problem given in Chandy and Misra [1984] and in Chandy and Misra [1988], which uses asynchronous FIFO message passing. Let each P_i^I be a philosopher, and let P_i^I and P_j^I share a fork if and only if $i I j$. For each shared variable x_{ij} in the pair-system $(S_{ij}^0, P_i^I \parallel P_j^I)$, we introduce an instance of the dining philosophers algorithm. By our assumption, the only assignments to x_{ij} in P_i^I are of the form $\bigvee_{j \in I(i)} x_{ij} := j$. We implement x_{ij} in terms of the *priority relation* of the dining philosophers algorithm. This relation determines which philosopher has priority in obtaining the shared fork. We encode the value of x_{ij} as follows:

$$\begin{aligned} x_{ij} = i & \text{ iff } P_i^I \text{ has priority} \\ x_{ij} = j & \text{ iff } P_j^I \text{ has priority} \end{aligned}$$

Now when P_i^I needs to execute $\bigvee_{j \in I(i)} x_{ij} := j$, it transitions to “hungry” and waits to “eat.” After P_i^I has eaten, all forks will be at P_i^I and will be dirty, i.e., P_j^I (for all $j \in I(i)$) will have priority, as required. Thus the large atomicity multiple assignment $\bigvee_{j \in I(i)} x_{ij} := j$ has been implemented in terms of the small atomicity operations of the dining philosophers algorithm.

8. DISCUSSION

8.1 Scope of the Synthesis Method

Our synthesis method requires the pair-systems to satisfy two technical assumptions: the wait-for-graph assumption and the liveness-assumption. A consequence of this is that the method is not *complete*: it is possible that a given pair-system may not satisfy these assumptions, and yet the I -system synthesized from it may still satisfy the MPCTL* specification from which the pair-system was produced. It is natural to ask to what extent the range of applicability of our method is limited by imposing these assumptions.

Roughly, the liveness assumption prevents a process from executing continuously while delaying another process that has a pending eventuality. This behavior should not occur in a correct program anyway, so the liveness assumption is not restrictive at all; it can be thought of as a *design rule*: any system permitting the behavior prevented by the liveness assumption should be considered to be badly designed, since it allows individual starvation. The wait-for-graph assumption is somewhat restrictive. Roughly, it requires that a process join a wait-chain either at its beginning or at its end. However, there are situations where processes enter wait-chains “in the middle,” e.g. priority queues, and where the I -system would still be deadlock-free. Essentially, the wait-for-graph assumption represents a particular strategy for *deadlock prevention*. We expect that a more general deadlock prevention strategy (e.g., acquire resources in a fixed linear order) should also be applicable to our synthesis method. Investigating this and generalizing the wait-for-graph assumption are two of our research goals.

Even with our current wait-for-graph assumption, we point out that many systems can be synthesized. We have successfully applied the method to the well-known resource contention problems from the distributed computing literature: mutual exclusion, readers-writers, dining philosophers, drinking philosophers, and the m -out-of- n mutual exclusion problem (in which n processes contend for m identical and unsharable resources). Thus, we see that our method already deals with a large and important class of concurrent and distributed systems, namely resource contention. Generalizing the wait-for-graph assumption will enlarge its scope even further.

8.2 Complexity of the Synthesis Method

There are three algorithmic steps in applying the method: (1) applying Definition 5.1.1 to synthesize the I -system from the given pair-system, (2) checking the wait-for-graph assumption, and (3) checking the liveness-assumption. These have worst-case time complexities of $O(|I| \cdot |P_k^\ell|)$, $O(b \cdot |P_k^\ell|^{b+4} \cdot 2^{(b+1)}|\mathcal{SH}_{k\ell}|)$, and $O(|P_k^\ell|^3 \cdot 2^{2|\mathcal{SH}_{k\ell}|})$, respectively (see Appendix E), where b is the maximum branching factor (i.e., maximum number of outgoing moves from a single node) of P_k^ℓ . Since b is at least 1, the second term dominates the third, and hence the overall complexity of the synthesis method is the sum of the first two terms, i.e., $O(|I| \cdot |P_k^\ell| + b \cdot |P_k^\ell|^{b+4} \cdot 2^{(b+1)}|\mathcal{SH}_{k\ell}|)$.

The first term clearly cannot be improved, since the given pair-system and interconnection relation that are the inputs to the method have to be examined at least once. The second term is due to the check of the wait-for-graph assumption. The

check given in this article is basically a brute-force approach: the structure M_J for the relevant J -system(s) is first constructed, and then essentially model checked against the wait-for-graph assumption. It may be possible to perform this check without exhaustively generating M_J , e.g., by using the BDD-based model-checking techniques of Browne et al. [1986; 1990] and Clarke et al. [1992]. In particular, Browne et al. [1990] state that symmetric functions have compact BDD representations. Since the wait-for-graph assumption is symmetric in most of its arguments, namely the processes ℓ_1, \dots, ℓ_n , this approach appears promising.

Devising a more efficient check may allow us to eliminate the exponential factor in b . We note that the exponential factor in $|\mathcal{SH}_{k\ell}|$ cannot be eliminated in general, since the number of reachable states of $M_{k\ell}$ is exponential in $|\mathcal{SH}_{k\ell}|$ in the worst case.

8.3 Efficiency of the Synthesized System

The systems currently synthesized are not very efficient, since they have poor *locality*: a process can wait for another process that is quite distant from it (a process that communicates with it indirectly via a long chain of other processes). Furthermore, a process can wait for another process repeatedly, i.e., the second process can acquire and release resources many times before the first process progresses. By regarding the pair-controllers as resources (they can be regarded as such, since their permission is required in order to make a transition), we can apply efficient distributed resource allocation strategies to enhance the efficiency of the distributed low-atomicity systems produced by the implementation technique outlined above. For example, in Lynch [1981] and Styer and Peterson [1988], distributed resource allocation algorithms are presented where a process' waiting time is a "small function of the number and speed of nearby processes" and where a process cannot repeatedly acquire and release resources ahead of its neighbors (I -neighbors in our context).

9. RELATED WORK

The only other temporal logic synthesis method which exploits similarity of processes that we are aware of is that of Emerson and Srinivasan [1990] and Srinivasan [1991]. Specifications in this method are expressed in a logic similar to MPCTL. The method constructs a model of a given (satisfiable) specification that can be regarded as a global state transition diagram in which the (global) states record the number of processes in a particular local state. From this model, a concurrent program with an arbitrary number of processes is extracted. By recording only the number of processes in each local state, instead of the explicit local states, the size of the model is kept independent of the number of processes, and so state explosion is avoided.

One potential disadvantage of this method is that all the processes share a single variable *dist*, which holds the identity of a current "distinguished" process. In our framework, this means that the interconnection relation I is always the complete graph. Thus, communication between certain pairs of processes (e.g., non-neighboring philosophers) will be required even when it is not really necessary for a solution. Variable *dist* is used to schedule in a round-robin fashion all processes that have pending eventualities. Since this may include pairs of processes that do

not “conflict” with each other (again the example of non-neighboring philosophers is illustrative) the resulting ordering would, in many cases, be overly restrictive and would result in poor performance. Additionally, *dist* constitutes a bottleneck, with all of the well-known performance and reliability problems that result.

Other temporal logic synthesis methods to date [Anuchitanukul and Manna 1994; Emerson and Clarke 1982; Manna and Wolper 1984; Pnueli and Rosner 1989a; 1989b] all rely on some form of explicit global state enumeration (either generating an explicit global state transition diagram or a tree automaton) and so have at least exponential time complexity.

Related to synthesis is the technique of *model checking*: given a program and a specification, determine mechanically whether the program satisfies the specification. In Clarke et al. [1986] and German and Sistla [1992], methods of model checking finite-state concurrent programs (with respect to a specification expressed in some temporal logic) that exploit process similarity are presented. Clarke et al. [1986] require the manual construction of a bisimulation between the two-process and K -process programs, and their method is therefore not completely mechanical. Clarke and Grumberg [1987] attempt to automate the method of Clarke et al. [1986], but require an intermediate construction that is exponentially large in the size of a single process. The method is also incomplete in the same sense that ours is: there is no guarantee that the required technical assumptions will be satisfied in every case. German and Sistla [1992] present a model-checking method for programs consisting of strictly identical processes (no differentiating process indices) that communicate via CCS-style [Milner 1989] synchronous actions. Their model-checking algorithm runs in time polynomial in the size of a single process and exponential in the size of the specification. Unfortunately, the program notation is so weak that a program satisfying mutual exclusion cannot be written. To accommodate mutual exclusion, a distinguished controller process must be introduced. However, the running time now becomes doubly exponential in both process and specification size.

Finally, we mention Clarke et al. [1993] and Emerson and Sistla [1993], where the symmetry of the global state transition diagram that results from process similarity is exploited. The basic idea is to model check a reduced form of the global state transition diagram that results by taking the quotient with respect to a symmetry group. Computing the symmetry group is expensive in general, so either manual techniques [Emerson and Sistla 1993] or approximation methods [Clarke et al. 1993] must be used.

10. FURTHER WORK

We believe our work is a significant step in the direction of producing a practical temporal logic synthesis technique although there are still many open problems for future research. Some of these are as follows:

- (1) The accommodation of multiple classes of processes, with each class containing an arbitrarily large number of similar processes. Let the classes be $\mathcal{C}_1, \dots, \mathcal{C}_N$. Now we require the following pair-systems as input:
 - One pair-system per class \mathcal{C}_n , which is a model for a pair of generic processes $P_{n,i}, P_{n,j}$ which are representative of that class. This model captures the

interactions of the processes within \mathcal{C}_n .

- One pair-system for every pair of classes $\mathcal{C}_n, \mathcal{C}_m$, which is a model for a pair of generic representative processes $P_{n,i}, P_{m,j}$, where $P_{n,i}$ represents \mathcal{C}_n and $P_{m,j}$ represents \mathcal{C}_m . This model captures the interactions of processes in \mathcal{C}_n with processes in \mathcal{C}_m .

The interconnection relation would now relate processes in different classes as well as processes in the same class. The method of computing I -processes from pair-processes by Definition 5.1.1 remains as before. However, the spatial modalities of MPCTL are modified slightly so that the indices that they quantify over range over a particular class. So, for example, we would have $\bigwedge_{j \in \mathcal{C}_n}$ quantify over the set of processes in class \mathcal{C}_n , and we would have $\bigwedge_{i \in \mathcal{C}_n, j \in \mathcal{C}_m}$ to quantify over the pairs of processes where one element of the pair is in class \mathcal{C}_n and the other process is in class \mathcal{C}_m . The generalized large-model theorem should then allow us to conclude that

$$M_{ij}, S_{ij}^0 \models \bigwedge_{k \in \mathcal{C}_n, \ell \in \mathcal{C}_m} f_{k\ell} \text{ implies } M_I, S_I^0 \models \bigwedge_{k \in \mathcal{C}_n, \ell \in \mathcal{C}_m} f_{k\ell},$$

given that some antecedents (the analogue of the wait-for-graph assumption, for example) hold. Here i and j are arbitrary elements of \mathcal{C}_n and \mathcal{C}_m , respectively.

- (2) We intend to formalize the implementation method outlined in Section 7 and prove its correctness. We will also improve the efficiency of the final implemented system using the resource allocation algorithms of Lynch [1981] and Styer and Peterson [1988] (see Section 8.3). Further improvements may also be obtained by devising special-purpose multiparty-interaction implementation algorithms that exploit the symmetric structure of the programs produced by MP-synthesis and/or the specific program semantics. For example, in some cases, the tight synchronization of multiparty interactions may not be needed, and it may be possible to directly replace a multiparty interaction by a much more efficient asynchronous broadcast [Birman and Clark 1994].
- (3) Another issue that bears investigation is the use of a “disjunctive” composition technique to synthesize many-process systems. The technique we presented is “conjunctive” and is suited for problems such as the mutual exclusion problem, where access to the critical section requires the “permission” of all the processes in the system. Some problems have a “disjunctive” nature, for example the diffusing computation problem [Dijkstra et al. 1983], and it would be worthwhile to attempt to produce a synthesis technique which addresses such problems.
- (4) There has been considerable success in the application of model checking to the verification of VLSI circuits [Browne et al. 1986; Clarke et al. 1989; 1992; Mishra and Clarke 1985]. It would be worthwhile to investigate the problem of synthesizing circuits from temporal logic specifications (this would require the extension to multiple-process classes as a prerequisite).
- (5) We intend to investigate the synthesis of real-time concurrent systems. Our approach is to use a real-time variant of CTL, such as RTCTL [Emerson et al. 1993], to express real-time specifications. Already, we have established a real-time version of the large-model theorem [Attie 1995]. It remains to do the same for the generalized large-model theorem.
- (6) Finally we shall investigate the synthesis of fault-tolerant concurrent systems.

11. CONCLUSIONS

In this article, we have shown how to extend the synthesis methods of Emerson and Clark [1982] and Manna and Wolper [1984] to deal with an arbitrary number of similar processes without incurring the exponential overhead due to the state explosion problem. Since we have reduced the problem of reasoning about K processes to the problem of reasoning about a small number of processes, the results of Emerson et al. [1992] imply that the synthesis problem can be solved in polynomial time. Our method works for arbitrary process interconnection schemes. We also outlined a way of implementing the synthesized programs in terms of interprocess communication and synchronization primitives that are available in practice. Thus, we see this contribution as a significant step toward producing a mechanical synthesis technique capable of generating solutions for realistic problems.

APPENDIX

A. GLOSSARY OF SYMBOLS

\models	Satisfies relation of CTL, CTL*
\models_{Φ}	Satisfies relation of CTL, CTL* relativized to fairness notion Φ
Φ	CTL* path formula that specifies weak blocking fairness
\otimes	“Conjunctive” guarded-command composition operator
\oplus	“Disjunctive” guarded-command composition operator
$\{\}$	State to formula operator
\bigwedge_i	Spatial modality (quantifies over individual processes)
\bigwedge_{ij}	Spatial modality (quantifies over pairs of processes related by I)
$\uparrow i$	State projection onto a single process i
$\uparrow ij$	State or path projection onto an ij -system
$\uparrow J$	State or path projection onto a J -system
$j_1/i_1, \dots, j_m/i_m$	Process index substitution operator
θ	Synonym for $j_1/i_1, \dots, j_m/i_m$
a_i^j	Pair-move (move in a compact pair-process)
a_i^I	I -move (move in a compact I -process)
$a_i^I.start$	Start state of the I -move a_i^I
$a_i^I.guard$	Guard of the I -move a_i^I
$a_i^I.guard_j$	Conjunct of $a_i^I.guard$ that tests shared state with process j
blk_i	Propositional formula true in exactly the sometimes-blocking local states
$dom(I)$	Synonym for i_1, \dots, i_K
i_1, \dots, i_K	Domain of I (i.e., the set of process indices in the synthesized system)
$t_k.moves$	The set of moves with start state t_k
$FLCTL_{kl}$	The sublogic Flat CTL
I	The interconnection relation
M_{ij}	The ij -structure
M_I	The I -structure
M_I^r	The reachable portion of M_I
P_i^j	Process i in the pair-system consisting of processes i and j
P_i^I	Process i in the system synthesized using interconnection

	relation I
R_{ij}	The transition relation of M_{ij}
R_I	The transition relation of M_I
R_I^r	The reachable portion of R_I
S_{ij}	The set of states of M_{ij}
S_{ij}^0	The set of initial states of M_{ij}
S_I	The set of states of M_I
S_I^0	The set of initial states of M_I
S_I^r	The set of reachable states of M_I
$W_I(s)$	The wait-for-graph for the I -system in global state s
$(S_{ij}^0, P_i^j \parallel P_j^i)$	The pair-system consisting of processes i and j
$(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$	The I -system
\mathcal{AP}	The family of sets of atomic propositions of the synthesized system
\mathcal{AP}_i	The set of atomic propositions of process i
\mathcal{L}	The sublogics constructor
\mathcal{SH}_{ij}	The set of variables shared by processes i and j

B. PROOF OF THE LARGE-MODEL THEOREM

THEOREM 6.6.2 (LARGE MODEL). *Let f_{kl} be an arbitrary formula of $FLCTL_{kl}$. Let $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$. Furthermore, let s be an arbitrary reachable I -state, and, for all i, j such that $i I j$, let $s_{ij} = s \upharpoonright ij$. If $M_I, S_I^0 \models AGEXtrue$, then*

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models \mathbf{\Lambda}_{kl} f_{kl}) \text{ implies } M_I, s \models \mathbf{\Lambda}_{kl} f_{kl}.$$

PROOF. By MPCTL* semantics (see Section 3), $M_{ij}, s_{ij} \models \mathbf{\Lambda}_{kl} f_{kl}$ is equivalent to $M_{ij}, s_{ij} \models f_{ij}$. Also by MPCTL* semantics, $M_I, s \models \mathbf{\Lambda}_{kl} f_{kl}$ is equivalent to $M_I, s \models \bigwedge (i, j) \in I. (f_{ij})$. This is equivalent, by CTL* semantics, to $\bigwedge (i, j) \in I. (M_I, s \models f_{ij})$. Hence the large-model theorem is established if we can prove that

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models f_{ij}) \text{ implies } \bigwedge (i, j) \in I. (M_I, s \models f_{ij}) \quad (\text{P})$$

given the assumptions of the large-model theorem. The proof is by induction on the structure of f_{ij} . In each case we assume the antecedent $\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models f_{ij})$, and we establish the consequent $\bigwedge (i, j) \in I. (M_I, s \models f_{ij})$. In the remainder of the proof, we make the convention that all free occurrences of i, j represent arbitrary elements of $dom(I)$ such that $i I j$. Thus it suffices to establish $M_I, s \models f_{ij}$. Note that in some cases (which do not invoke the inductive hypothesis), we will, for convenience, use $M_{ij}, s_{ij} \models f_{ij}$ as the antecedent, with i, j free. Hence we will, for these cases, actually be establishing that $\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models f_{ij} \text{ implies } M_I, s \models f_{ij})$, which implies (P). This is permissible, since we only do this for those cases which do not invoke the inductive hypothesis.

$f_{ij} = h_{ij}$. The antecedent is $M_{ij}, s_{ij} \models h_{ij}$. Now $s_{ij} = s \upharpoonright ij$ by assumption. Thus by the I -state projection proposition (6.3.1) with $J = \{\{i, j\}\}$, we have $M_I, s \models h_{ij}$.

$f_{ij} = f'_{ij} \wedge f''_{ij}$. The antecedent is $\bigwedge(i, j) \in I. (M_{ij}, s_{ij} \models (f'_{ij} \wedge f''_{ij}))$. So, by CTL* semantics, we get $\bigwedge(i, j) \in I. (M_{ij}, s_{ij} \models f'_{ij})$ and $\bigwedge(i, j) \in I. (M_{ij}, s_{ij} \models f''_{ij})$. From these and the induction hypothesis, we get $\bigwedge(i, j) \in I. (M_I, s \models f'_{ij})$ and $\bigwedge(i, j) \in I. (M_I, s \models f''_{ij})$. So by CTL* semantics we get $\bigwedge(i, j) \in I. (M_I, s \models (f'_{ij} \wedge f''_{ij}))$.

$f_{ij} = AGh_{ij}$. Let t be an arbitrary s -reachable state in M_I . By assumption $s \upharpoonright ij = s_{ij}$. Thus, by the relativized state-mapping corollary (6.4.6) with $J = \{\{i, j\}\}$, $t \upharpoonright ij$ is a s_{ij} -reachable state in M_{ij} . The antecedent is $M_{ij}, s_{ij} \models AGh_{ij}$. Hence $M_{ij}, t \upharpoonright ij \models h_{ij}$ by CTL* semantics. From this and the I -state projection proposition (6.3.1) with $J = \{\{i, j\}\}$, we get $M_I, t \models h_{ij}$. Since t is an arbitrary s -reachable state in M_I , we conclude $M_I, s \models AGh_{ij}$ by CTL* semantics.

$f_{ij} = AG(p_i \Rightarrow AY_i q_i)$. We will establish that

$$M_I, t \models (p_i \Rightarrow AY_i q_i) \quad (*)$$

where t is an arbitrary s -reachable state in M_I (remember that t is s -reachable if and only if there is a path from s to t). If $M_I, t \models \neg p_i$ then $(*)$ holds immediately, and we are done. Otherwise $M_I, t \models p_i$, and we proceed as follows. The antecedent is

$$M_{ij}, s_{ij} \models AG(p_i \Rightarrow AY_i q_i). \quad (a)$$

By assumption, $s \upharpoonright ij = s_{ij}$. Thus by the relativized state-mapping corollary (6.4.6) with $J = \{\{i, j\}\}$, $t \upharpoonright ij$ is a s_{ij} -reachable state in M_{ij} . So, by (a) we have

$$M_{ij}, t \upharpoonright ij \models (p_i \Rightarrow AY_i q_i). \quad (b)$$

Since $M_I, t \models p_i$, we have $M_{ij}, t \upharpoonright ij \models p_i$ by the I -state projection proposition (6.3.1) with $J = \{\{i, j\}\}$. Together with (b), this yields

$$M_{ij}, t \upharpoonright ij \models AY_i q_i. \quad (c)$$

Now let $t \xrightarrow{i} u$ be an arbitrary P_i^I -transition (in M_I) out of state t (if no such transition exists, then $(*)$ holds by CTL* semantics, and we are done). By the transition-mapping lemma (6.4.1), we have

$$t \upharpoonright ij \xrightarrow{i} u \upharpoonright ij \in R_{ij}. \quad (d)$$

By (c), (d), and CTL* semantics we have

$$u \upharpoonright ij \models q_i. \quad (e)$$

By (e) and the I -state projection proposition (6.3.1) with $J = \{\{i, j\}\}$ we have

$$u \models q_i.$$

Since $t \xrightarrow{i} u$ is an arbitrary P_i^I -transition out of state t , we have $M_I, t \models AY_i q_i$, and therefore $M_I, t \models (p_i \Rightarrow AY_i q_i)$.

We have therefore established $(*)$. Since t is an arbitrary s -reachable state in M_I , we conclude that $M_I, s \models AG(p_i \Rightarrow AY_i q_i)$.

$f_{ij} = AG(p_i \Rightarrow A[q_i Ur_i])$. We will establish $M_I, t \models (p_i \Rightarrow A[q_i Ur_i])$ where t is an arbitrary s -reachable state in M_I . If $M_I, t \models \neg p_i$ then $M_I, t \models (p_i \Rightarrow A[q_i Ur_i])$, and we are done. Otherwise we have $M_I, t \models p_i$, and we proceed as follows.

Let π be an arbitrary fullpath starting in t . By assumption, $M_I, S_I^0 \models AGEXtrue$, and so π is infinite. Since the number $K (= |dom(I)|)$ of I -processes is finite, π contains an infinite number of transitions of at least one I -process. Call this I -process P_k^I . We now select an I -process P_ℓ^I as follows. If $k = i$, then ℓ is an arbitrary member of $I(i)$. If $k \neq i$, then ℓ is k (note that k, ℓ depend on the choice of π , in general). Hence, we have

$$\begin{aligned} &\pi \text{ contains an infinite number of } P_i^I \text{ transitions or} \\ &\text{an infinite number of } P_\ell^I \text{ transitions.} \end{aligned} \quad (**)$$

The antecedent is

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models AG(p_i \Rightarrow A[q_i Ur_i])).$$

By assumption, $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$, so we have $(i, \ell) \in I$. Thus

$$M_{i\ell}, s_{i\ell} \models AG(p_i \Rightarrow A[q_i Ur_i]). \quad (a)$$

By assumption, $s \upharpoonright \ell = s_{i\ell}$. Thus by the relativized state-mapping corollary (6.4.6) with $J = \{\{i, \ell\}\}$, $t \upharpoonright \ell$ is a $s_{i\ell}$ -reachable state in $M_{i\ell}$. So by (a) we have

$$M_{i\ell}, t \upharpoonright \ell \models (p_i \Rightarrow A[q_i Ur_i]). \quad (b)$$

Since $M_I, t \models p_i$, we have $M_{i\ell}, t \upharpoonright \ell \models p_i$ by the I -state projection proposition (6.3.1) with $J = \{\{i, \ell\}\}$. Together with (b), this yields

$$M_{i\ell}, t \upharpoonright \ell \models A[q_i Ur_i]. \quad (c)$$

By the path-mapping lemma (6.4.3), $\pi \upharpoonright \ell$ is a path in $M_{i\ell}$. By (**) and the path projection definition (5.1.2.1), $\pi \upharpoonright \ell$ contains an infinite number of P_ℓ^I -transitions or an infinite number of P_i^I -transitions, and so $\pi \upharpoonright \ell$ is a fullpath. Also, by the path projection definition (5.1.2.1), $\pi \upharpoonright \ell$ starts in $t \upharpoonright \ell$, and therefore, by (c), we have $M_{i\ell}, \pi \upharpoonright \ell \models [q_i Ur_i]$. By the path projection lemma (6.3.3) with $J = \{\{i, \ell\}\}$, we have $M_I, \pi \models [q_i Ur_i]$.

Since π is an arbitrary fullpath starting in t , we conclude, by CTL* semantics, that $M_I, t \models A[q_i Ur_i]$, and hence $M_I, t \models (p_i \Rightarrow A[q_i Ur_i])$. Since t is an arbitrary s -reachable state in M_I , we conclude that $M_I, s \models AG(p_i \Rightarrow A[q_i Ur_i])$.

$f_{ij} = AG(a_i \Rightarrow EX_i b_i)$. We will establish $M_I, t \models (a_i \Rightarrow EX_i b_i)$ where t is an arbitrary s -reachable state in M_I . If $M_I, t \models \neg a_i$ then $M_I, t \models (a_i \Rightarrow EX_i b_i)$, and we are done. Otherwise $M_I, t \models a_i$, and we proceed as follows.

Since $a_i, b_i \in \{\{s_i\} \mid s_i \in S_i\}$, let $a_i = \{s_i^m\}, b_i = \{s_i^n\}$, respectively, for some s_i^m, s_i^n in S_i . By the antecedent, we have

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models AG(a_i \Rightarrow EX_i b_i)).$$

Hence, letting i be free (and remembering our convention for i, j) we have

$$\bigwedge j \in I(i). (M_{ij}, s_{ij} \models AG(a_i \Rightarrow EX_i b_i)).$$

By assumption, $s\uparrow ij = s_{ij}$. Thus, by the relativized state-mapping corollary (6.4.6) with $J = \{\{i, j\}\}$, $t\uparrow ij$ is a s_{ij} -reachable state in M_{ij} for all j in $I(i)$. So,

$$\bigwedge j \in I(i) . (M_{ij}, t\uparrow ij \models (a_i \Rightarrow EX_i b_i)).$$

Since $M_I, t \models a_i$, we conclude $M_{ij}, t\uparrow ij \models a_i$ by the I -state projection proposition (6.3.1). Together with the above this yields

$$\bigwedge j \in I(i) . (M_{ij}, t\uparrow ij \models EX_i b_i). \quad (a)$$

From this and CTL* semantics, we get

$$\bigwedge j \in I(i) . (t\uparrow ij \xrightarrow{i} u_{ij} \in R_{ij} \text{ for some } ij\text{-state } u_{ij} \text{ such that } u_{ij} \models b_i). \quad (b)$$

From $b_i = \{s_i^n\}$ and $u_{ij} \models b_i$, we have

$$\bigwedge j \in I(i) . (u_{ij} \models \{s_i^n\}). \quad (d)$$

By (d) and the local state projection proposition (6.3.2) we have

$$\bigwedge j \in I(i) . (u_{ij}\uparrow i \models \{s_i^n\}).$$

By this and the definition of $\{s_i^n\}$ (Section 3) we get

$$\bigwedge j \in I(i) . (u_{ij}\uparrow i = s_i^n). \quad (e)$$

Now let u be a global state such that

$$\begin{aligned} & u\uparrow i = s_i^n \\ & \bigwedge j \in I(i) . (u\uparrow j = u_{ij}\uparrow j \text{ and } u\uparrow \mathcal{SH}_{ij} = u_{ij}\uparrow \mathcal{SH}_{ij}) \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i) . (u\uparrow j = t\uparrow j) \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j \neq k . (u\uparrow \mathcal{SH}_{jk} = t\uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (f)$$

By (e) and (f), we have

$$\bigwedge j \in I(i) . (u\uparrow i = u_{ij}\uparrow i \text{ and } u\uparrow j = u_{ij}\uparrow j \text{ and } u\uparrow \mathcal{SH}_{ij} = u_{ij}\uparrow \mathcal{SH}_{ij}).$$

From this we have

$$\bigwedge j \in I(i) . ((u\uparrow i \cup u\uparrow j \cup u\uparrow \mathcal{SH}_{ij}) = (u_{ij}\uparrow i \cup u_{ij}\uparrow j \cup u_{ij}\uparrow \mathcal{SH}_{ij})).$$

Since $(u\uparrow i \cup u\uparrow j \cup u\uparrow \mathcal{SH}_{ij}) = u\uparrow ij$, and $(u_{ij}\uparrow i \cup u_{ij}\uparrow j \cup u_{ij}\uparrow \mathcal{SH}_{ij}) = u_{ij}$, we have

$$\bigwedge j \in I(i) . (u\uparrow ij = u_{ij}). \quad (g)$$

By (b) and (g) we have

$$\bigwedge j \in I(i) . (t\uparrow ij \xrightarrow{i} u\uparrow ij \in R_{ij}). \quad (h)$$

By (h) and (f) we have

$$\begin{aligned} & \bigwedge j \in I(i) . (t\uparrow ij \xrightarrow{i} u\uparrow ij \in R_{ij}) \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i) . (u\uparrow j = t\uparrow j) \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j \neq k . (u\uparrow \mathcal{SH}_{jk} = t\uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (i)$$

By (i) and the transition-mapping lemma (6.4.1), we have

$$t \xrightarrow{i} u \in R_I.$$

Now $u \uparrow ij = u_{ij}$, and $u_{ij} \models b_i$ (by (b)). So by the I -state projection proposition (6.3.1), $u \models b_i$. Thus by CTL* semantics we get

$$M_I, t \models EX_i b_i,$$

so $M_I, t \models (a_i \Rightarrow EX_i b_i)$. Since t is an arbitrary s -reachable state in M_I , we conclude $M_I, s \models AG(a_i \Rightarrow EX_i b_i)$.

This last case concludes the proof of (P), which in turn completes the proof of the large-model theorem. \square

COROLLARY 6.6.3 (LARGE MODEL). *Let f_{kl} be an arbitrary formula of $FLCTL_{kl}$. Let $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$, and let (i, j) be an arbitrary pair in I . If $M_I, S_I^0 \models AGEXtrue$, then*

$$M_{ij}, S_{ij}^0 \models \bigwedge_{kl} f_{kl} \text{ implies } M_I, S_I^0 \models \bigwedge_{kl} f_{kl}.$$

PROOF. We assume the antecedent $M_{ij}, S_{ij}^0 \models \bigwedge_{kl} f_{kl}$, and we establish the consequent $M_I, S_I^0 \models \bigwedge_{kl} f_{kl}$. From $M_{ij}, S_{ij}^0 \models \bigwedge_{kl} f_{kl}$ and Observation 6.2.2, we have

$$\bigwedge (i, j) \in I. (M_{ij}, S_{ij}^0 \models \bigwedge_{kl} f_{kl}). \quad (\text{a})$$

Let s_I^0 be an arbitrary initial state (i.e., member of S_I^0). By the MP-synthesis definition (5.1.1), $\bigwedge (i, j) \in I. (s_I^0 \uparrow ij \in S_{ij}^0)$. Thus by (a) we have

$$\bigwedge (i, j) \in I. (M_{ij}, s_I^0 \uparrow ij \models \bigwedge_{kl} f_{kl}).$$

Hence, by the large-model theorem (6.6.2), with $s = s_I^0$, we obtain

$$M_I, s_I^0 \models \bigwedge_{kl} f_{kl}.$$

Since s_I^0 is an arbitrarily chosen member of S_I^0 , we have

$$M_I, S_I^0 \models \bigwedge_{kl} f_{kl}. \quad \square$$

ONLINE-ONLY APPENDICES

C. PROOFS

D. A COMPACT REPRESENTATION FOR SYNCHRONIZATION SKELETONS

E. CHECKING THE WAIT-FOR-GRAPH ASSUMPTION AND THE LIVENESS ASSUMPTION

Appendices C-E are available only online. You should be able to get the online-only appendices from the citation page for this article:

<http://www.acm.org/pubs/citations/journals/toplas/1998-20-1/p51-attie/>

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ACKNOWLEDGMENTS

The authors would like to thank Prasad Sistla for very helpful discussions on viewing process indices in terms of spatial modalities. We also thank Mani Chandy, and Mike Evangelist for helpful comments. We are also indebted to the referees, whose comments helped to significantly improve the presentation.

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Received March 1996; revised August 1996; accepted July 1997

THIS DOCUMENT IS THE ONLINE-ONLY APPENDIX TO:

Synthesis of Concurrent Systems with Many Similar Processes

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ACM Transactions on Programming Languages and Systems, Vol. 20, No. 1, January 1998, Pages 51–115.

C. PROOFS

C.1 Process Similarity

The *process index substitution operator* $\theta = \{j_1/i_1, \dots, j_m/i_m\}$ denotes the simultaneous replacement of process indices i_1, \dots, i_m by process indices j_1, \dots, j_m respectively. We require that i_1, \dots, i_m be pairwise distinct, and j_1, \dots, j_m be pairwise distinct. θ can be applied to all of the pair syntactic constructs defined in the article, as well as any pair model (e.g., M_{ij}). We define θ in a bottom-up manner as follows.

Definition C.1.1 (Process Index Substitution Operator). The process index substitution operator $\theta = \{j_1/i_1 \dots j_m/i_m\}$ is defined as follows:

- (1) θ distributes through propositional logic connectives, $=$, $:=$, \rightarrow , $//$, and \cup .
- (2) For any process index i :

$$\begin{aligned} i\theta &= j_k \text{ if } i = i_k \text{ for some } k \in [1 : m] \\ i\theta &= i \text{ otherwise.} \end{aligned}$$

- (3) For any $Q_i \in \mathcal{AP}_i$:

$$Q_i\theta = Q_{i\theta}$$

- (4) For any $x_{ij} \in \mathcal{SH}_{ij}$:

$$\begin{aligned} x_{ij}\theta &= x_{i\theta, j\theta} \\ x_{ij}^i\theta &= x_{i\theta, j\theta}^{i\theta} \\ x_{ij}^j\theta &= x_{i\theta, j\theta}^{j\theta} \end{aligned}$$

- (5) For any i -state s_i :

$$s_i\theta = \{ \langle Q_{i\theta}, s_i(Q_i) \rangle \mid Q_i \in \mathcal{AP}_i \}$$

i.e., when $i = i_k$, $s_i\theta$ is a j_k -state. It satisfies the atomic propositions in \mathcal{AP}_{j_k} , which correspond to the atomic propositions in \mathcal{AP}_i that s_i satisfies (remember

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that the sets of atomic proposition form a uniform family, see Section 3 on MPCTL*). Note that $s_i\theta = s_i$ if $i \neq i_k$ for all $k \in [1 : m]$.

(6) For any pair-process P_i^j :

$$P_i^j\theta = \{(s_i\theta, B_i^j\theta \rightarrow A_i^j\theta, t_i\theta) \mid (s_i, B_i^j \rightarrow A_i^j, t_i) \in P_i^j\}$$

(7) For any ij -state s_{ij} :

$$\begin{aligned} s_{ij}\theta = & \{ \langle Q_{i\theta}, s_{ij}(Q_i) \rangle \mid Q_i \in \mathcal{AP}_i \} \cup \\ & \{ \langle Q_{j\theta}, s_{ij}(Q_j) \rangle \mid Q_j \in \mathcal{AP}_j \} \cup \\ & \{ \langle x_{ij}\theta, h_{ij}\theta \rangle \mid s_{ij}(x_{ij}) = h_{ij}, x_{ij} \in \mathcal{SH}_{ij} \} \end{aligned}$$

(8) For transition relation R_{ij} :

$$R_{ij}\theta = \{(s_{ij}\theta, h_{ij}\theta, t_{ij}\theta) \mid (s_{ij}, h_{ij}, t_{ij}) \in R_{ij}\}$$

(9) For the sets of initial ij -states S_{ij}^0 and all ij -states S_{ij} :

$$\begin{aligned} S_{ij}^0\theta &= \{s_{ij}\theta \mid s_{ij} \in S_{ij}^0\} \\ S_{ij}\theta &= \{s_{ij}\theta \mid s_{ij} \in S_{ij}\} \end{aligned}$$

(10) For the pair-structure $M_{ij} = (S_{ij}^0, S_{ij}, R_{ij})$:

$$M_{ij}\theta = (S_{ij}^0\theta, S_{ij}\theta, R_{ij}\theta)$$

PROPOSITION 6.2.1 (PAIR-STRUCTURE SIMILARITY). *Let i, j, i', j' be arbitrary elements of $\{i_1, \dots, i_K\}$ such that $i I j$ and $i' I j'$. Then we have*

$$M_{ij} = M_{i'j'}\{i/i', j/j'\}.$$

PROOF. From the pair-structure definition (5.2.1), we have $M_{i'j'}\{i/i', j/j'\} = (S_{i'j'}^0, S_{i'j'}, R_{i'j'})\{i/i', j/j'\}$. By the process index substitution operator definition (C.1.1), this is equal to $(S_{i'j'}^0\{i/i', j/j'\}, S_{i'j'}\{i/i', j/j'\}, R_{i'j'}\{i/i', j/j'\})$. Now $S_{i'j'}^0\{i/i', j/j'\} = S_{ij}^0$ by the initial-state assumption. $S_{i'j'}$ is the set of all possible $i'j'$ -states. Now $\mathcal{AP}_{i'}, \mathcal{AP}_{j'}$ are similar to $\mathcal{AP}_i, \mathcal{AP}_j$ by assumption (see Section 3.4), and since $P_i^j = P_{i'}^{j'}\{i/i', j/j'\}$, we must have $\mathcal{SH}_{ij} = \mathcal{SH}_{i'j'}\{i/i', j/j'\}$, since \mathcal{SH}_{ij} is merely the set of shared variables the occur in P_i^j, P_j^i . Thus, $S_{ij} = S_{i'j'}\{i/i', j/j'\}$, since their domains are related by $\{i/i', j/j'\}$. Finally, by the process similarity assumption and the pair structure definition (5.2.1), we can infer $R_{ij} = R_{i'j'}\{i/i', j/j'\}$, since similar arcs in $P_i^j, P_{i'}^{j'}$ give rise to similar transitions in $R_{ij}, R_{i'j'}$, respectively. \square

C.2 State and Path Projection Results

PROPOSITION 6.3.1 (I-STATE PROJECTION). *Let $J \subseteq I$, and let s be an I -state. Then*

$$s \models f \text{ iff } s \upharpoonright J \models f$$

where f is a formula of $\mathcal{L}(\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i, \neg, \wedge)$.

PROOF. The proof is by induction on the structure of f .

$f \in \bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$. $s \models f$ iff $s(f) = \text{true}$ iff (since $s \upharpoonright J = (\bigcup_{i \in \text{dom}(J)} s \upharpoonright i) \cup (\bigcup_{(i,j) \in J} s \upharpoonright \mathcal{SH}_{ij})$) $s \upharpoonright J(f) = \text{true}$ iff $s \upharpoonright J \models f$.

$f = g \wedge h$. $s \models (g \wedge h)$ iff $(s \models g \text{ and } s \models h)$ iff (by the inductive hypothesis) $(s \upharpoonright J \models g \text{ and } s \upharpoonright J \models h)$ iff $s \upharpoonright J \models (g \wedge h)$.

$f = \neg g$. $s \models \neg g$ iff not $(s \models g)$ iff (by the inductive hypothesis) not $(s \upharpoonright J \models g)$ iff $s \upharpoonright J \models \neg g$. \square

PROPOSITION 6.3.2 (LOCAL STATE PROJECTION). *Let $i \in \text{dom}(J)$, and let s_J be a J -state. Then*

$$s_J \models f_i \text{ iff } s_J \upharpoonright i \models f_i$$

where f_i is a formula of $\mathcal{L}(\mathcal{AP}_i, \neg, \wedge)$.

PROOF. Analogous to that of Proposition 6.3.1. \square

LEMMA 6.3.3 (PATH PROJECTION). *Let π be a path in M_I , and let $J \subseteq I$. Then*

$$\pi \models f \text{ iff } \pi \upharpoonright J \models f$$

where f is a formula of $\mathcal{L}(\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i, \neg, \wedge, U)$.

PROOF. The proof is by induction on the structure of f .

$f \in \bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$. Let s, s_J be the initial states of $\pi, \pi \upharpoonright J$ respectively. By the definition of path projection, $s_J = s \upharpoonright J$, and so s_J and s agree on all atomic propositions in $\bigcup_{i \in \text{dom}(J)} \mathcal{AP}_i$. Hence $s \models f$ iff $s_J \models f$. Also $\pi \models f$ iff $s \models f$ and $\pi \upharpoonright J \models f$ iff $s_J \models f$ by CTL* semantics (Section 3.1). These three equivalences yield $\pi \models f$ iff $\pi \upharpoonright J \models f$.

$f = g \wedge h$. $\pi \models (g \wedge h)$ iff $(\pi \models g \text{ and } \pi \models h)$ iff, by the inductive hypothesis, $(\pi \upharpoonright J \models g \text{ and } \pi \upharpoonright J \models h)$ iff $\pi \upharpoonright J \models (g \wedge h)$.

$f = \neg g$. $\pi \models \neg g$ iff not $(\pi \models g)$ iff, by the inductive hypothesis, not $(\pi \upharpoonright J \models g)$ iff $\pi \upharpoonright J \models \neg g$.

$f = gUh$. Proof by double implication.

Left to right: $\pi \models gUh$ implies $\pi \upharpoonright J \models gUh$.

Assume $\pi \models gUh$. Therefore, for some $n_0 \geq 1$, we have by CTL* semantics (Section 3.1):

$$\pi^{n_0} \models h \text{ and } \bigwedge n. (1 \leq n < n_0 \Rightarrow \pi^n \models g). \quad (\text{a})$$

By (a) and the induction hypothesis, we get

$$\pi^{n_0} \upharpoonright J \models h \text{ and } \bigwedge n. (1 \leq n < n_0 \Rightarrow \pi^n \upharpoonright J \models g). \quad (\text{b})$$

Let B^{m_0} denote the J -block of π that contains s^{n_0} (the first state of π^{n_0}). By the path projection definition (5.1.2.1), $\pi^{n_0} \upharpoonright J = (\pi \upharpoonright J)^{m_0}$, since the m_0 'th J -block of π corresponds to the m_0 'th state of $\pi \upharpoonright J$. So, by (b) we have

$$(\pi \upharpoonright J)^{m_0} \models h. \quad (\text{c})$$

Let m be an arbitrary integer in $[1 : (m_0 - 1)]$. The first state s_J^m of $(\pi \upharpoonright J)^m$ corresponds to the m th J -block of π (by the path projection definition (5.1.2.1)). Let s^l (the first state of π^l) be an arbitrary state of the m th J -block of π . Hence

$(\pi \uparrow J)^m = \pi^l \uparrow J$. Also, since $m < m_0$, s^l occurs in an earlier J -block of π than s^{n_0} , and therefore $l < n_0$. So by (b) we have

$$\pi^l \uparrow J \models g. \quad (d)$$

By (d) and $(\pi \uparrow J)^m = \pi^l \uparrow J$, we get

$$(\pi \uparrow J)^m \models g. \quad (e)$$

But m is an arbitrary integer in $[1 : (m_0 - 1)]$. So, by (c) and (e) we get

$$(\pi \uparrow J)^{m_0} \models h \text{ and } \bigwedge m. (1 \leq m < m_0 \Rightarrow (\pi \uparrow J)^m \models g).$$

By CTL* semantics (Section 3.1), this is equivalent to $\pi \uparrow J \models gUh$.

Right to left: $\pi \uparrow J \models gUh$ implies $\pi \models gUh$.

Assume $\pi \uparrow J \models gUh$. Therefore, for some $m_0 \geq 1$, we have by CTL* semantics (Section 3.1)

$$(\pi \uparrow J)^{m_0} \models h \text{ and } \bigwedge m. (1 \leq m < m_0 \Rightarrow (\pi \uparrow J)^m \models g). \quad (a)$$

Let s^{n_0} (the first state of π^{n_0}) be the first state of the m_0 th J -block of π . By the path projection definition (5.1.2.1), $\pi^{n_0} \uparrow J = (\pi \uparrow J)^{m_0}$. Hence, by (a) we have

$$\pi^{n_0} \uparrow J \models h. \quad (b)$$

Now let m be an arbitrary integer in $[1 : (m_0 - 1)]$. By (a), we get

$$(\pi \uparrow J)^m \models g. \quad (c)$$

Furthermore, let s^n (the first state of π^n) be an arbitrary state of the m th J -block of π . By the path projection definition (5.1.2.1), $\pi^n \uparrow J = (\pi \uparrow J)^m$. Also, since $m < m_0$, s^n occurs in an earlier J -block of π than s^{n_0} , and therefore $n < n_0$. Furthermore, since s^{n_0} is the *first* state of the m_0 th J -block of π , we see that n ranges over $[1 : (n_0 - 1)]$, since s^{n_0-1} occurs in an earlier J -block of π than s^{n_0} . Hence, by (c) and $\pi^n \uparrow J = (\pi \uparrow J)^m$ we have

$$\bigwedge n. (1 \leq n < n_0 \Rightarrow \pi^n \uparrow J \models g). \quad (d)$$

From (b) and (d) we get

$$\pi^{n_0} \uparrow J \models h \text{ and } \bigwedge n. (1 \leq n < n_0 \Rightarrow \pi^n \uparrow J \models g). \quad (e)$$

From (e) and the induction hypothesis applied to g, h , we get

$$\pi^{n_0} \models h \text{ and } \bigwedge n. (1 \leq n < n_0 \Rightarrow \pi^n \models g).$$

By CTL* semantics (Section 3.1), this is equivalent to $\pi \models gUh$. \square

C.3 Mapping of I -Structures into J -Structures

LEMMA 6.4.1 (TRANSITION MAPPING). *For all I -states $s, t \in S_I$ and $i \in \text{dom}(I)$,*

$$\begin{aligned} s \xrightarrow{i} t \in R_I \text{ iff :} \\ \bigwedge j \in I(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij}) \text{ and} \\ \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned}$$

PROOF. Let s, t be arbitrary I -states, and let i be an arbitrary element of $\text{dom}(I)$. By the I -structure definition (5.3.1), $s \xrightarrow{i} t \in R_I$ is equivalent to

$$\begin{aligned} & (s \uparrow i, \bigwedge_{j \in I(i)} B_i^j \rightarrow //_{j \in I(i)} A_i^j, t \uparrow i) \text{ is an arc in } P_i^I \text{ and} \\ & \bigwedge j \in I(i). (s \uparrow i j (B_i^j) = \text{true} \text{ and } < s \uparrow \mathcal{SH}_{ij} > A_i^j < t \uparrow \mathcal{SH}_{ij} >) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \{i\}. (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (a)$$

By the MP-synthesis definition (5.1.1), (a) is equivalent to

$$\begin{aligned} & \bigwedge j \in I(i). ((s \uparrow i, B_i^j \rightarrow A_i^j, t \uparrow i) \text{ is an arc in } P_i^j) \text{ and} \\ & \bigwedge j \in I(i). (s \uparrow i j (B_i^j) = \text{true} \text{ and } < s \uparrow \mathcal{SH}_{ij} > A_i^j < t \uparrow \mathcal{SH}_{ij} >) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \{i\}. (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (b)$$

Now (b), by rewriting the “ $\bigwedge j \in \text{dom}(I) - \{i\} \dots$ ” universal quantification as the conjunction of “for all j in $I(i) \dots$ ” and “ $\bigwedge j \in \text{dom}(I) - \hat{I}(i) \dots$ ”, is equivalent to

$$\begin{aligned} & \bigwedge j \in I(i). ((s \uparrow i, B_i^j \rightarrow A_i^j, t \uparrow i) \text{ is an arc in } P_i^j) \text{ and} \\ & \bigwedge j \in I(i). (s \uparrow i j (B_i^j) = \text{true} \text{ and } < s \uparrow \mathcal{SH}_{ij} > A_i^j < t \uparrow \mathcal{SH}_{ij} >) \text{ and} \\ & \bigwedge j \in I(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (c)$$

By merging all three “ $\bigwedge j \in I(i) \dots$ ” universal quantifications, (c) is equivalent to

$$\begin{aligned} & \bigwedge j \in I(i). (\\ & \quad (s \uparrow i, B_i^j \rightarrow A_i^j, t \uparrow i) \text{ is an arc in } P_i^j \text{ and} \\ & \quad s \uparrow i j (B_i^j) = \text{true} \text{ and} \\ & \quad < s \uparrow \mathcal{SH}_{ij} > A_i^j < t \uparrow \mathcal{SH}_{ij} > \text{ and} \\ & \quad s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (d)$$

By the obvious identities $s \uparrow i = (s \uparrow i j) \uparrow i$, $s \uparrow \mathcal{SH}_{ij} = (s \uparrow i j) \uparrow \mathcal{SH}_{ij}$, $s \uparrow j = (s \uparrow i j) \uparrow j$, $t \uparrow i = (t \uparrow i j) \uparrow i$, $t \uparrow \mathcal{SH}_{ij} = (t \uparrow i j) \uparrow \mathcal{SH}_{ij}$, $t \uparrow j = (t \uparrow i j) \uparrow j$, (d) is equivalent to

$$\begin{aligned} & \bigwedge j \in I(i). (\\ & \quad ((s \uparrow i j) \uparrow i, B_i^j \rightarrow A_i^j, (t \uparrow i j) \uparrow i) \text{ is an arc in } P_i^j \text{ and} \\ & \quad s \uparrow i j (B_i^j) = \text{true} \text{ and} \\ & \quad < (s \uparrow i j) \uparrow \mathcal{SH}_{ij} > A_i^j < (t \uparrow i j) \uparrow \mathcal{SH}_{ij} > \text{ and} \\ & \quad (s \uparrow i j) \uparrow j = (t \uparrow i j) \uparrow j) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned} \quad (e)$$

By the pair-structure definition (5.2.1), (e) is equivalent to

$$\begin{aligned} & \bigwedge j \in I(i). (s \uparrow i j \xrightarrow{i} t \uparrow i j \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}). \end{aligned}$$

The lemma then follows by transitivity of equivalence. \square

COROLLARY 6.4.2 (TRANSITION MAPPING). *Let $J \subseteq I$ and $i \in \text{dom}(J)$. If $s \xrightarrow{i} t \in R_I$, then $s \uparrow J \xrightarrow{i} t \uparrow J \in R_J$.*

PROOF. Assume $s \xrightarrow{i} t \in R_I$. Then, by the transition-mapping lemma (6.4.1), we have

$$\begin{aligned} & \bigwedge j \in I(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \hat{I}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{H}_{jk} = t \uparrow \mathcal{H}_{jk}). \end{aligned} \quad (a)$$

From $\bigwedge j \in I(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij})$ and the pair-structure definition (5.2.1), we get $\bigwedge j \in I(i). (s \uparrow j = t \uparrow j)$. Together with (a), this yields

$$\begin{aligned} & \bigwedge j \in I(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \{i\}. (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(I) - \{i\}, j I k. (s \uparrow \mathcal{H}_{jk} = t \uparrow \mathcal{H}_{jk}). \end{aligned} \quad (b)$$

Since $J \subseteq I$, we have $\text{dom}(J) \subseteq \text{dom}(I)$. Now $i \in \text{dom}(J)$ by assumption, and so $i \in \text{dom}(I)$. Thus, by $J \subseteq I$, we have $J(i) \subseteq I(i)$ and $\bigwedge j, k. (j J k \text{ implies } j I k)$. Thus, from (b) we get

$$\begin{aligned} & \bigwedge j \in J(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(I) - \{i\}. (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(J) - \{i\}, j J k. (s \uparrow \mathcal{H}_{jk} = t \uparrow \mathcal{H}_{jk}). \end{aligned} \quad (c)$$

Since $J \subseteq I$, we have $\text{dom}(J) \subseteq \text{dom}(I)$. Also, $i \in \hat{J}(i)$, and hence $\text{dom}(J) - \hat{J}(i) \subseteq \text{dom}(I) - \{i\}$. Thus, by (c) we have

$$\begin{aligned} & \bigwedge j \in J(i). (s \uparrow ij \xrightarrow{i} t \uparrow ij \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(J) - \hat{J}(i). (s \uparrow j = t \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(J) - \{i\}, j J k. (s \uparrow \mathcal{H}_{jk} = t \uparrow \mathcal{H}_{jk}). \end{aligned} \quad (d)$$

Now $s \uparrow ij = (s \uparrow J) \uparrow ij$ when $j \in I(i)$, and $s \uparrow j = (s \uparrow J) \uparrow j$ when $j \in \text{dom}(J)$, and $s \uparrow \mathcal{H}_{jk} = (s \uparrow J) \uparrow \mathcal{H}_{jk}$ when $j J k$. Thus, (d) can be rewritten as

$$\begin{aligned} & \bigwedge j \in J(i). ((s \uparrow J) \uparrow ij \xrightarrow{i} (t \uparrow J) \uparrow ij \in R_{ij}) \text{ and} \\ & \bigwedge j \in \text{dom}(J) - \hat{J}(i). ((s \uparrow J) \uparrow j = (t \uparrow J) \uparrow j) \text{ and} \\ & \bigwedge j, k \in \text{dom}(J) - \{i\}, j J k. ((s \uparrow J) \uparrow \mathcal{H}_{jk} = (t \uparrow J) \uparrow \mathcal{H}_{jk}). \end{aligned} \quad (e)$$

By the transition-mapping lemma (6.4.1) with $I := J$, and (e), we get $s \uparrow J \xrightarrow{i} t \uparrow J \in R_J$ as required. \square

LEMMA 6.4.3 (PATH MAPPING). *Let $J \subseteq I$. If π is a path in M_I , then $\pi \uparrow J$ is a path in M_J .*

PROOF. Let n be an arbitrary integer greater than zero, and let $u_J^n \xrightarrow{d_n} u_J^{n+1}$ be the n th transition along $\pi \uparrow J$. Thus $d_n \in \text{dom}(J)$, and u_J^n, u_J^{n+1} are the n th, $(n+1)$ st states respectively along $\pi \uparrow J$. By the path projection definition (5.1.2.1), $u_J^n = B^n \uparrow J$, $u_J^{n+1} = B^{n+1} \uparrow J$, where B^n, B^{n+1} are the n th, $(n+1)$ st, J -blocks respectively along π . Let s, t , be the last, first states of B^n, B^{n+1} , respectively, so $s \uparrow J = B^n \uparrow J = u_J^n$, $t \uparrow J = B^{n+1} \uparrow J = u_J^{n+1}$. Also, by the path projection definition (5.1.2.1), $s \xrightarrow{d_n} t$

is a transition along π , and therefore $s \xrightarrow{d_n} t \in R_I$, since π is a path in M_I . Hence, by the transition-mapping corollary (6.4.2), $s \uparrow J \xrightarrow{d_n} t \uparrow J \in R_J$, so $u_J^n \xrightarrow{d_n} u_J^{n+1} \in R_J$.

Since every transition of $\pi \uparrow J$ is a transition in R_J , it follows that $\pi \uparrow J$ is a path in M_J . \square

COROLLARY 6.4.4 (PATH MAPPING). *Let $J \subseteq I$. If π is an initialized path in M_I then $\pi \uparrow J$ is an initialized path in M_J .*

PROOF. Let B^0 be the first J -block of π . By the path projection definition (5.1.2.1), the first state of $\pi \uparrow J$ is $B^0 \uparrow J$. Since π is initialized, B^0 contains some initial state $s_I^0 \in S_I^0$. Hence $B^0 \uparrow J = s_I^0 \uparrow J$.

Now by the MP-synthesis definition (5.1.1), we have $\bigwedge(i, j) \in I. (s_I^0 \uparrow ij \in S_{ij}^0)$. Since $J \subseteq I$, we have $\bigwedge(i, j) \in J. (s_I^0 \uparrow ij \in S_{ij}^0)$. Also, by definition of $\uparrow J$ (Section 6), $\bigwedge(i, j) \in J. ((s_I^0 \uparrow J) \uparrow ij = s_I^0 \uparrow ij)$. Hence, we have $\bigwedge(i, j) \in J. ((s_I^0 \uparrow J) \uparrow ij \in S_{ij}^0)$. Now, by the MP-synthesis definition (5.1.1) with J replacing I we have $S_J^0 = \{s_J \mid \bigwedge(i, j) \in J. (s_J \uparrow ij \in S_{ij}^0)\}$. Hence we conclude $s_I^0 \uparrow J \in S_J^0$. Thus, the first state of $\pi \uparrow J$ is an initial state of M_J . By the path-mapping lemma (6.4.3), $\pi \uparrow J$ is a path in M_J . It follows that $\pi \uparrow J$ is an initialized path in M_J . \square

COROLLARY 6.4.5 (STATE MAPPING). *Let $J \subseteq I$. If t is a reachable state in M_I , then $t \uparrow J$ is a reachable state in M_J .*

PROOF. Since t is reachable in M_I , there must exist at least one initialized path π in M_I which ends in state t . By the path-mapping corollary (6.4.4), $\pi \uparrow J$ is an initialized path in M_J . By the path projection definition (5.1.2.1), $\pi \uparrow J$ ends in state $t \uparrow J$. Hence $t \uparrow J$ is reachable in M_J . \square

COROLLARY 6.4.6 (RELATIVIZED STATE MAPPING). *Let $J \subseteq I$. If t is a s -reachable state in M_I , then $t \uparrow J$ is a $s \uparrow J$ -reachable state in M_J .*

PROOF. Since t is s -reachable in M_I , there must exist at least one path π in M_I which starts in state s and ends in state t . By the path-mapping lemma (6.4.3), $\pi \uparrow J$ is a path in M_J . By the path projection definition (5.1.2.1), $\pi \uparrow J$ starts in state $s \uparrow J$ and ends in state $t \uparrow J$. Hence $t \uparrow J$ is $s \uparrow J$ -reachable in M_J . \square

C.4 Deadlock Freedom of the Many-Process System

PROPOSITION 6.5.4.1 (WAIT-FOR-GRAPH PROJECTION). *Let $J \subseteq I$ and $i J j$. Furthermore, let s_I be an arbitrary I -state. Then*

- (1) $P_i^I \longrightarrow a_i^I \in W_I(s_I)$ iff $P_i^J \longrightarrow a_i^J \in W_J(s_I \uparrow J)$, and
- (2) $a_i^I \longrightarrow P_j^I \in W_I(s_I)$ iff $a_i^J \longrightarrow P_j^J \in W_J(s_I \uparrow J)$.

PROOF. By assumption, $i J j$ and $J \subseteq I$. Hence $i I j$.

Proof of clause (1). By the wait-for-graph definition (6.5.2.1), $P_i^I \longrightarrow a_i^I \in W_I(s_I)$ iff $s_I \uparrow i = a_i^I.start$. Since $i \in dom(J)$, we have $(s_I \uparrow J) \uparrow i = s_I \uparrow i$ by definition of $\uparrow J$. Thus $s_I \uparrow i = a_i^I.start$ iff $(s_I \uparrow J) \uparrow i = a_i^I.start$ (since $a_i^I.start = a_i^J.start = s_i$). Finally, by the wait-for-graph definition (6.5.2.1) and $i J j$, $(s_I \uparrow J) \uparrow i = a_i^I.start$ iff

$P_i^J \longrightarrow a_i^J \in W_J(s_I \uparrow J)$. These three equivalences together yield clause (1) (using transitivity of equivalence).

Proof of clause (2). By the wait-for-graph definition (6.5.2.1), $a_i^I \longrightarrow P_j^I \in W_I(s_I)$ iff $s \uparrow ij \not\models a_i^I.\text{guard}_j$. Since $i J j$, we have $(s_I \uparrow J) \uparrow ij = s_I \uparrow ij$ by definition of $\uparrow J$. Also, $a_i^I.\text{guard}_j = a_i^J.\text{guard}_j = \bigvee_{\ell \in [1:n]} B_{i,\ell}^j$. Thus $s_I \uparrow ij \not\models a_i^I.\text{guard}_j$ iff $(s_I \uparrow J) \uparrow ij \not\models a_i^J.\text{guard}_j$. Finally, by the wait-for-graph definition (6.5.2.1) and $i J j$, $(s_I \uparrow J) \uparrow ij \not\models a_i^J.\text{guard}_j$ iff $a_i^J \longrightarrow P_j^J \in W_J(s_I \uparrow J)$. These three equivalences together yield clause (2), (using transitivity of equivalence, and noting that $s \not\models B$ and $s(B) = \text{false}$ have identical meaning). \square

THEOREM 6.5.4.3 (SUPERCYCLE-FREE WAIT-FOR-GRAPH). *If the wait-for-graph assumption WG holds, and $W_I(s_I^0)$ is supercycle-free for every initial state $s_I^0 \in S_I^0$, then for every reachable state t of M_I , $W_I(t)$ is supercycle-free.*

PROOF. Let t be an arbitrary reachable state of M_I , and let s be an arbitrary reachable state of M_I such that $s \xrightarrow{k} t$ for some $k \in \text{dom}(I)$. We shall establish that

$$\text{if } W_I(t) \text{ is supercyclic, then } W_I(s) \text{ is supercyclic.} \quad (\text{P1})$$

The contrapositive of P1 together with the assumption that $W_I(s_I^0)$ is supercycle-free for all $s_I^0 \in S_I^0$ is sufficient to establish the conclusion of the theorem (by induction on the length of a path from some $s_I^0 \in S_I^0$ to t).

We say that an edge is *k-incident* iff at least one of its vertices is P_k^I or a_k^I . The following (P2) will be useful in proving P1

$$\text{if edge } e \text{ is not } k\text{-incident, then } e \in W_I(t) \text{ iff } e \in W_I(s). \quad (\text{P2})$$

Proof of P2. If e is not k -incident, then, by the wait-for-graph definition (6.5.2.1), either $e = P_h^I \longrightarrow a_h^I$, or $e = a_h^I \longrightarrow P_\ell^I$, for some h, ℓ such that $h \neq k, \ell \neq k$. From $h \neq k, \ell \neq k$ and $s \xrightarrow{k} t \in R_I$, we have $s \uparrow h = t \uparrow h$ and $s \uparrow h\ell = t \uparrow h\ell$ by the wait-for-graph definition (6.5.2.1). Since $e \in W_I(t), e \in W_I(s)$ are determined solely by $t \uparrow h\ell, s \uparrow h\ell$ respectively, (see the wait-for-graph definition (6.5.2.1)), P2 follows. (End proof of P2.)

Let v be a vertex in a supercycle SC . We define $\text{depth}_{SC}(v)$ to be the length of the longest backward path in SC which starts in v . If there exists an infinite backward path (i.e., one that traverses a cycle) in SC starting in v , then $\text{depth}_{SC}(v) = \omega$ (for “infinity”). We now establish that

$$\text{every supercycle } SC \text{ contains at least one cycle.} \quad (\text{P3})$$

Proof of P3. Suppose P3 does not hold, and SC is a supercycle containing no cycles. Therefore, all backward paths in SC are finite, and so by definition of depth_{SC} all vertices of SC have finite depth. Thus, there is at least one vertex v in SC with maximal depth. But, by definition of depth_{SC} , v has no successors in SC , which, by the supercycle definition (6.5.3.1), contradicts the assumption that SC is a supercycle. (End proof of P3.)

Our final prerequisite for the proof of P1 is

if SC is a supercycle in $W_I(s)$, then the graph SC' obtained from SC by removing all vertices of finite depth from SC (along with incident edges) is also a supercycle in $W_I(s)$. (P4)

Proof of P4. By P3, $SC' \neq \emptyset$. Thus SC' satisfies clause (1) of the supercycle definition (6.5.3.1). Let v be an arbitrary vertex of SC' . Thus $v \in SC$ and $depth_{SC}(v) = \omega$ by definition of SC' . Let w be an arbitrary successor of v in SC . $depth_{SC}(w) = \omega$ by definition of $depth$. Hence $w \in SC'$. Furthermore, w is a successor of v in SC' , by definition of SC' . Thus every vertex v of SC' is also a vertex of SC , and the successors of v in SC' are the same as the successors of v in SC . Now since SC is a supercycle, every vertex v in SC has enough successors in SC to satisfy clauses (2) and (3) of the supercycle definition (6.5.3.1). It follows that every vertex v in SC' has enough successors in SC' to satisfy clauses (2) and (3) of the supercycle definition (6.5.3.1). (End proof of P4.)

We now present the proof of (P1). We assume the antecedent of P1 and establish the consequent. Let SC be some supercycle in $W_I(t)$. Let SC' be the graph obtained from SC by removing all vertices of finite depth from SC (along with incident edges). We now show that $P_k^I \notin SC'$ and that SC' contains no move vertex of the form a_k^I . There are two cases.

Case 1: $P_k^I \notin SC$. Then obviously $P_k^I \notin SC'$. Now suppose some node of the form a_k^I is in SC' . By definition of SC' , we have $a_k^I \in SC$ and $depth_{SC}(a_k^I) = \omega$. Hence, by definition of $depth$, there exists an infinite backward path in SC starting in a_k^I . Thus a_k^I must have a predecessor in SC . By the supercycle definition (6.5.3.1), P_k^I is the only possible predecessor of a_k^I in SC , and hence $P_k^I \in SC$, contrary to the case assumption. We therefore conclude that SC' contains no vertices of the form a_k^I . (End of case 1.)

Case 2: $P_k^I \in SC$. By the supercycle definition (6.5.3.1),

$$\bigwedge a_k^I \in W_I(t) . (\bigvee \ell . (a_k^I \longrightarrow P_\ell^I \in W_I(t))). \quad (a)$$

Since there are exactly n moves a_k^I of process P_k^I in $W_I(t)$ ($n = |t_k.moves|$), we can select ℓ_1, \dots, ℓ_n (where ℓ_1, \dots, ℓ_n are not necessarily pairwise distinct) such that

$$\bigwedge a_k^I \in W_I(t) . (\bigvee \ell \in \{\ell_1, \dots, \ell_n\} . (a_k^I \longrightarrow P_\ell^I \in W_I(t))). \quad (b)$$

Now let $J = \{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_n\}\}$ where j is an arbitrary element of $I(k)$. Applying the wait-for-graph projection proposition (6.5.4.1) to (b) gives us

$$\bigwedge a_k^J \in W_J(t \uparrow J) . (\bigvee \ell \in \{\ell_1, \dots, \ell_n\} . (a_k^J \longrightarrow P_\ell^J \in W_J(t \uparrow J))). \quad (c)$$

Now $s \xrightarrow{k} t \in R_I$ by assumption. Hence $s \uparrow J \xrightarrow{k} t \uparrow J \in R_J$ by the transition-mapping corollary (6.4.2). Also, by the state-mapping corollary (6.4.5) $s \uparrow J$ is reachable in M_J , since s is reachable in M_I . Thus we can apply the wait-for-graph assumption to $t \uparrow J$ to get

$$\bigwedge a_j^J . (a_j^J \longrightarrow P_k^J \notin W_J(t \uparrow J))$$

or

$$\bigvee a_k^J \in W_J(t \uparrow J) . (\bigwedge \ell \in \{\ell_1, \dots, \ell_n\} . (a_k^J \longrightarrow P_\ell^J \notin W_J(t \uparrow J))). \quad (d)$$

Now (c) contradicts the second disjunct of (d). Hence

$$\bigwedge a_j^J . (a_j^J \longrightarrow P_k^J \notin W_J(t \uparrow J)),$$

and applying the wait-for-graph projection proposition (6.5.4.1) to this gives us

$$\bigwedge a_j^J . (a_j^J \longrightarrow P_k^I \notin W_I(t)).$$

Since j is an arbitrary element of $I(k)$, we conclude that P_k^I has no incoming edges in $W_I(t)$. Thus, by definition of $depth$, $depth_{SC}(P_k^I) = 0$, and so $P_k^I \notin SC'$.

Now suppose some node of the form a_k^I is in SC' . By definition of SC' , we have $a_k^I \in SC$ and $depth_{SC}(a_k^I) = \omega$. Hence, by definition of $depth$, there exists an infinite backward path in SC starting in a_k^I . Thus a_k^I must have a predecessor in SC . By the supercycle definition (6.5.3.1), P_k^I is the only possible predecessor of a_k^I in SC , and hence there exists an infinite backward path in SC starting in P_k^I . Thus $depth_{SC}(P_k^I) = \omega$ by definition of $depth$. But we have established $depth_{SC}(P_k^I) = 0$, so we conclude that SC' contains no vertices of the form a_k^I . (End of case 2.)

In both cases, $P_k^I \notin SC'$, and SC' contains no move vertex of the form a_k^I . Thus every edge of SC' is not k -incident. Hence, by P2, every edge of SC' is an edge of $W_I(s)$ (since $SC' \subseteq W_I(t)$). By P4, SC' is a supercycle, so $W_I(s)$ is supercyclic. Thus P1 is established, which establishes the theorem. \square

C.5 Liveness Properties

PROPOSITION 6.7.3.4 (SOMETIMES-BLOCKING). *Let s_{ij} be a reachable state of M_{ij} and $r_i \in \mathcal{L}(\mathcal{AP}_i, \neg, \wedge)$. If $M_{ij}, s_{ij} \models \neg r_i \wedge AFr_i$, then $s_{ij} \uparrow i$ is sometimes-blocking.*

PROOF. We assume the antecedent and establish the consequent. Define a P_j^i -path to be a path in M_{ij} composed entirely of P_j^i -transitions. For a pair move $a_j^i = (s_i, \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$, let $a_j^i.start, a_j^i.guard$ denote $s_i, \bigvee_{\ell \in [1:n]} B_{i,\ell}^j$ respectively.

By assumption, $s_{ij} \not\models r_i$. Also, by $M_{ij}, s_{ij} \models AFr_i$ and CTL* semantics, every path starting in s_{ij} must lead to a state t_{ij} which fulfills AFr_i , i.e., $t_{ij} \models r_i$. Since $s_{ij} \uparrow i \neq t_{ij} \uparrow i$ (otherwise we could not have $s_{ij} \not\models r_i$ and $t_{ij} \models r_i$), we conclude, by the pair-structure definition (5.2.1), that every maximal path starting in s_{ij} must eventually contain a P_j^i -transition, because a P_j^i -transition cannot change any atomic proposition in \mathcal{AP}_i . Thus, every maximal P_j^i -path starting in s_{ij} (if any) is finite and ends in a state which has no outgoing P_j^i -transitions.

We now demonstrate the existence of a reachable state u_{ij} in M_{ij} such that $u_{ij} \uparrow i = s_{ij} \uparrow i$ and such that u_{ij} has no outgoing P_j^i -transitions. There are two cases.

Case 1: There are no P_j^i -paths starting in s_{ij} . Therefore s_{ij} has no outgoing P_j^i -transitions, so let u_{ij} be s_{ij} . (End of case 1.)

Case 2: There is at least one P_j^i -path starting in s_{ij} . Hence there is at least one maximal P_j^i -path starting in s_{ij} . By the above discussion, this path is finite and ends in a state with no outgoing P_j^i -transitions. Let u_{ij} be this state. Since there is a P_j^i -path starting in s_{ij} and ending in u_{ij} we conclude that $u_{ij} \uparrow i = s_{ij} \uparrow i$, since, by the pair-structure definition (5.2.1) a P_j^i -transition cannot change the truth value assigned to any atomic proposition in \mathcal{AP}_i . Also, u_{ij} is reachable, since s_{ij} is reachable and there is a path from s_{ij} to u_{ij} . (End of case 2.)

Since u_{ij} has no outgoing P_j^i -transitions, all of the moves in P_j^i (we are assuming compact notation here) are disabled in state u_{ij} , so we have

$$u_{ij} \models \bigwedge a_j^i \in P_j^i . (\neg \{a_j^i.start\} \vee \neg a_j^i.guard), \quad (a)$$

since a move is disabled iff control is not at its start state or its guard evaluates to false. Since every local state of a process has at least one outgoing arc (Section 2), there exists at least one move a_j^i in P_j^i such that $u_{ij} \models \{a_j^i.start\}$. From this and (a), we have

$$u_{ij} \models \bigvee a_j^i \in P_j^i . (\{a_j^i.start\} \wedge \neg a_j^i.guard). \quad (b)$$

Finally, since u_{ij} is reachable and $u_{ij} \uparrow i = s_{ij} \uparrow i$, we obtain from (b)

$$\bigvee s_{ij}^0 \in S_{ij}^0 . (M_{ij}, s_{ij}^0 \models EF(\{s_{ij} \uparrow i\} \wedge (\bigvee a_j^i \in P_j^i . (\{a_j^i.start\} \wedge \neg a_j^i.guard))))).$$

By the sometimes-blocking state definition (6.7.3.1), we conclude that $s_{ij} \uparrow i$ is sometimes-blocking. \square

LEMMA 6.7.4.1 (PROGRESS). *If*

- (1) *the liveness assumption LV holds,*
- (2) *for every reachable I-state s , $W_I(s)$ is supercycle-free, and*
- (3) *v is a reachable I-state of M_I such that $\bigwedge k \in I(\ell) . (M_{k\ell}, v \uparrow k \ell \models \neg r_\ell \wedge AFr_\ell)$ for some $\ell \in \text{dom}(I)$ and $r_\ell \in \mathcal{L}(\mathcal{AP}_\ell, \neg, \wedge)$, then*

$$M_I, v \models_\Phi AFex_\ell.$$

PROOF. We prove $\pi \models Fex_\ell$ for π an arbitrary fullpath in M_I such that v is the first state of π and $\pi \models \Phi$. By the definition of \models_Φ , this establishes $M_I, v \models_\Phi AFex_\ell$. Now $W_I(s)$ is supercycle-free for every reachable I -state s , by assumption. Hence, by the deadlock freedom theorem (6.5.5.2), we have $M_I, S_I^0 \models AGEXtrue$. Therefore, π is infinite. Also, by propositional and temporal logic reasoning (and using $nblk_i \equiv \neg blk_i$)

$$\pi \models \bigwedge i \in \text{dom}(I) . (\overset{\infty}{G}nblk_i \vee (\overset{\infty}{F}nblk_i \wedge \overset{\infty}{F}blk_i) \vee \overset{\infty}{G}blk_i).$$

Hence we can partition $\text{dom}(I)$ into three sets $\psi_{nblk}, \psi_{ex}, \psi_{blk}$ such that

$$\begin{aligned} \pi &\models \bigwedge i \in \psi_{nblk} . \overset{\infty}{G}nblk_i \\ \pi &\models \bigwedge i \in \psi_{ex} . (\overset{\infty}{F}nblk_i \wedge \overset{\infty}{F}blk_i) \\ \pi &\models \bigwedge i \in \psi_{blk} . \overset{\infty}{G}blk_i. \end{aligned}$$

So, by CTL* semantics, π has a suffix ρ such that

$$\rho \models \bigwedge i \in \psi_{nblk} . Gnblk_i \quad (a)$$

$$\rho \models \bigwedge i \in \psi_{ex} . (\overset{\infty}{F}nblk_i \wedge \overset{\infty}{F}blk_i) \quad (b)$$

$$\rho \models \bigwedge i \in \psi_{blk} . Gblk_i. \quad (c)$$

We now have three cases, depending on which of $\psi_{nblk}, \psi_{ex}, \psi_{blk}$ contains ℓ .

Case 1: $\ell \in \psi_{nblk}$. Since $\bigwedge k \in I(\ell) . (M_{k\ell}, v \uparrow k \ell \models \neg r_\ell \wedge AFr_\ell)$ by assumption, $v \uparrow \ell$ is sometimes-blocking by the sometimes-blocking proposition (6.7.3.4). Thus

$v \models blk_\ell$. Since $\rho \models nblk_\ell$, P_ℓ must have changed state along π , because v is the first state of π , and ρ is a suffix of π (remember that blk_ℓ is a purely propositional formula). By the I -structure definition (5.3.1), P_ℓ must have been executed along π . Hence $\pi \models Fex_\ell$. (End of case 1.)

Case 2: $\ell \in \psi_{ex}$. Since $\pi \models \overset{\infty}{F}nblk_\ell \wedge \overset{\infty}{F}blk_\ell$, P_ℓ must change state infinitely often along π . Hence by the I -structure definition (5.3.1), P_ℓ must be executed infinitely often along π . Thus $\pi \models \overset{\infty}{F}ex_\ell$, which implies $\pi \models Fex_\ell$. (End of case 2.)

Case 3: $\ell \in \psi_{blk}$. First, a few definitions are needed. A subset ψ of $dom(I)$ is I -connected if and only if $\bigwedge i, j \in \psi, i \neq j. (i I^+ j)$ where I^+ is the transitive closure of I . An I -process P_k borders ψ if and only if $k \in I(i) - \psi$ for some $i \in \psi$. We let $border(\psi)$ denote the set of all bordering I -processes of ψ . Let η be a maximal I -connected subset of ψ_{blk} . We call η a blk -region. Consider the I -processes that border η . Clearly, no such I -process can be in ψ_{blk} , as η would not be maximal. Hence, every bordering I -process must be in ψ_{nblk} or in ψ_{ex} . It is clear that ψ_{blk} can be partitioned into (I -disconnected) blk -regions. Let θ be a maximal I -connected subset of $dom(I)$ such that $\pi \models \bigwedge i \in \theta. (\overset{\infty}{F}ex_i)$. We call θ an inf -region. For an I -process P_i such that $\pi \models \overset{\infty}{F}ex_i$, we define $inf(i)$ to be the inf -region which P_i is a member of. For ψ an arbitrary subset of $dom(I)$, if wait-for-graph $W_I(s)$ contains an edge $\alpha_i^I \rightarrow P_j^I$ such that $i \in \psi$ and $j \notin \psi$, then we say that $W_I(s)$ contains an edge out of ψ .

We now establish a series of assertions P2, P3, P4, which together allow us to establish $\pi \models Fex_\ell$.

Let P_j border some inf -region θ . Then there exists a suffix ρ' of ρ such that, for every state s along ρ' ,

$$W_I(s) \text{ contains no edges into } P_j. \quad (\text{P2})$$

Proof of P2. Since P_j borders inf -region θ , there exists $k \in \theta$ such that $j I k$. Since $j \notin \theta$, we have $\rho \models \overset{\infty}{G}\neg ex_j$, as otherwise θ would not be maximal. Thus there exists a suffix ρ'' of ρ such that

$$\rho'' \models (\overset{\infty}{F}ex_k \wedge G\neg ex_j).$$

Now consider $\rho'' \upharpoonright jk$. By the path projection definition (5.1.2.1) and $\rho'' \models (\overset{\infty}{F}ex_k \wedge G\neg ex_j)$, we have $\rho'' \upharpoonright jk \models (\overset{\infty}{F}ex_k \wedge G\neg ex_j)$. Thus, by the path-mapping lemma (6.4.3), $\rho'' \upharpoonright jk$ is a fullpath in M_{jk} . Also, $\rho'' \upharpoonright jk \models Gex_k$, since all transitions of $\rho'' \upharpoonright jk$ are either P_k^j -transitions or P_j^k -transitions. Hence

$$M_{jk}, s \upharpoonright jk \models EGex_k, \quad (\text{a})$$

where s is an arbitrary state along ρ'' . Note that s is a reachable state in M_I , since it is reachable from v , and v is, by assumption, a reachable state in M_I . Now let i be an arbitrary element of $I(j)$, and let $J = \{\{i, j\}, \{j, k\}\}$. By the state-mapping corollary (6.4.5), $s \upharpoonright J$ is reachable in M_J . Since $(s \upharpoonright J) \upharpoonright jk = s \upharpoonright jk$, we have, by (a)

$$M_{jk}, (s \upharpoonright J) \upharpoonright jk \models EGex_k. \quad (\text{b})$$

Letting $s_J = s \uparrow J$ in LV (Definition 6.7.2.1), we get from (b)

$$\bigwedge a_i^J . (a_i^J \longrightarrow P_j^J \notin W_J(s \uparrow J)), \quad (c)$$

so there is no edge in $W_J(s \uparrow J)$ from i to j . Since $i I j$ and $j I k$, we have $J \subseteq I$. So, by applying the wait-for-graph projection proposition (6.5.4.1) to (c), we obtain $\bigwedge a_i^I . (a_i^I \longrightarrow P_j^I \notin W_I(s))$. Since i is an arbitrarily chosen element of $I(j)$, we conclude that $W_I(s)$ contains no edges into P_j^I . As s is an arbitrary state along ρ'' , setting ρ' to ρ'' concludes the proof of P2. (End proof of P2.)

Let ψ be an arbitrary subset of ψ_{blk} . If there exists a suffix ρ' of ρ such that,

for every state s along ρ' , $W_I(s)$ contains no edge out of ψ ,

then,

$$\rho \models \bigvee i \in \psi . \overset{\infty}{Fex}_i. \quad (P3)$$

Proof of P3. Assume otherwise. Thus there exists a suffix ρ'' of ρ such that $\rho'' \models \bigwedge i \in \psi . G \neg ex_i$. Let s be an arbitrary state along ρ'' . By assumption, $W_I(s)$ is supercycle-free. Hence $W_I(s) \uparrow \psi$ is supercycle-free. Thus, by the supercycle proposition (6.5.5.1) with $I := I \uparrow \psi$, some move node $a_{i_s}^I$ such that $i_s \in \psi$, must have no outgoing edges in $W_I(s) \uparrow \psi$. By assumption, $W_I(s)$ contains no edge from a move in ψ to an I -process outside ψ . Thus $a_{i_s}^I$ must have no outgoing edges in $W_I(s)$. Hence, by Observation 6.5.2.2, $s(a_{i_s}^I.guard) = true$. So, by the compact I -structure definition (5.4.3), $a_{i_s}^I$ can be executed, and thus $M_I, s \models en_{i_s}$.

Now let t be the successor state to s along ρ'' . By assumption, $W_I(t)$ contains no edge out of ψ . Also, $W_I(t) \uparrow \psi = W_I(s) \uparrow \psi$ since no I -process in ψ is executed along ρ'' . Thus $W_I(t)$ contains no edges out of $a_{i_s}^I$. Hence, by Observation 6.5.2.2, $t(a_{i_s}^I.guard) = true$. We can inductively repeat this argument (e.g., for the successor state to t , and then the successor state to that state, etc...) to conclude that $u(a_{i_s}^I.guard) = true$ for every state u which occurs after s along ρ'' . Thus $\rho'' \models \overset{\infty}{Gen}_{i_s}$.

Now $\rho'' \models \overset{\infty}{Gblk}_{i_s}$, since $i_s \in \psi$ and $\psi \subseteq \psi_{blk}$. Since $\rho'' \models G \neg ex_{i_s}$, we have $\rho'' \models \overset{\infty}{G}(blk_{i_s} \wedge en_{i_s}) \wedge G \neg ex_{i_s}$. Hence, by the weak blocking fairness definition (6.7.3.3), $\rho'' \models \neg \Phi$. Hence $\pi \models \neg \Phi$, since ρ'' is a suffix of π . But π was chosen so that $\pi \models \Phi$. Hence the original assumption is false, and P3 is established. (End proof of P3.)

Let j be an arbitrary element of ψ_{blk} . Then either

P_j is a member of some *inf*-region, or

P_j borders some *inf*-region.

(P4)

Proof of P4. Assume otherwise. Since $j \in \psi_{blk}$, we have $j \in \eta$ for some *blk*-region η . Hence

$$\zeta = \eta - (\bigcup_{\theta \text{ is an } \textit{inf}\text{-region}} (\theta \cup border(\theta)))$$

is nonempty, since $j \notin \theta \cup border(\theta)$ for any *inf*-region θ by assumption, and thus $j \in \zeta$. By (P2), we have that there exists a suffix ρ' of ρ such that

for any I -process P_k which borders an *inf*-region,
 for every state s along ρ' ,
 $W_I(s)$ contains no edges into P_k . (a)

Also, if for a state s and an I -process P_k , $s \models nblk_k$, then $W_I(s)$ contains no edge into P_k (see Observation 6.7.3.2). Thus, by definition of ψ_{nblk} we have

for every state s along ρ' ,
 $W_I(s)$ contains no edge into any I -process P_k in ψ_{nblk} . (b)

Now consider an arbitrary member P_k of $border(\zeta)$. Since every I -process is a member of exactly one of ψ_{blk} , ψ_{ex} , ψ_{nblk} , we have three (sub-)cases.

$k \in \psi_{blk}$. Since $\zeta \subseteq \eta$, and $k \in border(\zeta)$, we have $k \in \eta$, since η is a *blk*-region, (and $k \in \psi_{blk}$ by the case assumption) and all *blk*-regions are maximal, by definition. By assumption, $k \in border(\zeta)$, so $k \notin \zeta$ by definition of *border*. Therefore, $k \in \eta - \zeta$. Since $\zeta = \eta - (\bigcup \theta \text{ is an } \textit{inf}\text{-region}(\theta \cup border(\theta)))$ by definition, we have $k \in \theta \cup border(\theta)$ for some *inf*-region θ . Since $k \in border(\zeta)$, there exists P_m such that $k I m$ and $m \in \zeta$. However, if $k \in \theta$, then $m \in border(\theta)$, contrary to the definition of ζ . Thus we conclude that $k \in border(\theta)$.

$k \in \psi_{ex}$. Therefore $\pi \models \overset{\infty}{F}ex_k$ (see case 2). So k is a member of some *inf*-region θ , by definition of *inf*-region. Since $k \in border(\zeta)$, there exists m such that $k I m$ and $m \in \zeta$. Thus $m \in border(\theta)$, contrary to the definition of ζ . We conclude that k cannot be a member of any *inf*-region, and hence k cannot be a member of ψ_{ex} .

$k \in \psi_{nblk}$. We do not need to infer anything for this case other than $k \in \psi_{nblk}$.

Considering the above three (sub-)cases, we have shown that every member of $border(\zeta)$ either borders some *inf*-region or is a member of ψ_{nblk} . Therefore, in $W_I(s)$, every edge out of ζ must have a target P_k such that P_k borders an *inf*-region or such that P_k is in ψ_{nblk} . Thus, by (a) and (b), we conclude that

for every state s along ρ' , $W_I(s)$ contains no edge out of ζ .

Since $\zeta \subseteq \eta$, ζ is a subset of ψ_{blk} by definition of *blk*-region. Thus, by (P3), $\rho' \models \overset{\infty}{F}ex_{k'}$ for some $P_{k'}$ in ζ . Hence $k' \in \theta$ for some *inf*-region θ . But this contradicts $k' \in \zeta$, by definition of ζ . Hence the original assumption is false, and P4 is established. (End proof of P4.)

We can now establish $\pi \models Fex_\ell$. By the case 3 assumption, $\ell \in \psi_{blk}$. By (P4), we have

P_ℓ is a member of some *inf*-region θ or P_ℓ borders some *inf*-region θ .

If P_ℓ is a member of θ , then $\pi \models \overset{\infty}{F}ex_\ell$, and therefore $\pi \models Fex_\ell$. If P_ℓ borders θ , then there exists $j \in I(\ell) \cap \theta$. Hence $\pi \models \overset{\infty}{F}ex_j$, and since π is an infinite fullpath, we conclude by the path projection definition (5.1.2.1) that $\pi \upharpoonright \ell j$ is an infinite path (and therefore is a fullpath) and $\pi \upharpoonright \ell j \models \overset{\infty}{F}ex_j$. Moreover, the first state of $\pi \upharpoonright \ell j$ is $v \upharpoonright \ell j$. Since $\bigwedge k \in I(\ell). (M_{k\ell}, v \upharpoonright k \ell \models \neg r_\ell \wedge AFr_\ell)$ by assumption (3) of the lemma, and $j \in I(\ell)$, we have $M_{\ell j}, v \upharpoonright \ell j \models \neg r_\ell \wedge AFr_\ell$. Thus, $M_{\ell j}, v \upharpoonright \ell j \models AFex_\ell$, and so

$\pi \upharpoonright \ell j \models \text{Fex}_\ell$ since $\pi \upharpoonright \ell j$ is a fullpath and the first state of $\pi \upharpoonright \ell j$ is $v \upharpoonright \ell j$. Hence, by the path projection definition (5.1.2.1), we conclude $\pi \models \text{Fex}_\ell$. (End of case 3.)

Since we have established $\pi \models \text{Fex}_\ell$ in all three cases, the lemma is established. \square

C.6 The Generalized Large Model Theorem

THEOREM 6.8.1 (GENERALIZED LARGE MODEL). *Let f_{kl} be an arbitrary formula of FLCTL_{kl} (Definition 6.6.1). Let s be an arbitrary reachable I -state, and, for all i, j such that $i I j$, let $s_{ij} = s \upharpoonright ij$. If the liveness assumption LV holds, and $W_I(u)$ is supercycle-free for every reachable I -state u , then*

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models \bigwedge_{kl} f_{kl}) \text{ implies } M_I, s \models_{\Phi} \bigwedge_{kl} f_{kl}.$$

PROOF. By MPCTL^* semantics, $M_{ij}, s_{ij} \models \bigwedge_{kl} f_{kl}$ is equivalent to $M_{ij}, s_{ij} \models f_{ij}$. Also by MPCTL^* semantics, $M_I, s \models_{\Phi} \bigwedge_{kl} f_{kl}$ is equivalent to $M_I, s \models_{\Phi} \bigwedge (i, j) \in I. (f_{ij})$. This is equivalent, by CTL^* semantics, to $\bigwedge (i, j) \in I. (M_I, s \models_{\Phi} f_{ij})$. Hence, the generalized large-model theorem is established if we can prove that

$$\bigwedge (i, j) \in I. (M_{ij}, s_{ij} \models f_{ij}) \text{ implies } \bigwedge (i, j) \in I. (M_I, s \models_{\Phi} f_{ij}) \quad (*)$$

given the assumptions of the generalized large-model theorem. The proof is by induction on the structure of f_{ij} .

The proofs of the cases $f_{ij} = h_{ij}$, $f_{ij} = f'_{ij} \wedge f''_{ij}$, $f_{ij} = AGh_{ij}$, $f_{ij} = AG(p_i \Rightarrow AY_i q_i)$, $f_{ij} = AG(a_i \Rightarrow EX_i b_i)$, are verbatim identical to the proofs for the same cases respectively in the large-model theorem (6.6.2), and are thereby omitted here. Note that only the proof for the remaining case of $f_{ij} = AG(p_i \Rightarrow A[q_i Ur_i])$ in Theorem 6.6.2 appealed to the assumptions $I = \{i_1, \dots, i_K\} \times \{i_1, \dots, i_K\} - \{(i, i) \mid i \in \{i_1, \dots, i_K\}\}$ and $M_I, S_I^0 \models AGEX \text{true}$ in the antecedent of Theorem 6.6.2. We now give the proof for $f_{ij} = AG(p_i \Rightarrow A[q_i Ur_i])$ in the generalized case.

$f_{ij} = AG(p_i \Rightarrow A[q_i Ur_i])$. We will establish $M_I, t \models_{\Phi} (p_i \Rightarrow A[q_i Ur_i])$ where t is an arbitrary s -reachable state in M_I . If $M_I, t \models \neg p_i$ then $M_I, t \models_{\Phi} (p_i \Rightarrow A[q_i Ur_i])$, and we are done. Otherwise $M_I, t \models p_i$, and we must establish $M_I, t \models_{\Phi} A[q_i Ur_i]$. If $M_I, t \models r_i$ then we are done. Otherwise $M_I, t \models \neg r_i$, so $t \upharpoonright ij \models \neg r_i$ by the I -state projection proposition (6.3.1). Let π be an arbitrary fullpath of M_I starting in t such that $\pi \models \Phi$. The antecedent is

$$M_{ij}, s_{ij} \models AG(p_i \Rightarrow A[q_i Ur_i]). \quad (a)$$

By assumption, $s \upharpoonright ij = s_{ij}$. Thus by the relativized state-mapping corollary (6.4.6) with $J = \{\{i, j\}\}$, $t \upharpoonright ij$ is a s_{ij} -reachable state in M_{ij} . So, by (a) we have

$$M_{ij}, t \upharpoonright ij \models (p_i \Rightarrow A[q_i Ur_i]). \quad (b)$$

Since $M_I, t \models p_i$, we conclude $M_{ij}, t \upharpoonright ij \models p_i$ by the I -state projection proposition (6.3.1). Together with (b), this yields

$$M_{ij}, t \upharpoonright ij \models A[q_i Ur_i], \quad (c)$$

and since $t \upharpoonright ij \models \neg r_i$, we have by (c) and CTL^* semantics

$$M_{ij}, t \upharpoonright ij \models \neg r_i \wedge AF r_i. \quad (d)$$

By the path-mapping lemma (6.4.3), $\pi \upharpoonright ij$ is a path in M_{ij} . Also, $\pi \upharpoonright ij$ starts in $t \upharpoonright ij$, and therefore, by (c) and CTL* semantics, we have $M_{ij}, \pi \upharpoonright ij \models [q_i U_w r_i]$ (we cannot conclude $M_{ij}, \pi \upharpoonright ij \models [q_i U r_i]$ because we have not shown that $\pi \upharpoonright ij$ is a fullpath). By the path projection lemma (6.3.3) with $J = \{\{i, j\}\}$, we have $M_I, \pi \models [q_i U_w r_i]$. It remains for us to establish $M_I, \pi \models Fr_i$.

Since j ranges over $I(i)$, we have, from (d), $\bigwedge j \in I(i). (M_{ij}, t \upharpoonright ij \models \neg r_i \wedge A Fr_i)$. By assumption, s is reachable, and t is s -reachable, so t is reachable. Together with the assumptions of the theorem, we have satisfied the antecedent of the progress lemma (6.7.4.1) for $v = t$ and $\ell = i$. Thus, by the progress lemma (6.7.4.1), we have $M_I, t \models_{\Phi} A F ex_i$. Therefore, P_i^I is eventually executed along π . By repeating this argument inductively, we conclude that either P_i^I is executed repeatedly along π until a global state t' (along π) is reached such that $M_I, t' \models r_i$, (and hence $M_I, \pi \models Fr_i$), or P_i^I is executed infinitely often along π (since $\neg r_i \wedge A Fr_i$ continues to hold, and therefore the progress lemma can be applied repeatedly). In the latter case, we have that π contains an infinite number of P_i^I -transitions, so by definition of path projection, $\pi \upharpoonright ij$ contains an infinite number of P_i^j -transitions, so $\pi \upharpoonright ij$ is a fullpath. Since $\pi \upharpoonright ij$ is a fullpath starting in $t \upharpoonright ij$, we have, by (c), $M_{ij}, \pi \upharpoonright ij \models [q_i U r_i]$. By the path projection lemma (6.3.3), we have $M_I, \pi \models [q_i U r_i]$, which implies $M_I, \pi \models Fr_i$.

Since $M_I, \pi \models Fr_i$ in both cases, and $M_I, \pi \models [q_i U_w r_i]$ has been established above, we have $M_I, \pi \models [q_i U r_i]$. Since π is an arbitrary fullpath starting in t such that $\pi \models \Phi$, we conclude $M_I, t \models_{\Phi} A[q_i U r_i]$. Hence $M_I, t \models_{\Phi} (p_i \Rightarrow A[q_i U r_i])$. Since t is an arbitrary s -reachable state in M_I , we conclude $M_I, s \models_{\Phi} AG(p_i \Rightarrow A[q_i U r_i])$. Since i ranges over $\{i_1, \dots, i_K\}$, and j ranges over $I(i)$, we have $\bigwedge (i, j) \in I. (M_I, s \models_{\Phi} AG(p_i \Rightarrow A[q_i U r_i]))$. Thus, we have established (*), which concludes the proof of the generalized large-model theorem. \square

D. A COMPACT REPRESENTATION FOR SYNCHRONIZATION SKELETONS

In this appendix, we provide a full discussion of the compact representation introduced in Section 5.4. Our discussion here is self-contained and so repeats some of the material in Section 5.4 (in particular, the definition of compact MP-synthesis is repeated, and so retains the number, 5.4.1, that it has in the main text).

Suppose P_i^j (for every $j \in I(i)$) contains two arcs from i -state s_i to i -state s'_i , e.g., $a_{i,1}^j = (s_i, B_{i,1}^j \rightarrow A_{i,1}^j, s'_i)$ and $a_{i,2}^j = (s_i, B_{i,2}^j \rightarrow A_{i,2}^j, s'_i)$. Then, by the MP-synthesis definition (5.1.1), P_i^I contains $2^{|I(i)|}$ arcs from i -state s_i to i -state s'_i , one arc for each element of the cartesian product

$$\{a_{i,1}^{j_1}, a_{i,2}^{j_1}\} \times \dots \times \{a_{i,1}^{j_n}, a_{i,2}^{j_n}\}$$

where $\{j_1, \dots, j_n\} = I(i)$. Thus P_i^I is exponentially large in K ($= |dom(I)|$) in the worst case (since $|I(i)| = K - 1$ when $i I j$ for every j in $dom(I) - \{i\}$), which defeats the purpose of MP-synthesis. We deal with this by defining a compact representation for processes in which there is at most one arc between any pair of (local) states, thereby avoiding the exponential blowup illustrated above.

Consider a pair-process P_i^j which has two arcs from state s_i to state s'_i , labeled with the synchronization commands $B_{i,1}^j \rightarrow A_{i,1}^j, B_{i,2}^j \rightarrow A_{i,2}^j$. In compact notation, we replace these two arcs by a single arc whose label is $B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j$.

The symbol \oplus is a binary operator which takes a pair of guarded commands as arguments. It is defined as follows:

$$(B_1 \rightarrow A_1) \oplus (B_2 \rightarrow A_2) \stackrel{\text{df}}{=} B_1 \vee B_2 \rightarrow \begin{array}{l} \text{if } B_1 \rightarrow A_1 \\ \square B_2 \rightarrow A_2 \\ \text{fi} \end{array}$$

Roughly, the operational semantics of $B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j$ is that if one of the guards $B_{i,1}^j, B_{i,2}^j$ evaluates to true, then the corresponding body $A_{i,1}^j, A_{i,2}^j$, respectively, can be executed. If neither $B_{i,1}^j$ nor $B_{i,2}^j$ evaluates to true, then the command “blocks,” i.e., waits until one of $B_{i,1}^j, B_{i,2}^j$ evaluates to true. Note that the guarded commands which we use as arc labels are, in general, *partial commands*, i.e., they cannot be executed in every global state. A full treatment of the calculus of partial guarded commands is beyond our scope here. The reader is referred to Nelson [1989]. It is easily seen that \oplus is commutative. To see that it is also associative, we note that

$$(B_1 \rightarrow A_1 \oplus B_2 \rightarrow A_2) \oplus B_3 \rightarrow A_3$$

and

$$B_1 \rightarrow A_1 \oplus (B_2 \rightarrow A_2 \oplus B_3 \rightarrow A_3)$$

have the same semantics, namely, if one of B_1, B_2, B_3 is true, then the corresponding body can be executed, and if none of B_1, B_2, B_3 are true, then the command blocks. Since \oplus is commutative and associative, it can be extended to n arguments using the indexed notation $\oplus_{\ell \in [1:n]}$. Thus, if P_i^j contains n arcs from i -state s_i to i -state s'_i , with labels $B_{i,1}^j \rightarrow A_{i,1}^j, \dots, B_{i,n}^j \rightarrow A_{i,n}^j$, then these n arcs can be replaced by a single arc whose label is $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$.

We call an arc whose label has the form $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$ a (*pair*) *move*. In compact notation, a pair-process has at most one move between any pair of local states. The translation from compact notation back to normal notation is straightforward: simply replace every move by the corresponding set of arcs. The operational semantics is as follows. Assume that the current state is $(s_1, \dots, s_i, \dots, s_K, x_1, \dots, x_m)$, and that P_i contains a move from s_i to s'_i labeled by $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$. If a guard $B_{i,\ell}^j$ (for some ℓ in $[1 : n]$) evaluates to true in the current state, then $(s_1, \dots, s'_i, \dots, s_K, x'_1, \dots, x'_m)$ is a permissible next-state where x'_1, \dots, x'_m is a list of updated shared variables resulting from action $A_{i,\ell}^j$. A computation path is an infinite sequence of states where successive pairs of states are related by the above next-state relation. As before, omission of the guard $B_{i,\ell}^j$ from a guarded command means that $B_{i,\ell}^j$ is interpreted as *true*, and we write the command as $A_{i,\ell}^j$, while omission of the action $A_{i,\ell}^j$ from a guarded command means that the shared variables are unaltered, and we write the command as $B_{i,\ell}^j$. This is formalized by the following definition, which is a consequence of the pair structure definition (5.2.1) and the translation between compact and normal notation given above.

Definition 5.4.2 (Compact Pair-Structure). Let $i I j$. The semantics of $(S_{ij}^0, P_i^j \| P_j^i)$ in compact notation is given by the pair-structure $M_{ij} = (S_{ij}^0, S_{ij}, R_{ij})$ where

- (1) S_{ij} is a set of ij -states,
- (2) $S_{ij}^0 \subseteq S_{ij}$ gives the initial states of $P_i^j \| P_j^i$, and
- (3) $R_{ij} \subseteq S_{ij} \times \{i, j\} \times S_{ij}$ is a transition relation giving the transitions of $P_i^j \| P_j^i$.
A transition (s_{ij}, h, t_{ij}) by $P_h^{\bar{h}}$ is in R_{ij} if and only if
 - (a) $h \in \{i, j\}$,
 - (b) s_{ij} and t_{ij} are ij -states, and
 - (c) there exists a move $(s_{ij} \uparrow h, \oplus_{\ell \in [1:n]} B_{h,\ell}^{\bar{h}} \rightarrow A_{h,\ell}^{\bar{h}}, t_{ij} \uparrow h)$ in $P_h^{\bar{h}}$ such that there exists $m \in [1 : n]$:
 - (i) $s_{ij}(B_{h,m}^{\bar{h}}) = \text{true}$,
 - (ii) $\langle s_{ij} \uparrow \mathcal{H}_{ij} \rangle A_{h,m}^{\bar{h}} \langle t_{ij} \uparrow \mathcal{H}_{ij} \rangle$, and
 - (iii) $s_{ij} \uparrow \bar{h} = t_{ij} \uparrow \bar{h}$.

Here $\bar{h} = i$ if $h = j$ and $\bar{h} = j$ if $h = i$.

Now an I -process P_i^I is derived by “composing” all the moves of P_i^j (as j varies over $I(i)$) which have the same start and end states. Toward this end, we define the binary composition operator \otimes on guarded commands. We say that a guarded command is *simple* iff it has the form $B \rightarrow A$ where B is a guard and A is a parallel assignment statement. Applied to a pair of simple guarded commands, \otimes returns a simple guarded command whose guard is the conjunction of the guards of the operands and whose assignment statement is the parallel composition of the assignment statements of the operands.

Definition D.1 (Simple Guarded-Command Composition). The composition of two simple guarded commands is given by

$$(B_1 \rightarrow A_1) \otimes (B_2 \rightarrow A_2) = B_1 \wedge B_2 \rightarrow A_1 // A_2.$$

That is, if both guards are true, then both bodies can be executed in parallel. We note that, in general, it is possible that A_1, A_2 have a variable in common on their left-hand sides. Since \otimes is a *syntactic* operator, this presents no problems in the definition of \otimes . However, the semantics of $A_1 // A_2$ is problematic in general. But, the only time when we need to assign a semantics to guarded commands containing \otimes is in Definition 5.4.3 below, and we shall see that in this case, the possibility of assigning twice to the same variable within a parallel assignment does not arise. It is clear that \otimes , when applied to simple guarded commands, is commutative and associative, since both \wedge and $//$ are. We now define general guarded commands as those that can be built up from simple guarded commands by applying \oplus and \otimes .

Definition D.2 (General Guarded Command). General guarded commands are inductively defined as follows.

- (1) A simple guarded command is a general guarded command
- (2) If G_1, G_2 are general guarded commands, then so are $G_1 \oplus G_2$ and $G_1 \otimes G_2$
- (3) The only general guarded commands are those that are generated by rules (1) and (2) above

In order to define the operational semantics of a general guarded command, we introduce a *normal form* for general guarded commands.

Definition D.3 (Normal Form). A general guarded command is in normal form iff it has the form $\oplus_{\ell \in [1:n]} G_\ell$, where each G_ℓ is a simple guarded command.

We note that the operational semantics of normal forms has been described informally above. The compact pair-structure definition (5.4.2) formalizes the operational semantics of normal forms by defining the pair-structures that are generated by pair-processes expressed in compact notation (whose arc labels are general guarded commands expressed in normal form). We can now define the application of \otimes to general guarded commands in normal form as follows.

Definition D.4 (Normal Form Composition). The composition of two general guarded commands in normal form is given by

$$(\oplus_{\ell \in \varphi} G_\ell) \otimes (\oplus_{k \in \psi} G_k) = \oplus_{\ell \in \varphi, k \in \psi} (G_\ell \otimes G_k).$$

We recall that $G_\ell \otimes G_k$ is given by the simple guarded-command composition definition (D.1), since G_ℓ, G_k are simple guarded commands.

It remains to show that every guarded command can be expressed in normal form.

PROPOSITION D.5 (NORMAL FORM). *Every general guarded command can be expressed in normal form.*

PROOF. By structural induction over the definition of a general guarded command G .

Base Case. G is simple. Then $G = \oplus_{\ell \in \varphi} G_\ell$ where φ is a singleton set. This is in normal form.

Induction Step. $G = G_1 \oplus G_2$. By the induction hypothesis, G_1 and G_2 are in normal form. Let $G_1 = \oplus_{\ell \in \varphi} G_\ell$, $G_2 = \oplus_{k \in \psi} G_k$. Hence G_ℓ for all $\ell \in \varphi$ and G_k for all $k \in \psi$ are simple guarded commands, by the normal-form definition (D.3). Since \oplus is associative and commutative over simple guarded commands, we have $G_1 \oplus G_2 = (\oplus_{\ell \in \varphi} G_\ell) \oplus (\oplus_{k \in \psi} G_k) = (\oplus_{m \in \varphi \cup \psi} G_m)$. This last expression is in normal form.

Induction Step. $G = G_1 \otimes G_2$. By the induction hypothesis, G_1 and G_2 are in normal form. Let $G_1 = \oplus_{\ell \in \varphi} G_\ell$, $G_2 = \oplus_{k \in \psi} G_k$. Hence, by the normal-form composition definition (D.4), $G_1 \otimes G_2 = \oplus_{\ell \in \varphi, k \in \psi} (G_\ell \otimes G_k)$. Now G_ℓ for all $\ell \in \varphi$ and G_k for all $k \in \psi$ are simple guarded commands, by the normal-form definition (D.3). Thus, by the simple guarded-command composition definition (D.1), $G_\ell \otimes G_k$ can be rewritten as a simple guarded command. Thus by the normal-form definition (D.3), $G_1 \otimes G_2$ can be expressed in normal form. \square

We finally recast the MP-synthesis definition (5.1.1) into compact form as follows.

Definition 5.4.1 (Compact MP-Synthesis). A compact I -process P_i^I is derived from the compact pair-process P_i^j as follows:

P_i^I contains a move from s_i to t_i with label $\otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$
iff
for every j in $I(i)$: P_i^j contains a move from s_i to t_i with label $\oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$.

The initial state set S_I^0 of the I -system is derived from the initial state S_{ij}^0 of the pair-system as follows:

$$S_I^0 = \{s \mid \bigwedge (i, j) \in I. (s \upharpoonright ij \in S_{ij}^0)\}.$$

Thus an I -process in compact notation has at most one move of the form $(s_i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t_i)$ between any pair of i -states s_i, t_i . Note that a move in a pair-process is simply a special case of a move in an I -process when I is a single pair, e.g., if $I = \{\{i, k\}\}$, then $I(i)$ in $\otimes_{j \in I(i)}$ expands to the singleton $\{k\}$ giving a move in the pair-process P_i^k . In the sequel, we shall use a_i^j, a_i^I to denote moves in P_i^j, P_i^I , respectively.

Returning to the example given at the beginning of this section, the pair of arcs $(s_i, B_{i,1}^j \rightarrow A_{i,1}^j, s'_i), (s_i, B_{i,2}^j \rightarrow A_{i,2}^j, s'_i)$ in P_i^j are replaced by the single move $(s_i, (B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j), s'_i)$, and the $2^{|I(i)|}$ arcs in P_i^I are replaced by the single move $(s_i, \otimes_{j \in I(i)} (B_{i,1}^j \rightarrow A_{i,1}^j \oplus B_{i,2}^j \rightarrow A_{i,2}^j), s'_i)$.

The translation of I -processes from normal to compact notation is most easily achieved via the translation given above for pair-processes, i.e., derive the corresponding pair-process (essentially applying MP-synthesis “in reverse”), perform the translation to obtain the compact pair-processes, and then derive the compact I -process using the compact MP-synthesis definition (5.4.1). Translating from compact to normal notation is achieved by applying the definition of the \otimes operator, in effect expanding the general guarded command $\otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_\ell \rightarrow A_\ell$ into normal form, and then creating one arc for each simple guarded command in the normal form. For example, if $I = \{\{i, k\}, \{k, \ell\}, \{\ell, i\}\}$, then

$$\otimes_{j \in I(i)} (T_j \rightarrow x_{ij} := j \oplus N_j \vee C_j \rightarrow skip)$$

expands into normal form as follows. First, we expand the bound variable j over its range $\{k, \ell\}$, thereby replacing the indexed form $\otimes_{j \in I(i)}$ by the infix form \otimes , to obtain

$$\begin{aligned} & (T_k \rightarrow x_{ik} := k \oplus N_k \vee C_k \rightarrow skip) \otimes \\ & (T_\ell \rightarrow x_{i\ell} := \ell \oplus N_\ell \vee C_\ell \rightarrow skip). \end{aligned}$$

Now we apply the normal-form composition definition (D.4) to obtain

$$\begin{aligned} & (T_k \rightarrow x_{ik} := k \otimes T_\ell \rightarrow x_{i\ell} := \ell) \oplus \\ & (N_k \vee C_k \rightarrow skip \otimes T_\ell \rightarrow x_{i\ell} := \ell) \oplus \\ & (T_k \rightarrow x_{ik} := k \otimes N_\ell \vee C_\ell \rightarrow skip) \oplus \\ & (N_k \vee C_k \rightarrow skip \otimes N_\ell \vee C_\ell \rightarrow skip). \end{aligned}$$

Finally, we apply the simple guarded-command composition definition (D.1) to all four occurrences of \otimes , to obtain

$$T_k \wedge T_\ell \rightarrow x_{ik} := k / x_{i\ell} := \ell \oplus$$

$$\begin{aligned}
 (N_k \vee C_k) \wedge T_\ell &\rightarrow \text{skip} // x_{i\ell} := \ell \oplus \\
 T_k \wedge (N_\ell \vee C_\ell) &\rightarrow x_{ik} := k // \text{skip} \oplus \\
 (N_k \vee C_k) \wedge (N_\ell \vee C_\ell) &\rightarrow \text{skip} // \text{skip}.
 \end{aligned}$$

It is easy to see that if P_i^I had three neighbors instead of two, that the size of the final result would be eight instead of four. Generalizing, we see that if $|I(i)| = m$, then

$$\otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j$$

expands into a term with size on the order of n^m . Thus the compact form provides an exponential savings in the size of the representation of the I -processes.

The operational semantics of compact I -processes is given by the following definition, which is a consequence of the I -structure definition (5.3.1) and the translation between compact and normal notation given above.

Definition 5.4.3 (Compact I-Structure). The semantics of $(S_I^0, P_{i_1}^I \parallel \dots \parallel P_{i_K}^I)$ in compact notation is given by the I -structure $M_I = (S_I^0, S_I, R_I)$ where

- (1) S_I is a set of I -states,
- (2) $S_I^0 \subseteq S_I$ gives the initial states of $P_{i_1}^I \parallel \dots \parallel P_{i_K}^I$, and
- (3) $R_I \subseteq S_I \times \text{dom}(I) \times S_I$ is a transition relation giving the transitions of $P_{i_1}^I \parallel \dots \parallel P_{i_K}^I$. A transition (s, i, t) by P_i^I is in R_I if and only if
 - (a) $i \in \text{dom}(I)$,
 - (b) s and t are I -states, and
 - (c) there exists a move $(s \uparrow i, \otimes_{j \in I(i)} \oplus_{\ell \in [1:n]} B_{i,\ell}^j \rightarrow A_{i,\ell}^j, t \uparrow i)$ in P_i^I such that all of the following hold:
 - (i) for all j in $I(i)$, there exists $m \in [1 : n]$:
 $s \uparrow ij (B_{i,m}^j) = \text{true}$ and $< s \uparrow \mathcal{SH}_{ij} > A_{i,m}^j < t \uparrow \mathcal{SH}_{ij} >$,
 - (ii) for all j in $\text{dom}(I) - \{i\}$: $s \uparrow j = t \uparrow j$, and
 - (iii) for all j, k in $\text{dom}(I) - \{i\}$, $j I k$: $s \uparrow \mathcal{SH}_{jk} = t \uparrow \mathcal{SH}_{jk}$.

Given a pair-system $(S_{ij}^0, P_i^j \parallel P_j^i)$ in normal notation, we can apply Definitions 5.2.1, 5.1.1, and 5.3.1 to generate the pair-structure, the I -system, and the I -structure (for the generated I -system) respectively which correspond to $(S_{ij}^0, P_i^j \parallel P_j^i)$. Alternatively, if $(S_{ij}^0, P_i^j \parallel P_j^i)$ is given in compact notation, then we can apply Definitions 5.4.2, 5.4.1, and 5.4.3 to generate the pair-structure, the I -system (in compact notation), and the I -structure (for the generated I -system), respectively which correspond to $(S_{ij}^0, P_i^j \parallel P_j^i)$. It should be apparent that the pair, K -process structures generated by one set of definitions are identical to those generated by the other set of definitions. Thus, any result established in the sequel using one set of definitions will be equally applicable to the other set of definitions. Hence, in establishing a particular result, we will use whichever set of definitions is more convenient at the time. In particular, although some results will be established using normal notation, the implementation of MP-synthesis will be in compact notation, so as to avoid the exponential size I -processes which the normal notation may produce.

E. CHECKING THE WAIT-FOR-GRAPH ASSUMPTION AND THE LIVENESS ASSUMPTION

E.1 Checking the Wait-for-Graph Assumption

The wait-for-graph assumption WG is mechanically checked as follows. As stated in Section 5 we are initially given a pair-system $(S_{k\ell}^0, P_k^\ell \| P_\ell^k)$. For technical convenience, we use the process indices k, ℓ rather than i, j . Using the pair-structure definition (5.2.1) we generate $M_{k\ell}^r$ (recall that a superscript r denotes reachable states/structures—see Section 6.1) from $(S_{k\ell}^0, P_k^\ell \| P_\ell^k)$. We also translate P_k^ℓ, P_ℓ^k to compact notation as shown in Section 5.4. From $M_{k\ell}^r$, we determine the set S_k^r of reachable k -states in $M_{k\ell}^r$.¹⁴ For every $t_k \in S_k^r$, we determine $|t_k.moves|$, using the compact form of P_k^ℓ . Now the wait-for-graph assumption (Definition 6.5.4.2) specifies that J has the form $\{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_n\}\}$ where $n = |t_k.moves|$ and $k \notin \{j, \ell_1, \dots, \ell_n\}$. Since no constraint is placed on the equality of members of $\{j, \ell_1, \dots, \ell_n\}$, we must consider all possible cases for the form of J . We therefore define:

$$\begin{aligned} \mathcal{J}(t_k) &= \{J \mid J = \{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_m\}\} \text{ and} \\ &\quad m \in [1 : n] \text{ and} \\ &\quad j, k, \ell_1, \dots, \ell_m \text{ are pairwise distinct}\} \\ \mathcal{J}'(t_k) &= \{J \mid J = \{\{k, \ell_1\}, \dots, \{k, \ell_m\}\} \text{ and} \\ &\quad m \in [1 : n] \text{ and} \\ &\quad k, \ell_1, \dots, \ell_m \text{ are pairwise distinct}\} \end{aligned}$$

$\mathcal{J}(t_k) \cup \mathcal{J}'(t_k)$ is the set of all the distinct forms of J that must be considered when checking WG for the k -state t_k , with $\mathcal{J}(t_k)$ containing all the forms of J in which $j \notin \{\ell_1, \dots, \ell_m\}$, and $\mathcal{J}'(t_k)$ containing all the forms of J in which $j \in \{\ell_1, \dots, \ell_m\}$. Note that m is really the number of distinct indices in $\{\ell_1, \dots, \ell_n\}$, so it ranges over $[1 : n]$. For every J in $\mathcal{J}(t_k)$, we generate M_J^r , using the compact MP-synthesis definition (5.4.1) and the compact I -structure definition (5.4.3). Then, for every J -state t_J such that $t_J \uparrow k = t_k$ and $s_J \xrightarrow{k} t_J \in R_J$ for some reachable J -state s_J of M_J , we evaluate

$$\bigwedge a_j^J . (a_j^J \longrightarrow P_k^J \notin W_J(t_J))$$

or

$$\bigvee a_k^J \in W_J(t_J) . (\bigwedge \ell \in \{\ell_1, \dots, \ell_m\} . (a_k^J \rightarrow P_\ell^J \notin W_J(t_J))).$$

Using the wait-for-graph definition (6.5.2.1) and CTL* semantics, we can rewrite this as follows:

$$\begin{aligned} t_J \models & \left(\bigwedge a_j^J . (\{a_j^J.start\} \Rightarrow a_j^J.guard_k) \right. \\ & \bigvee \\ & \left. \bigvee a_k^J . (\{a_k^J.start\} \wedge (\bigwedge \ell \in \{\ell_1, \dots, \ell_m\} . a_k^J.guard_\ell)) \right) \end{aligned} \quad (a)$$

Since the formula on the right-hand side of the \models in (a) is purely propositional,

¹⁴This step, in effect, enforces our assumption (made in Section 6.1) that every i -state s_i of P_i^j is reachable in M_{ij} .

it is straightforward to evaluate (a) using the inductive definition for \models supplied in Section 1. If, for any t_k , J , and t_J , (a) evaluates to false, then WG does not hold. Likewise, for every J in $\mathcal{J}'(t_k)$, we generate M_J^r . Then, we set j equal to an arbitrarily chosen member of $\{\ell_1, \dots, \ell_m\}$. Since J is “radially symmetric” with respect to the ℓ_1, \dots, ℓ_m (i.e., J is a “star” with k as the center and ℓ_1, \dots, ℓ_m as the points), the particular choice of member of $\{\ell_1, \dots, \ell_m\}$ makes no difference to the final outcome of the check. As in the case of $J \in \mathcal{J}(t_k)$, we evaluate (a), and if, for any t_k , J and t_J , (a) evaluates to false, then WG does not hold.

If, during the execution of the procedure outlined above, no t_k , J , and t_J are found for which (a) evaluates to false, then WG holds, and we can thus conclude that all I -systems are deadlock-free (provided that $W_I(s_I^0)$ is supercycle-free for every $s_I^0 \in S_I^0$). We summarize the procedure given above in Figure 16, and compute its time complexity. The procedure clearly terminates, since the range of all loop variables is finite. Furthermore, upon termination, WG holds if and only if $WGflag$ is set to “true.”

We use the notation $|S|$ to denote the number of bits needed to represent S using a “straightforward” encoding scheme. By definition of $M_{k\ell}^r$, we have that $|M_{k\ell}^r|$ is $O(|S_{k\ell}^0| + |S_{k\ell}^r| + |R_{k\ell}^r|)$. Since the pair-system is assumed to be nonterminating (Section 6.5), $R_{k\ell}^r$ is total, so we have $|S_{k\ell}^0| \leq |R_{k\ell}^r|$ and $|S_{k\ell}^r| \leq |R_{k\ell}^r|$. Hence $|M_{k\ell}^r|$ is $O(|R_{k\ell}^r|)$. We consider each step of the procedure and compute its worst-case time complexity. Step 0 can be performed in constant time. Step 1 requires the construction of $R_{k\ell}^r$. We generate $R_{k\ell}^r$ incrementally by starting with $S_{k\ell}^0$, selecting some state $s_{k\ell} \in S_{k\ell}^0$, and “expanding” $s_{k\ell}$ by computing all the transitions with source state $s_{k\ell}$ using the compact pair structure definition (5.4.2). We then mark $s_{k\ell}$ as “expanded” and repeat the process until all reachable states have been marked. Each reachable state (and each reachable transition) need be inspected only once. The “expansion” of each state using Definition 5.2.1 requires access to P_k^ℓ and P_ℓ^k . Hence the time complexity of step 1 is $O(|P_k^\ell| \cdot |R_{k\ell}^r|)$.

From the definition of the compact notation given in Section 5.4, we see that the complexity of step 2 (converting P_k^ℓ, P_ℓ^k to compact notation) is $O(|P_k^\ell|)$.

Step 3 can be performed by a single traversal of $R_{k\ell}^r$, and so its complexity is $O(|R_{k\ell}^r|)$.

The complexity of step 4.1 is $O(n)$, as we must count all the moves in $t_k.moves$. The complexity of step 4.2 is simply $O(\sum_{J \in \mathcal{J}(t_k)} |J|)$. Since $|J| = m + 1$, this can be rewritten as $O(\sum_{1 \leq m \leq n} m)$, i.e., $O(n^2)$.

We evaluate the complexity of step 4.3 starting with the most deeply nested iterations and proceeding outward. Now (a) can be rewritten as

$$(t_J \uparrow j k \models \bigwedge a_j^J \cdot (\{a_j^J.start\} \Rightarrow a_j^J.guard_k))$$

or

$$\bigvee a_k^J \cdot ((t_J \uparrow j \models \{a_j^J.start\}) \text{ and } (\bigwedge \ell \in \{\ell_1, \dots, \ell_m\} \cdot (t_J \uparrow j \ell \models a_k^J.guard_\ell))).$$

Hence the complexity of evaluating (a) is $O(sz2 \cdot |P_k^\ell| + m \cdot sz2 \cdot |P_k^\ell|)$, where $sz2$ is the number of bits needed to represent a single $k\ell$ -state. This is just $O(m \cdot sz2 \cdot |P_k^\ell|)$. Since step 4.3.2 requires the inspection of all states in S_J^r (but *not* all transitions in R_J^r), the complexity of step 4.3.2 is $O(m \cdot |S_J^r| \cdot |P_k^\ell|)$. Step 4.3.1 is carried out in an analogous manner to step 1. Each reachable state (and each reachable transition) in R_j^r is inspected only once. The “expansion” of each state (using the compact

0. $WGflag := true$;
1. generate $M_{k\ell}^r$ from $(S_{k\ell}^0, P_k^\ell \parallel P_\ell^k)$;
2. translate P_k^ℓ, P_ℓ^k into compact notation;
3. $S_k^r := \{t_k \mid \bigvee s_{k\ell} \in S_{k\ell}^r \cdot (s_{k\ell} \upharpoonright k = t_k)\}$;
4. for all t_k in S_k^r
 - 4.1 $n := |t_k.moves|$;
 - 4.2 $\mathcal{J}(t_k) := \{J \mid J = \{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_m\}\} \text{ and } m \in [1 : n] \text{ and } j, k, \ell_1, \dots, \ell_m \text{ are pairwise distinct}\}$;
 - 4.3 for all J in $\mathcal{J}(t_k)$
 - 4.3.1 generate M_J^r ;
 - 4.3.2 for all t_J such that $t_J \upharpoonright k = t_k$ and $s_J \xrightarrow{k} t_J \in R_J^r$ for some s_J
 - if (a) evaluates to false, then $WGflag := false$;
 - 4.4 $\mathcal{J}'(t_k) := \{J \mid J = \{\{k, \ell_1\}, \dots, \{k, \ell_m\}\} \text{ and } m \in [1 : n] \text{ and } k, \ell_1, \dots, \ell_m \text{ are pairwise distinct}\}$;
 - 4.5 for all J in $\mathcal{J}'(t_k)$
 - 4.5.1 generate M_J^r ;
 - 4.5.2 set j equal to an arbitrarily selected member of $\{\ell_1, \dots, \ell_m\}$;
 - 4.5.3 for all t_J such that $t_J \upharpoonright k = t_k$ and $s_J \xrightarrow{k} t_J \in R_J^r$ for some s_J
 - if (a) evaluates to false, then $WGflag := false$

Fig. 16. Procedure to check the wait-for-graph assumption.

I -structure definition (5.4.3) with $I := J$ now) requires access to P_k^J and P_ℓ^J (for $\ell \in \{j, \ell_1, \dots, \ell_m\}$). Thus the complexity of step 4.3.1 is $O((|P_k^J| + m \cdot |P_\ell^J|) \cdot |R_J^r|)$. In compact notation, P_k^J is derived by applying the operator “ $\otimes_{\ell \in J(k)}$ ” to the label of every move in P_k^ℓ . Hence $|P_k^J|$ is $O(|J| \cdot |P_k^\ell|)$, which is $O(m \cdot |P_k^\ell|)$ since $|J|$ is $O(m)$. $|P_\ell^J|$ (for $\ell \in \{j, \ell_1, \dots, \ell_m\}$) is $O(|P_\ell^k|)$, since P_ℓ^J has only one J -neighbor, namely P_k^J . Hence the complexity of step 4.3.1 can be rewritten as $O(m \cdot |P_k^\ell| \cdot |R_J^r|)$. So the complexity of steps 4.3.1 and 4.3.2 combined is $O(m \cdot |P_k^\ell| \cdot |R_J^r| + m \cdot |P_k^\ell| \cdot |S_J^r|)$. Since $|S_J^r| \leq |R_J^r|$, the first summand is not less than the second, so this is $O(m \cdot |P_k^\ell| \cdot |R_J^r|)$.

Now R_J^r can be regarded as a “product” of the $m + 1$ transition sets $R_{k\ell}$ for $\ell \in \{j, \ell_1, \dots, \ell_n\}$, as given by the transition-mapping lemma (6.4.1). Thus, $|R_J^r|$ is $O(|R_{k\ell}^r|^{m+1})$. However, this counts the process P_k^J $m + 1$ times instead of once, so we can improve this upper bound to $O(|R_{k\ell}^r|^{m+1} / |P_k^\ell|^m)$. We rewrite this as $O(|R_{k\ell}^r| \cdot \alpha^m)$, where $\alpha = |R_{k\ell}^r| / |P_k^\ell|$. So the complexity of steps 4.3.1 and 4.3.2 combined is rewritten as $O(m \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^m)$.

Step 4.3 iterates steps 4.3.1 and 4.3.2 over all J in $\mathcal{J}(t_k)$, so its complexity is $O(\sum_{J \in \mathcal{J}(t_k)} m \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^m)$. Now J determines m , since $m = (\text{the number of pairs in } J) - 1$. So as J varies from $\{\{j, k\}, \{k, \ell_1\}\}$ to $\{\{j, k\}, \{k, \ell_1\}, \dots, \{k, \ell_n\}\}$, m will vary from 1 to n . Thus the complexity of step 4.3 is $O(|P_k^\ell| \cdot |R_{k\ell}^r| \cdot \sum_{1 \leq m \leq n} m \cdot \alpha^m)$, since $|P_k^\ell|$ and $|R_{k\ell}^r|$ are independent of J . $\sum_{1 \leq m \leq n} m \cdot \alpha^m$ is bounded from above by $n \cdot \sum_{1 \leq m \leq n} \alpha^m$, and $\sum_{1 \leq m \leq n} \alpha^m = (\alpha^{n+1} - 1) / (\alpha - 1)$, which is approximately $O(\alpha^n)$ when $\alpha \gg 1$, as is usually the case. Hence the complexity of step 4.3 is $O(n \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^n)$. Finally, the complexity of steps 4.4 and 4.5 is easily seen to be at most equal to the complexity of steps 4.2, 4.3 respectively, since the only difference is that J is smaller by one pair. Thus, the overall complexity of the body of step 4 (i.e., steps 4.1–4.5) is $O(n + n^2 + (n \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^n))$. Now n is bounded by the maximum branching within a single pair-process, so $|P_k^\ell| \geq n$. Hence the above is $O(n \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^n)$.

The body of step 4 is repeated for every k -state t_k in S_k^r . Let b denote the maximum value of n ($= |t_k.moves|$) as t_k ranges over S_k^r . Thus the complexity of step 4 is $O(b \cdot |P_k^\ell| \cdot |R_{k\ell}^r| \cdot \alpha^b \cdot |S_k^r|)$. The complexities of steps 1, 2, and 3 computed above are $O(|P_k^\ell| \cdot |R_{k\ell}^r|)$, $O(|P_k^\ell|)$, and $O(|R_{k\ell}^r|)$, respectively. These are all subsumed by the complexity of step 4, which therefore gives the overall complexity of the procedure for checking WG . Now $|S_k^r|$ is $O(|P_k^\ell|)$, since S_k^r is a subset of the k -states of P_k^ℓ . Thus, the overall time complexity can be rewritten as $O(b \cdot |P_k^\ell|^2 \cdot |R_{k\ell}^r| \cdot \alpha^b)$. Replacing α by $|R_{k\ell}^r|/|P_k^\ell|$, we obtain $O(b \cdot |R_{k\ell}^r|^{b+1}/|P_k^\ell|^{b-2})$ for the worst-case time complexity of the procedure for checking the wait-for-graph assumption.

Now each $|R_{k\ell}^r|$ is $O(|P_k^\ell|^2 \cdot 2^{|\mathcal{SH}_{k\ell}|})$, since a pair-system contains two pair-processes (which contribute $|P_k^\ell|^2$ to the size of $R_{k\ell}^r$), and a set of pairwise shared variables (which contributes $2^{|\mathcal{SH}_{k\ell}|}$ to the size of $R_{k\ell}^r$). Hence, the overall time complexity can also be written as $O(b \cdot (|P_k^\ell|^{2b+2} \cdot 2^{(b+1)|\mathcal{SH}_{k\ell}|})/|P_k^\ell|^{b-2})$, that is, $O(b \cdot |P_k^\ell|^{b+4} \cdot 2^{(b+1)|\mathcal{SH}_{k\ell}|})$.

Also, the space complexity of the procedure is $O(|R_J^r|)$, since R_J^r is the largest of a fixed set of data structures that are used. This is $O(|R_{k\ell}^r|^{b+1}/|P_k^\ell|^b)$, or, alternatively, $O(|P_k^\ell|^{b+2} \cdot 2^{(b+1)|\mathcal{SH}_{k\ell}|})$.

E.2 Checking the Liveness Assumption

The liveness assumption (6.7.2.1) is mechanically checked as follows. We introduce a “new”¹⁵ atomic proposition Q to M_J . We set Q to *true* in every state s_J of M_J such that

$$M_{jk}, s_J \uparrow jk \models EGex_k$$

and to *false* in all other states of M_J .

Now $\bigwedge a_i^J \cdot (a_i^J \longrightarrow P_j^J \not\subseteq W_J(s_J))$ is equivalent to $s_J \models noblock(i, j)$, where $noblock(i, j) \stackrel{\text{df}}{=} \bigwedge a_i^J \cdot (\{a_i^J.start\} \Rightarrow a_i^J.guard_j)$. Thus, the liveness assumption can be rewritten as

for every reachable state s_J in M_J ,

$$M_{jk}, s_J \uparrow jk \models EGex_k \text{ implies } M_J, s_J \models noblock(i, j).$$

By definition of Q , we have $M_{jk}, s_J \uparrow jk \models EGex_k$ iff $M_J, s_J \models Q$. Hence the above is equivalent to

for every reachable state s_J in M_J ,

$$M_J, s_J \models Q \text{ implies } M_J, s_J \models noblock(i, j).$$

By CTL semantics, this is equivalent to

for every reachable state s_J in M_J ,

$$M_J, s_J \models (Q \Rightarrow noblock(i, j)).$$

Finally, we translate the quantification over all reachable states into CTL by prefixing the formula with the AG modality, and evaluating it in all initial states:

¹⁵i.e., $Q \notin \mathcal{AP}_i \cup \mathcal{AP}_j \cup \mathcal{AP}_k$.

$$M_J, S_J^0 \models AG(Q \Rightarrow noblock(i, j)).$$

To determine the truth assignment to Q , we model-check M_{jk}^r (the reachable part of M_{jk}) for the formula $AGEGex_k$. We employ the CTL model-checking algorithm of Clarke et al. [1986], which marks all states in M_{jk} with all subformulae of $AGEGex_k$ that are true in the state. In particular, all states of M_{jk} that satisfy $EGex_k$ will be so marked. We can then use this marking to assign the appropriate truth value to Q in each reachable state s_J of M_J . Finally, we model-check M_J with respect to $AG(Q \Rightarrow noblock(i, j))$ to determine whether the liveness assumption holds or not, again using the model-checking algorithm of Clarke et al. [1986].

The algorithm of Clarke et al. [1986] has time complexity linear in both the structure and the formula being checked. Here we invoke it twice, once for M_{jk}^r with respect to $AGEGex_k$ and once for M_J^r with respect to $AG(Q \Rightarrow noblock(i, j))$. Since $|AGEGex_k|$ is constant, and $|M_{jk}^r| < |M_J^r|$ is easily seen to hold, the time complexity of model-checking M_J^r dominates. Since $noblock(i, j)$ is quantified over all the moves of P_i^J , its size is $O(|P_i^J|)$. Hence, the liveness assumption can be model-checked in time $O(|M_J^r| \cdot |P_i^J|)$. Now $|M_J^r|$ is $O(|P_k^\ell|^3 \cdot 2^{2|\mathcal{SH}_{k\ell}|})$, since the J -system contains three processes and two sets of pairwise shared variables. Also $|P_i^J|$ is $O(|P_\ell^k|)$, since $|J|$ is fixed. Hence the overall complexity can be rewritten as $O(|P_k^\ell|^4 \cdot 2^{2|\mathcal{SH}_{k\ell}|})$. Note that the check must be made for the case $i \neq k$ and also for the case $i = k$.