ON THE MECHANICAL TRANSFORMATION
OF PROGRAM EXECUTIONS TO
DERIVE CONCURRENCY

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TR-83-20 October 1983
Abstract

A recent trend in program development is to derive correct implementations from program specifications by the application of a formal calculus, a "programming methodology". The application of formal rules lends itself to automation.

We investigate the automation of one part of a methodology for programming with concurrency. In this methodology, concurrency is derived by transforming the sequential execution of a program into an equivalent concurrent execution on the basis of formal transformation rules. This paper is an initial investigation of the automation of such transformations for a certain class of programs: sorting networks. We present an implementation of a part of the underlying semantic theory in Boyer & Moore's mechanical logic, and report on the mechanical proof of a transformation that derives concurrency for a bubble sort.
Table of Contents

1 Introduction
2 A Methodology for Programming with Concurrency
3 On the Automation of Trace Transformations
4 Proving the Semantic Equivalence of Two Traces: A Case Study
   4.1 Boyer & Moore's Theorem Prover
   4.2 Trace Representation
   4.3 Trace Semantics
   4.4 Trace Composition
   4.5 Trace Transformations
   4.6 Non-Neighbors
   4.7 Semantic Equivalence of TAU and TAU-
   4.8 Evaluation
5 Conclusions
6 References

Appendix A: Events of Proof Session
Appendix B: Proof Outline of TAU's Transformation
1 Introduction

In the last decade, few "buzzwords" have gained as much popularity in computing as structured programming [7]. It refers to a disciplined approach to program development by a set of informal derivation rules (stepwise refinement, divide-and-conquer, case analysis, data abstraction, etc.). The application of such rules is encouraged by suitable programming constructs in "structured" programming languages. Structured programming (a more appropriate term would be "disciplined" programming) aims to provide guidance and diminish the need for intuition during the process of programming. The need for intuition makes programming enjoyable and rewarding, but it also makes it difficult. Foremost, programming should be reliable, not enjoyable. Software production aims to reward the customer first, the programmer second.

The reliability of disciplined programming methods is greatly enhanced when they can be applied on a formal basis. Program verification [9, 13] refers to formal methods of proving properties of programs, most importantly, that the program conforms to its specifications. Program verification provides an exact description of programs and thereby their correctness or incorrectness. But it does not support correct program development.

Programming methodology [8, 11, 12] is a new area of research that attempts to merge structured programming and program verification. A programming methodology is a set of formal program development rules, a "programming calculus", that can be applied to transform the specification of a program into a correct
implementation. It replaces some of the programmer's intuition by formal guidance.

The programmer is left with the same responsibility but with less reward. Where the pleasure of invention and creativity has been replaced by the boredom of applying the intricate rules of a formal logic, there remains the responsibility to apply these rules correctly. We can ease the programmer's plight by automating the most monotonous and technical parts of this task. We are therefore interested in a connection between programming methodology and automatic theorem proving [5], the mechanical application of formal logics.

Concurrency is the execution of several program parts in parallel. Rigor in program development is especially desirable in the presence of concurrency. If the input data are known, the outcome of a program that performs operations in some defined sequence is totally predictable. But the outcome of a program that performs operations concurrently depends on the unpredictable order of its concurrent activities. Such dependencies may be buried deep inside the program's operations, or may even be determined by the architecture of the machine on which the program is executed. Simple inspection of the program text cannot reveal them. In addition, the number of possible orderings explodes exponentially with the number and length of concurrent activities. To describe the properties of a concurrent execution, a formal treatment is absolutely essential. The verification of programs with concurrency is now quite well developed [18, 23, 25]. However, programming methodologies are only just emerging. This paper deals with one such methodology [19].
2 A Methodology for Programming with Concurrency

Our goal is to mechanize parts of a particular methodology for the derivation of concurrency in programs [19]. This section describes that methodology.

Two different motives may lead to the application of concurrency:

(1) The desire for a specific program behaviour.

For example, one might wish to run an experiment which involves certain processes executed by designated processors that communicate and synchronize with each other in some fashion. Such applications are to ensure the correct functioning of some machine configuration with a specific concurrency structure.

(2) The desire for fast program results.

For example, one might wish to execute a numerical or data processing algorithm with concurrency in order to obtain a result faster. Such applications do not refer to a specific machine configuration or concurrency structure.

The programming methodology described here takes the second approach: concurrency is viewed as a tool for accelerating the acquisition of results, not as a basic characteristic of a program. Consequently, concurrency will not be part of the problem specification, but will be determined after the derivation of the program. We would like to perform this determination mechanically.
The specification of a programming problem has three parts:

(a) the input constraints under which the program shall operate,
(b) the results which the program is supposed to achieve, and
(c) the time limit imposed on the program's execution.

The program development then proceeds along the following lines:

(1) Perform a formal stepwise refinement of a program that achieves
the desired result under the given input constraints. The
program does not address the question of sequencing. It may not
require a total order of the execution of its operations. But an
easy, sequential execution can, at this point, serve as a first
execution time estimate.

(2) Declare simple relations between program components, so-called
"semantic relations", that allow relaxations in sequencing, e.g.,
concurrency. Do so until the execution time of the program
satisfies the specified time limit.

A refinement is, for instance, C1;C2. This refinement says that
program component C2 is applied to the results of program component
C1. The semicolon denotes "application". It may be implemented by
executing C1 and then C2, but need not be in all cases.

Semantic relations are, for instance, the commutativity of the
components C1 and C2 (written C1&C2), and the independence of C1 and
C2 (written C1||C2). C1 and C2 are commutative, i.e., C1&C2 may be
declared if the execution of C1 and then C2 has the same effect as the
execution of C2 and then C1. If C1&C2 is declared, C1;C2 may also be
implemented by executing C2 and then C1. C1 and C2 are independent,
i.e., C1 || C2 may be declared if the execution of C1 and C2 in parallel has the same effect as their execution in order. If C1 || C2 is declared, C1; C2 may also be implemented by executing C1 and C2 in parallel. Independence implies commutativity.

Let us illustrate the use of this methodology with an example.

Example: Sorting

The problem is to sort an array a[0..n] of real numbers into ascending order in time O(n). Our refinement is an insertion sort adapted from Knuth [15]:

\[
\begin{align*}
\text{sort}(n): & \quad \text{for } i = 1 \text{ to } n \text{ do } S(i) \\
S(0): & \quad \text{skip} \\
(i > 0) & \quad S(i): \quad cs(i); S(i-1) \\
|i-j| > 1 & \Rightarrow cs(i) || cs(j)
\end{align*}
\]

Statement cs(i) represents a restricted version of Knuth's "comparator module" [15]: cs(i) compares adjacent array elements a[i-1] and a[i] and, if necessary, swaps them into order. Knuth's comparator module can deal with not only adjacent but any two array elements. Knuth calls programs whose primitive components are comparator modules sorting networks. Let us call our restricted version simple sorting networks.

Note that, for |i-j| > 1, i.e., if i and j are not "neighbors", cs(i) and cs(j) are disjoint: they do not share any variables. Components that do not share variables may be declared independent. (For the underlying theory see [20].)
For, say, a five-element array (n=4), the refinement has the following sequential execution, if we interpret composition (denoted by ';'') as execution in order (denoted '->') and expand components S(i) of sort(4):

\[
\text{tau}(4) = \text{cs } 1 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1 \rightarrow \text{cs } 3 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1 \rightarrow \text{cs } 4 \rightarrow \text{cs } 3 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1
\]

If we count the number of comparator modules cs, \text{tau}(4) has length 10. In general, \text{tau}(n) has length \(n(n+1)/2\), i.e., is quadratic in \(n\). To derive a linear execution, we have to exploit the independence declaration for sort(\(n\)) and compress \text{tau}(n) into a trace with concurrency. Let us lay out the sequential trace \text{tau}(4) in a form which suggest how this can be done:

\[
\text{tau}(4) = \text{cs } 1 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1
\]

\[
\rightarrow \text{cs } 3 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1
\]

\[
\rightarrow \text{cs } 4 \rightarrow \text{cs } 3 \rightarrow \text{cs } 2 \rightarrow \text{cs } 1
\]

We commute comparator modules in \text{tau}(4) left, and then merge adjacent modules whose indices differ by 2 into a parallel:

\[
\text{tau}^\sim(4) = \text{cs } 1 \rightarrow \text{cs } 2 \rightarrow < \rightarrow < \rightarrow < \rightarrow \text{cs } 2 \rightarrow \text{cs } 1
\]

\[
\text{cs } 3/ \rightarrow \text{cs } 4/ \rightarrow \text{cs } 3/
\]

This execution is of length 7. In general, \text{tau}^\sim(n) is of length 2n-1, i.e., linear in \(n\). The degree of concurrency increases as we add inputs. We are not limited to a fixed number of concurrent actions. However, if only a fixed number \(k\) of processors is available, the independence declaration may not be exploited to generate a degree of concurrency higher than \(k\).
To declare semantic relations for some refinement, one does not need to understand the refinement as a whole. A local understanding of the components appearing in the declared relation is sufficient. Note also the simplicity of concurrency: there is no need for synchronization of concurrent components other than at the point of termination. Most semantic declarations come easily to mind and have a simple proof.

But the foremost benefit of this approach to the derivation of fast programs is that the more important and better understood question of program refinement is resolved before the less important and more complex question of concurrency arises. Concurrency is later added in isolated steps (by invoking semantic relations) without changing the approved meaning of the program. Thus the development of programs with concurrency is divided into two stages:

Stage 1: The development and formal semantic description of a program that achieves the desired result. This requires a formal refinement and the declaration of semantic relations.

Stage 2: The derivation of a fast execution of the program produced at Stage 1. (An execution of a program is also called a trace.) This is conceptually simple but computationally complex. It involves the computation of execution times and the invocation of semantic relations to transform traces and improve execution time.

Either of the two stages has the potential for automation. Automation of Stage 1 would yield a mechanical system for program refinement. Research along these lines is under way elsewhere [2,
22]. Automation of Stage 2 would yield a very powerful optimizing compiler (since we view concurrency as optimization). Early work in this area [17] has been without a formal semantic basis, because at that time formal semantics was in its infancy. Our objective is the automation of Stage 2 on a formal semantic basis.

The most common approach to programming in which the specification of concurrency is shifted from program development to program translation or execution is data flow programming [1]. A data flow program makes no explicit reference to the order of execution. It is executed on a special machine architecture that follows the sequencing imposed by the data dependencies of the program's variables. Data flow languages are "referentially transparent": they do not permit the re-assignment of variables. This simplifies the identification of data dependencies so much that, commonly, no programmer assistance is needed to identify concurrency. Our approach is "referentially opaque", i.e., permits the re-assignment of variables and, consequently, requires a more complicated data flow analysis. We have to explicitly declare and subsequently exploit data dependencies (in our formalism, semantic declarations).

The vast majority of software that exists today and is currently being produced is referentially opaque. The vast majority of today's machine architectures support the referentially opaque programming style. While we must strive for new programming styles and machine architectures, we must also continue to increase our understanding of the present technology.
We shall focus on the automation of Stage 2. (For initial ideas see [21, Sect. 6.2.1].)

Let us consider the previous example sort(n). Stage 2 derives a parallel trace \( \tau^*(n) \) that satisfies the execution time limit from sequential trace \( \tau(n) \) that is suggested by the refinement of sort(n) and does not satisfy the execution time limit. The semantic equivalence of \( \tau(n) \) and \( \tau^*(n) \) ensures that the meaning of the program is preserved. It is proved by a recursive application of a sequence of trace transformations that exploit no semantic relations other than those declared for sort(n). Although such transformations are in many cases, as in this one, quite simply described in informal English (see Sect. 2), their formal application is extremely tedious (see [21, Sect. 5.4]). We do not want to rely on the informal description but would like some mechanical aid in the formal application.

We might be tempted to view a transformation as an algorithm. For example, we might propose an algorithm that transforms \( \tau(n) \) into \( \tau^*(n) \). We would need a set of transformation commands [21, Sect. 6.2.1]. The algorithm we propose uses commands:

- `commute x left` to identify the next trace element of form \( x \) to the right of the current cursor position and commute it with its left neighbor, and
- `merge x left` to identify the next trace element of form \( x \) to the right of the current cursor position and merge it and its left neighbor into a parallel command.
Let us assume that, at the start of any transformation, the cursor is placed at the left end of the trace to be transformed. For the following algorithm, trans(n), the cursor is initially placed to the left of trace tau(n):

\[
\text{trans(n): } \begin{array}{c}
\text{for } i:=3 \text{ to } n \\
\text{do for } j:=0 \text{ to } i-3 \\
\text{do for } k:=1 \text{ to } i-3-j \\
\text{do commute cs(i-j) left} \quad \text{od} \\
\text{merge cs(i-j) left} \\
\text{od} \\
\text{od}
\end{array}
\]

trans(n) derives tau^{-}(n) from tau(n) in one left-to-right pass of right-to-left commutations and merges in cubic time (O(n) for every nested for). For n<3, tau(n)=tau^{-}(n) and trans(n) does nothing.

This approach has the problem that if the refinement contains a recursion (or repetition), as sort(n) does, the transformation algorithm refers to the depth of the recursion: for fixed n, trans(n) makes sort(n) linear, but not sort(n+1). We are trying to "execute" a recursive transformation, but an unbounded recursion can never be executed to completion.

A better approach is to treat trace transformations as what they are: theorems, not algorithms. In particular, recursive transformations are inductive theorems. For our transformation of sort(n), the theorem is:

TAU.MAIN: For all n>0, semantics of tau(n) = semantics of tau^{-}(n)

and is proved by induction on n.
With this view, the automation of Stage 2 constitutes a system that identifies, following a set of heuristics and, most likely, with the help of some interactive dialogue, valid and useful transformation theorems. Our present approach is to split this task into two parts:

(a) The identification and (recursive) formulation of trace transformations, and

(b) the proof of the corresponding (inductive) theorems.

This way, we can make use of existing results on (b) to guide us in the research of (a). Note that we prove single trace transformations correct. Once we have a set of algorithms that identify trace transformations, i.e., perform step (a), we may want to establish the correctness of these algorithms rather than verifying their output individually. But, for the time being, we prefer to deal with the semantics of trace transformations, not with the semantics of algorithms that yield trace transformations.

In the mechanical development and proof of theorems, simplicity is of overriding concern—much more so than in the proof of theorems alone. More often than not, mechanical proofs are messy and full of hints to the theorem prover, e.g., hints on the use of specific lemmas, induction schemes, or variable instantiations. These hints are supplied by the user in an interactive dialogue with the theorem prover. But a mechanical development of theorems must include the generation of the necessary hints for its proof. Therefore we must take special care to minimize proof hints and keep those hints that are indispensable systematic. In our case, the key is a simple representation of the semantic theory and a straightforward development of lemmas relevant to the equivalence of traces. The
following section develops the theory of simple sorting networks in more detail and, at the end, reflects on our efforts to keep things clean.

4 Proving the Semantic Equivalence of Two Traces: A Case Study

We are applying Boyer & Moore's mechanical treatment of recursion and induction [3]. This section describes the implementation of the part of the semantic theory that is necessary to prove theorem TAU.MAIN in Boyer & Moore's logic, and the proof of TAU.MAIN. For newcomers to the Boyer & Moore prover, Sect. 4.1 provides a short introduction adapted from [26]. App. A explains the types of commands (so-called "events") that formed the proof session and contains a transcript of all functions, axioms, and lemmas established during the session. App. B elaborates on the central steps of the mechanical proof.

Boyer & Moore express terms functionally in the prefix notation of Church's lambda calculus [6] and LISP [24]. To be shorter, we shall, in this document, keep basic arithmetic and logical operations in traditional infix notation.

4.1 Boyer & Moore's Theorem Prover

In Boyer & Moore's systems as described in [3], proofs are constructed in a quantifier-free logical theory which is built on the propositional calculus with equality, variables, and function symbols. The basic theory includes four functions:
TRUE and FALSE are both functions of zero arguments. The constants (TRUE) and (FALSE) are abbreviated T and F, respectively. The axiom T\#F allows them to be considered as distinct truth values.

EQUAL is a function of two arguments, axiomatized to require that (EQUAL l r) have the value T or F, depending on whether l=r.

IF is a function of three arguments. Axioms insure that the value (IF t u v) is that of v if t=F and that of u otherwise.

The function IF allows the use of conditionals in axioms which define new functions, as in the definition

\[(AND \: P \: Q) = (IF \: P \: (IF \: Q \: T \: F) \: F)\]

A function capturing the semantics of each of the other logical connectives is similarly defined. Thus, corresponding to any formula w, a term p can be constructed such that the formulas p\#F and w have the same truth values. In this situation, the term p is said to be a "theorem" if w is a theorem. All formulas may thus be represented as terms.

The theory also includes

a principle which permits the introduction of axioms specifying new types of inductively constructed objects,

a principle for admitting axioms which define new recursive functions,

a principle of induction (a rule of inference) based on the notion of a well-founded relation which is well-suited for inferring theorems about these objects and functions.

Two types of inductively constructed objects are relevant to the proofs presented below:

The natural numbers are formalized by the type "number". Peano arithmetic is realized through axioms about the functions NUMBERP, ADD1, SUB1, and the constant 0. NUMBERP is the "recognizer" for this type, i.e., (NUMBERP t) returns T or F.
depending on whether the value of term \( t \) is a number. \( \text{ADD1} \) is the usual successor function, and \( \text{SUB1} \) is its inverse function.

Ordered pairs are formalized by the type "list". The functions \( \text{CONS} \), \( \text{CAR} \), and \( \text{CDR} \) are axiomatized to have the properties of the familiar LISP functions, and \( \text{LISTP} \) is the recognizer for this type. Finite sequences are represented by means of \( \text{CONS} \) and the special constant \( \text{NIL} \); the sequence of terms \( t_1, t_2, \ldots, t_n \) is represented by the term

\[
(\text{CONS} \ t_1 \ (\text{CONS} \ t_2 \ (\ldots \ (\text{CONS} \ t_n \ \text{NIL}) \ldots)))
\]

which is abbreviated \( (\text{LIST} \ t_1 \ t_2 \ \ldots \ t_n) \).

The theorem proving program naturally relies heavily on induction. Several heuristics are employed to formulate induction schemes based on analysis of the structure of the recursive function definitions and the inductively constructed types which are involved in a conjecture. A proof by induction is only attempted, however, after a simplification procedure has been followed and has failed to establish the theorem.

Every conjecture which is presented to or generated by the theorem prover is first written as a single term and then represented internally as a conjunction of simpler "clauses", each of which is a disjunction of atomic formulas 'called "literals". In order to establish a clause, the prover first attempts to simplify each of its literals in turn, assuming the others to have the value \( F \). A variety of heuristics is employed, including the use of previously proved lemmas as rewrite rules. A term \( l \) is replaced by the term \( r \) if a rewrite rule of the form

\[
(\text{IMPLIES} \ h \ (\text{EQUAL} \ l \ r))
\]

is encountered and the hypothesis \( h \) can be established. The manner of use of a lemma therefore depends on its precise syntactic form.
In this paper (except App. A), we write (EQUAL l r) as \(1=r\) and (IMPLIES h c) as \(h\Rightarrow c\). We also write (ADD1 n) as \(n+1\), and (SUB1 n) as \(n-1\).

4.2 Trace Representation

Our goal is to prove the semantic equivalence of traces \(\tau_a(n)\) and \(\tau_a^-(n)\). E.g., for \(n=4\),

\[
\text{cs 1 } \rightarrow \text{ cs 2 } \rightarrow \text{ cs 1 } \rightarrow \text{ cs 3 } \rightarrow \text{ cs 2 } \rightarrow \text{ cs 1 } \rightarrow \text{ cs 4 } \rightarrow \text{ cs 3 } \rightarrow \text{ cs 2 } \rightarrow \text{ cs 1 }
\]

must prove semantically equal to

\[
\tau_a^-(4) = \text{cs 1 } \rightarrow \text{ cs 2 } \rightarrow \text{cs 1}\ \text{/} \text{cs 3/} \rightarrow \text{cs 2/} \rightarrow \text{cs 1}
\]

We represent a trace as a LISP list. The elements of the list are executed in sequence. If a list element is itself a list, it is called a "parallel command" and its elements are executed in parallel. If an element of a parallel command is again a list, its elements are executed in sequence, etc. Thus, a trace is a multi-level list whose odd levels reflect sequential execution, and whose even levels reflect parallel execution. In the realm of sorting networks, we can represent traces as multi-level lists of integers. For example,

\[
\begin{align*}
\tau_a(5) & = '(1 2 1 3 2 1 4 3 2 1 5 4 3 2 1) \\
\tau_a^-(5) & = '(1 2 (3 1) (4 2) (5 3 1) (4 2) (3 1) 2 1)
\end{align*}
\]

Note that the parallel trace \(\tau_a^-(5)\) has a phase \(T_1^-\) of growing concurrency

\[
(T_1^- 5) = '(1 2 (3 1) (4 2))
\]
followed by a phase $T_2^-$ of shrinking concurrency

$$(T_2^- 5) = '((5 3 1) (4 2) (3 1) 2 1)$$

In our formalism [20], parallel commands are binary, i.e., can have at most two parallel components. An $n$-ary parallel command is expressed as nested binary parallel commands. This coincides with LISP's (and Boyer & Moore's) representation of a list of $n$ elements as a nesting of pairs. For example, the parallel command $(5 3 1)$ of $(T_A^- 5)$ is really $(5 . (3 . (1 . NIL)))$.

4.3 Trace Semantics

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Traces have weakest precondition semantics [20]. Since a weakest precondition is a function from programs and predicates to predicates [8], the weakest precondition calculus can be directly implemented in Boyer & Moore's logic.

Our methodology divides the development of programs into two stages. Stage 1 is concerned with the derivation of program semantics, i.e., the derivation of a refinement. Stage 2 is concerned with the preservation of program semantics, i.e., the transformation of sequential executions into concurrent executions. Consequently, we need not implement a complete weakest precondition generator in order to implement Stage 2. A weakest precondition that is not affected by the trace transformations is not relevant to Stage 2; it need not be spelt out but may be provided as a "black box". In Boyer & Moore's logic, a black box is represented by a function that has been declared (without a function body) rather than defined (with a function body). The primitive components of sorting networks are comparator modules.
For the purpose of trace transformations, we are not interested in the inside of a comparator module. Therefore we declare the weakest precondition of comparator module cs(i) as a function

\[(CS \ I \ S)\]

where S denotes the postcondition (or "poststate"). Since function CS is declared, not defined, we must provide by axiom some essential information about CS that is not evident from the declaration. We add two axioms. One restricts the domain of comparator modules to the positive integers:

\[\text{CS.TAKES.NUMBERS:} \quad (\text{NOT} \ (\text{NUMBERP} \ I)) \Rightarrow ((CS \ I \ S) = \text{F})\]

The name of the axiom is CS.TAKES.NUMBERS. Axiom CS.TAKES.NUMBERS states that the prestate of CS for any non-number and poststate is false, i.e., that such a CS is not permitted. The other axiom expresses the "rule of the excluded miracle" (Dijkstra's first healthiness criterion [8]) for comparator modules:

\[\text{CS.TERMINATES:} \quad (CS \ I \ F) = \text{F}\]

Axiom CS.TERMINATES states that the prestate of any CS with false poststate is false. In the weakest precondition calculus, this means comparator modules must terminate.

To determine the weakest precondition of some trace L that is composed of comparator modules CS for poststate S, we define a "cs-machine", a function

\[(M.CS \ FLAG \ L \ S)\]
that composes calls to CS as prescribed by trace L. Beside L and S, M.CS takes a FLAG that signals whether the trace is to be executed in sequence (FLAG='SEQ) or in parallel (FLAG='PAR). In accordance with our trace representation, FLAG='SEQ in top-level calls and FLAG alternates with every recursive call.

When FLAG='PAR, the trace represents a parallel command and its elements must be checked for independence. We can make use of the semantic declarations provided at Stage 1. The smallest component that a semantic declaration for a sorting network will mention is the comparator module. We may therefore, from Stage 1, assume knowledge about the independence of comparator modules and may express this knowledge by a declared function

$$(\text{IND.CS I J})$$

that evaluates the independence of comparator modules cs(i) and cs(j). We then define a function

$$\text{(ARE.IND.CS L1 L2)}$$

that uses IND.CS to determine the mutual independence of all atoms of trace L1 with all atoms of trace L2. ARE.IND.CS is only interested in the atoms of a trace, not in the trace's structure. Therefore, ARE.IND.CS expects traces in a flattened form. The flattening is performed by function ALL.ATOMS, for example,

$$(\text{ALL.ATOMS (TAU- 5)})$$

$$= '(1 2 3 1 4 2 5 3 1 4 2 3 1 2 1)$$

If the two members of a parallel command (remember the restriction to binary parallel commands) pass test ARE.IND.CS their execution has
identical semantics in parallel as in sequence — only their execution time differs.

The execution time of traces plays a role in the selection of proper transformation theorems. At present, we take transformation theorems as given and only prove them by mechanical means. Therefore, execution time is left out of the current implementation.

The semantic equivalence of TAU⁻ and TAU is formally expressed by the equality

\[(\text{M.CS 'SEQ (TAU⁻ N)} \not\in \text{S}) = (\text{M.CS 'SEQ (TAU N)} \not\in \text{S})\]

### 4.4 Trace Composition

For the composition of lists, we provide two functions: APPEND and APPEND2. APPEND is the regular LISP (and Boyer & Moore) append function. It handles lists that end with NIL. However, we do not require a trace to end with NIL; it may end with any atom. If trace L1 does not end with NIL, (APPEND L1 L2) drops the last atom, whereas (APPEND2 L1 L2) picks it up. For example,

\[
\text{(APPEND '(1 . 2) '(3 . 4)) = '(1 3 . 4)}
\]

\[
\text{(APPEND2 '(1 . 2) '(3 . 4)) = '(1 2 3 . 4)}
\]

Therefore, traces are composed with APPEND2, not with APPEND.

The following theorem expresses the composition rule for weakest preconditions [8]:
M.CS.APPEND2.SEQ:

(FLAG='SEQ) => ( (M.CS FLAG (APPEND2 L1 L2) S) = (M.CS FLAG L1 (M.CS FLAG L2 S)) )

We made M.CS.APPEND2.SEQ part of a larger theorem (M.CS.APPEND2).

For lists that end with NIL, APPEND2 equals APPEND. We use APPEND for flattened traces, i.e., traces that have been treated with ALL.ATOMS.

4.5 Trace Transformations

Independence declarations are exploited via transformation rules that express commutations and parallel merges of independent program components.

The theorem for parallel merges corresponds to transformation rule (G31) of [20, Sect. 5.2]:

G31: (ARE.IND.CS (ALL.ATOMS L1) (ALL.ATOMS L2)) => ( (M.CS 'PAR (CONS L1 L2) S) = (M.CS 'SEQ (APPEND2 L1 (LIST L2)) S) )

To express commutations, we must be more specific about the meaning of "independence". The declaration of IND.CS does not provide any clues. We do not need to know everything about independence; otherwise we had defined, not declared IND.CS. But we must be able to conclude that independent comparator modules may be commuted. As we did with CS, we characterize IND.CS by axiom:

GLOBAL.IND.CS:

(IND.CS I J) => (((CS J (CS I S)) = (CS I (CS J S))))
If we instantiate both FLAG1 and FLAG2 to 'SEQ, the following theorem enables commutations:

\[
\begin{align*}
\text{ARE.IND.CS.IMPLIES.COMMUTATIVITY:} \\
(\text{ARE.IND.CS (ALL.ATOMS L1) (ALL.ATOMS L2)}) \\
\Rightarrow \left( (\text{M.CS FLAG1 L1 (M.CS FLAG2 L2 S)}) \\
= (\text{M.CS FLAG2 L2 (M.CS FLAG1 L1 S)}) \right)
\end{align*}
\]

4.6 Non-Neighbors

Stage 1 establishes that, in simple sorting networks, non-neighboring comparator modules are independent. We may provide this known fact by axiom:

\[
\begin{align*}
\text{NON.NEIGHBORS.ARE.IND.CS:} \\
(\text{NON.NEIGHBORS I J}) \Rightarrow (\text{IND.CS I J})
\end{align*}
\]

where function NON.NEIGHBORS identifies non-neighbors. NON.NEIGHBORS is defined while IND.CS is declared. With IND.CS alone we could not decide the independence of anything; with this axiom we can decide the independence of comparator modules. We may, for example, apply theorem G3i with 5 for L1 and '(3 1) for L2, since 5 is not neighbor of 3 and 1:

\[
(\text{M.CS 'PAR '(5 3 1) S}) = (\text{M.CS 'SEQ '(5 3 1)) S})
\]

Two more applications of G3i, exploiting also the non-neighborhood of 3 and 1, yield:

\[
(\text{M.CS 'PAR '(5 3 1) S}) = (\text{M.CS 'SEQ '(5 3 1) S})
\]
This formula expresses the equivalence of the parallel and sequential execution of comparator modules CS(5), CS(3), and CS(1).

4.7 Semantic Equivalence of TAU and TAU^-

The formal derivation of TAU^- from TAU requires three inductions on n. The main induction transforms TAU into TAU^- as described in Sect. 2.

TAU.MAIN: 0<N => ( (M.CS 'SEQ (TAU^- N) S)  
= (M.CS 'SEQ (TAU N) S) )

Notice the premise that trace argument N "makes sense", i.e., is a positive number. (A further theorem, TAU.MAIN.COMplete, extends the result to meaningless N.) The prover must be made aware of the induction scheme by a hint (function RAVEL). Let us explain the induction with an example. For N=5, the following traces are semantically equivalent:

(TAU 5) = '1 2 1 3 2 1 4 3 2 1 5 4 3 2 1)  
= '1 2 (3 1) (4 2) (3 1) 2 1 5 4 3 2 1) (ind. hyp.)  
= '1 2 (3 1) (4 2) (5 3 1) (4 2) (3 1) 2 1) (ind. step)  
= (TAU^- 5)

The induction hypothesis is that (TAU 4) is equivalent to (TAU^- 4). The induction step "ravels" the remaining sequential tail (SEQJ 5) to the left into (TAU^- 4). Note that the only part of TAU^- affected by this transformation is the part of shrinking concurrency, T2^- . Thus we may disregard T1^- in the further proof of the induction step. To perform the transformation of the induction step, we need two more inductions. The first, named TAU.RAVEL1, uses commutations (ARE.IND.CS.IMPLIES.COMMUTATIVITY) to place comparator modules beside
the parallel command that they then join (G3i) by the second induction, named TAU.RAVEL2. The intermediate stage of the trace has name UNRAVEL.T2^-.

TAU.RAVEL1:

1<N ⇒ (M.CS 'SEQ (UNRAVEL.T2^- N) S) = (M.CS 'SEQ (APPEND2 (T2^- N-2) (SEQJ N)) S)

TAU.RAVEL2:

0<N ⇒ (M.CS 'SEQ (T2^- N) S) = (M.CS 'SEQ (UNRAVEL.T2^- N) S)

For example, for N=5, the following traces are semantically equivalent:

(APPEND2 (T2^- 3) (SEQJ 5))
= '((3 1) 2 1 5 4 3 2 1) (by TAU.RAVEL1)
= '((5 3 1) (4 2) 3) (1) 2 1) (by TAU.RAVEL2)
= '((5 3 1) (4 2) (3 1) 2 1)

This concludes the main part of the proof. However, the prover needs some preliminary information to succeed. To enable the application of our general transformation rules G3i and ARE.IND.CS.IMPLIES.COMMUTATIVITY, we have to establish their prerequisites: the independence of certain trace parts. We also have to advise the prover of some trace identities that it cannot determine simply by opening functions. The interested reader is referred to App. A and B.
4.8 Evaluation

The main objective of the mechanical proof of theorem TAU.MAIN has been to substantiate that the trace calculus of [20] has a natural representation in Boyer & Moore's logic. The functions, axioms, and theorems of this representation fall into two classes:

(a) the basic semantic theory that will be required in every proof of a trace transformation theorem, and
(b) the functions and lemmas pertaining to the proof of a specific transformation theorem.

To prove TAU.MAIN, we had to develop a part of the theory of sorting networks (the theory of simple sorting networks, where comparator modules deal only with adjacent elements). It includes the semantics of traces of comparator modules, the general trace transformation rules, and the theory of non-neighbors. While the basic semantic theory is constant, the theory specific to the transformation of TAU, based on the definitions of TAU and TAU⁻ and culminating in theorem TAU.MAIN, will recur in modified form in other trace transformations.

In the original calculus, a transformation sequence starts with a sequential trace and ends with a parallel trace. In Boyer & Moore's logic, we rewrite in the reverse direction, from parallel to sequential, in order to avoid non-termination problems. We need not be aware that the prover actually transforms TAU⁻ into TAU. The direction of transformation is of no consequence to the proof of a semantic equivalence, even though our methodology prescribes a direction of transformation for the derivation of a semantic equivalence.
Because we want to mechanically derive and not only prove theorems, we are especially interested in a simple representation. While the basic semantic theory of sorting networks is constant, the theory specific to TAU will recur in modified form in other trace transformations. Let us explain our efforts to proceed in the most straight-forward manner to gain insight into the potential for a mechanical generation of trace transformations.

We have reduced the number of intermediate stages of TAU's transformation to a minimum: just one intermediate stage between a step of recursive commutations and a step of recursive parallel merges. Boyer & Moore's prover cannot be expected to perform a particular succession of inductions without advice. We had to communicate three traces to the prover: the initial trace TAU, the final trace TAU', and the intermediate trace UNRAVEL.T2'. The prover must also be made aware of the substitutions in the main induction scheme (by hint RAVEL).

Further we had to provide the prerequisites for the transformation, and some trace identities used in its proof. Say, we want to apply some theorem with certain instantiations of its variables. In Boyer & Moore's proof system, we can

(a) prescribe the use of the theorem with appropriate variable instantiations as a hint for the proof in which it is needed, or

(b) provide the hypothesis of the theorem with appropriate variable instantiations as a separate lemma.

Because it is easier to establish facts independently than to tie them to specific proofs, we prefer (b) and establish as lemmas the hypotheses of the general transformation theorems being used, where
free trace variables are instantiated by the particular trace parts that are subject to the transformation. We have to establish precisely these hypotheses, or the prover will not invoke the transformation theorems. For the same reason, the trace identities are expressed as semantic identities (with M.CS), although they hold even as literal identities (without M.CS). The prover would not be able to use the literal identities, because the theory deals with trace semantics, not with traces per se.

This problem is known as the "Knuth-Bendix" problem [16]. A fact appears in Boyer & Moore's logic as a composition of functions. The same fact may be represented in many different ways depending on which function calls are opened up. Some of those representations will match hypotheses of theorems known to the prover, others will not. We have to choose a representation that is useful to the prover.

The Knuth-Bendix problem appears again in the proofs of the actual transformation theorems for TAU. The prover must be prevented from opening functions beyond the level at which the prerequisites are expressed, or it will not establish a match. In addition, since at points where a parallel merge may be applied a commutation applies as well, we have to disable the application of the first transformation (TAU.RAVEL1) in the proof of the second (TAU.RAVEL2).

The worst consequence of the Knuth-Bendix problem in the proof of TAU.MAIN is our only "dirty" lemma: CLOSE.SEQJ. This lemma is not relevant to our theory, but it is needed to make the prover at a specific point in the proof of TAU.RAVEL1 "close" an expansion of function SEQJ (see App. B). To deal with problems like the necessity
of rewrite rule CLOSE SEQA will be one of the challenges of the
mechanical derivation of theorems.

We required two declarations and four axioms to express the
knowledge inherited from Stage 1. Of course, they are all part of the
basic theory. Relative to that theory, the transformation of TAU is
completely defined and certified.

5 Conclusions
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We are exploring the usefulness of techniques in automatic
theorem proving for the formal derivation of programs, specifically,
the derivation of concurrency in programs.

Currently we are focussing on a very specific class of programs:
sorting networks. We chose sorting networks for three reasons.
First, they are well-suited for our methodology: they terminate and
only their results, not their behaviors matter. Second, they have a
simple semantic structure with only one basic component: the
comparator module. Third, they are of more than academic interest:
many data processing applications require sorting.

We have developed a weakest precondition generator function WP,
for programs written in the language RL of our original methodology
[19, 20]. The implementation of the semantic theory of RL (with
function WP) would be significantly more complex than that of sorting
networks presented here (with function M.CS). We want to assess the
practicality of our approach on a simple semantic theory first.
Future research shall extend the mechanical theory to general sorting networks with comparator modules that deal with arbitrary rather than adjacent elements. In that theory, traces are composed of pairs of integers, not single integers. We plan to study a collection of trace transformations of sorting networks, and to derive from this study heuristics for the mechanical generation of such trace transformations.

Proving that a given trace cannot be derived is equally interesting as proving that it can be derived. For example, TAU does not represent the fastest execution that sorts in place. There are sorts with comparator networks that perform in linear time with a smaller constant and even in sublinear time [15]. However, these executions cannot be derived from TAU by the present independence declaration. The proof would have to resort to a meta theory about semantic declarations and the trace transformations they permit.

Acknowledgements
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I am indebted to J Moore and Bob Boyer who responded patiently to my countless questions about their prover. J Moore really bore the burden of introducing me to theorem proving. He also helped me getting started with the implementation of my theory. Thanks J!
References


Appendix A: Events of Proof Session

This appendix presents all events in the order in which they have been accepted by the theorem prover. We use four kinds of events: declaration of a function, definition of a function, addition of an axiom, and proof of a lemma. We shall briefly review the input command format of each. The User's Manual [4] explains how to run proof sessions, in general.

(1) Function Declaration: (DCL name args)

DCL declares "name" to be an undefined function with formal arguments "args".

(2) Function Definition: (DEFN name args body hints)

DEFN defines a function named "name" with formal arguments "args" and with body "body". Before admission of the function, the prover attempts to certify its termination by identifying a well-founded relation such that some measure of "args" gets smaller in every recursive call. In some cases, this relation and measure must be provided in the fifth argument "hints".

(3) Add Axiom: (ADD.AXIOM name types term)

ADD.AXIOM adds a new axiom. The name of the axiom is "name". "types" specifies the ways in which the axiom is used by the prover, and the statement of the axiom is "term". All of our axioms are of type REWRITE, i.e., are used as rewrite rules.

(4) Prove Lemma: (PROVE.LEMMA name types term hints)

PROVE.LEMMA attempts to prove the conjecture "term" and remember it as a lemma named "name". Only successfully proved lemmas are admitted as events. Lemma "name" will be used according to "types". Our lemmas are all used as rewrite rules. The fourth argument "hints" may contain several kinds of directives to aid the proof. We use the following:

(INDUCT (name args)) Use the induction scheme reflected by the recursive definition of function (name args),

(DISABLE ev1 ... evn) Prevent the use of events "evi" to "evn" in the proof.

The following list of events contains: 2 declared functions, 22 defined functions, 4 axioms, and 39 lemmas, i.e., 67 events.
SEMANTIC THEORY OF SIMPLE SORTING NETWORKS

Trace Composition

(DEFN APPEND (X Y)
  (IF (NLISTP X)
      Y
      (CONS (CAR X) (APPEND (CDR X) Y)))))

(DEFN APPEND2 (X Y)
  (IF (NLISTP X)
    (IF (EQUAL X NIL)
      Y
      (CONS X Y))
    (CONS (CAR X)
      (APPEND2 (CDR X) Y)))))

(PROVE.LEMMA APPEND2.NIL (REWRITE)
  (AND (IMPLIED (NOT (EQUAL X NIL))
    (NOT (EQUAL (APPEND2 X Y) NIL)))
  (IMPLIED (NOT (EQUAL Y NIL))
    (NOT (EQUAL (APPEND2 X Y) NIL)))))

(DEFN PLISTP (X)
  (IF (NLISTP X)
    (EQUAL X NIL)
    (PLISTP (CDR X)))))

(PROVE.LEMMA PLISTP.APPEND (REWRITE)
  (IMPLIED (AND (PLISTP X) (PLISTP Y))
    (PLISTP (APPEND X Y))))

(PROVE.LEMMA APPEND2.APPEND (REWRITE)
  (IMPLIED (PLISTP X)
    (EQUAL (APPEND2 X Y) (APPEND X Y))))

(DEFN ALL.ATOMS (L)
  (IF (LISTP L)
    (IF (EQUAL (CDR L) NIL)
      (ALL.ATOMS (CAR L))
      (APPEND (ALL.ATOMS (CAR L))
        (ALL.ATOMS (CDR L))))
    (IF (EQUAL L NIL)
      NIL
      (LIST L)))))

(PROVE.LEMMA ALL.ATOMS.PLISTP (REWRITE GENERALIZE)
  (PLISTP (ALL.ATOMS L)))

(PROVE.LEMMA ALL.ATOMS.APPEND2 (REWRITE)
  (EQUAL (ALL.ATOMS (APPEND2 X Y))
    (APPEND2 (ALL.ATOMS X) (ALL.ATOMS Y))))
Trace Semantics

(DCL CS (I S))

(DCL IND.CS (I J))

(DEFN IS.IND.CS (I L)
  (IF (NILSTP L)
    T
    (AND (IND.CS I (CAR L))
     (IS.IND.CS I (CDR L)))))

(PROVE.LEMMA IS.IND.CS.APPEND (REWRITE)
  (EQUAL (IS.IND.CS I (APPEND L1 L2))
     (AND (IS.IND.CS I L1) (IS.IND.CS I L2))))

(DEFN ARE.IND.CS (L1 L2)
  (IF (NILSTP L1)
    T
    (AND (IS.IND.CS (CAR L1) L2)
     (ARE.IND.CS (CDR L1) L2)))))

(PROVE.LEMMA ARE.IND.CS.APPEND.RIGHT (REWRITE)
  (EQUAL (ARE.IND.CS L1 (APPEND L2 L3))
     (AND (ARE.IND.CS L1 L2) (ARE.IND.CS L1 L3))))

(PROVE.LEMMA ARE.IND.CS.APPEND.LEFT (REWRITE)
  (EQUAL (ARE.IND.CS (APPEND L1 L2) L3)
     (AND (ARE.IND.CS L1 L3) (ARE.IND.CS L2 L3))))

(PROVE.LEMMA ARE.IND.CS.NIL (REWRITE)
  (ARE.IND.CS L NIL))

(DEFN M.CS (FLAG L S)
  (IF (NILSTP L)
    (IF (EQUAL L NIL) S (CS L S))
    (IF (EQUAL FLAG 'PAR)
      (IF (EQUAL (CDR L) NIL)
       (M.CS 'SEQ (CAR L) S)
       (M.CS 'SEQ (CAR L)
        (M.CS 'PAR (CDR L) S))
       (M.CS 'PAR (CAR L)
        (IF (EQUAL (CDR L) NIL)
         S
         (M.CS 'SEQ (CDR L) S))))))

(ADD.AXIOM CS.TAKES.NUMBERS (REWRITE)
  (IMPLIES (NOT (NUMBERP I))
    (EQUAL (CS I S) F))))
(ADD.AXIOM CS.TERMINATES (REWRITE)
  (EQUAL (CS I F) F))

(PROVE.LEMMA M.CS.TERMINATES (REWRITE)
  (EQUAL (M.CS FLAG L F) F)
  ((INDUCT (M.CS FLAG L S)))))

(PROVE.LEMMA M.CS.IDENTITY (REWRITE)
  (EQUAL (M.CS FLAG (LIST (LIST L)) S)
  (M.CS FLAG L S))))

(PROVE.LEMMA M.CS.CONS (REWRITE)
  (IMPLIES (AND (NUMBERP I)
                (OR (EQUAL FLAG 'SEQ)
                    (AND (EQUAL FLAG 'PAR)
                        (IS.IND.CS I (ALL.ATOMS L))))
                (EQUAL (M.CS FLAG (CONS I L) S)
                (CS I (M.CS FLAG L S))))))

(PROVE.LEMMA M.CS.APPEND2.NIL (REWRITE)
  (EQUAL (M.CS FLAG (APPEND2 L NIL) S)
  (M.CS FLAG L S))
  ((INDUCT (M.CS FLAG L S)))))

(PROVE.LEMMA M.CS.APPEND2 (REWRITE)
  (IMPLIES (OR (EQUAL FLAG 'SEQ)
               (AND (EQUAL FLAG 'PAR)
                   (ARE.IND.CS (ALL.ATOMS L1)
                   (ALL.ATOMS L2)))
               (EQUAL (M.CS FLAG (APPEND2 L1 L2) S)
               (M.CS FLAG L1 (M.CS FLAG L2 S))))
  ((INDUCT (M.CS FLAG L1 S))))

Trace Transformation Rules

(PROVE.LEMMA G3i (REWRITE)
  (IMPLIES (ARE.IND.CS (ALL.ATOMS L1) (ALL.ATOMS L2))
  (EQUAL (M.CS 'PAR (CONS L1 L2) S)
  (M.CS 'SEQ (APPEND2 L1 (LIST L2)) S))))

(PROVE.LEMMA G3ii (REWRITE)
  (IMPLIES (AND (NOT (EQUAL L NIL))
               (ARE.IND.CS (ALL.ATOMS L1) (ALL.ATOMS L2)))
               (EQUAL (M.CS 'PAR (CONS (APPEND2 L1 L) L2) S)
               (M.CS 'SEQ (APPEND2 L1 (LIST (CONS L L2)) S)))
  ((INDUCT (APPEND2 L1 L1)))))

(ADD.AXIOM GLOBAL.IND.CS (REWRITE)
  (IMPLIES (IND.CS I J)
  (EQUAL (CS J (CS I S))
  (CS I (CS J S)))))
(PROVE.LEMMA IS.IND.CS.IMPLIES.COMMUTATIVITY (REWRITE)
  (IMPLIES (IS.IND.CS I (ALL.ATOMS L))
    (EQUAL (CS I (M.CS FLAG L S))
      (M.CS FLAG L (CS I S))))
  ((INDUCT (M.CS FLAG L S))))

(PROVE.LEMMA ARE.IND.CS.IMPLIES.COMMUTATIVITY (REWRITE)
  (IMPLIES (ARE.IND.CS (ALL.ATOMS L1) (ALL.ATOMS L2))
    (EQUAL (M.CS FLAG1 L1 (M.CS FLAG2 L2 S))
      (M.CS FLAG2 L2 (M.CS FLAG1 L1 S))))
  ((INDUCT (M.CS FLAG1 L1 S))))

Theory of Maximum

(DEFN MAX (L)
  (IF (NLISTP L)
    O
    (IF (LESSP (CAR L) (MAX (CDR L)))
      (MAX (CDR L))
      (CAR L))))

(DEFN NUMBER.LISTP (L)
  (IF (NLISTP L)
    (EQUAL L NIL)
    (AND (NUMBERP (CAR L))
      (NUMBER.LISTP (CDR L))))))

(PROVE.LEMMA NUMBER.LISTP.PLISTP (REWRITE)
  (IMPLIES (NUMBER.LISTP X) (PLISTP X)))

(PROVE.LEMMA MAX.NUMBER.LISTP (REWRITE)
  (IMPLIES (NUMBER.LISTP X) (NUMBERP (MAX L))))

(PROVE.LEMMA MAX.APPEND (REWRITE)
  (IMPLIES (AND (NUMBER.LISTP X) (NUMBER.LISTP Y))
    (EQUAL (MAX (APPEND X Y))
      (MAX (LIST (MAX X) (MAX Y))))))

(PROVE.LEMMA NUMBER.LISTP.APPEND (REWRITE)
  (IMPLIES (AND (PLISTP X) (PLISTP Y))
    (EQUAL (NUMBER.LISTP (APPEND X Y))
      (AND (NUMBER.LISTP X) (NUMBER.LISTP Y))))))

Theory of Non-Neighbors

(DEFN NON.NEIGHBORS (I J)
  (AND (NUMBERP I)
    (NUMBERP J)
    (NOT (EQUAL I J))
    (NOT (EQUAL I (ADD1 J)))
    (NOT (EQUAL (ADD1 I) J))))
(ADD.AXIOM NON.NEIGHBORS.ARE.IND.CS (REWRITE)
  (IMPLIES (NON.NEIGHBORS I J) (IND.CS I J)))

(DEFN HAS.NO.NEIGHBOR (I L)
  (IF (NLISTP L)
    T
    (AND (NON.NEIGHBORS I (CAR L))
      (HAS.NO.NEIGHBOR I (CDR L)))))

(PROVE.LEMMA HAS.NO.NEIGHBOR.IS.IND.CS (REWRITE)
  (IMPLIES (HAS.NO.NEIGHBOR I L) (IS.IND.CS I L)))

(DEEFN HAVE.NO.NEIGHBOR (L1 L2)
  (IF (NLISTP L1)
    T
    (AND (HAS.NO.NEIGHBOR (CAR L1) L2)
      (HAVE.NO.NEIGHBOR (CDR L1) L2))))

(PROVE.LEMMA HAVE.NO.NEIGHBOR.ARE.IND.CS (REWRITE)
  (IMPLIES (HAVE.NO.NEIGHBOR L1 L2) (ARE.IND.CS L1 L2)))

(PROVE.LEMMA MAX.NON.NEIGHBOR (REWRITE)
  (IMPLIES (AND (NUMBER.LISTP L)
    (LESSL (ADDI (MAX L)) I))
    (EQUAL (HAS.NO.NEIGHBOR L)
      (NON.NEIGHBORS I (MAX L))))
APPLICATION: EQUVALENCE OF TAU AND TAU^~

Trace Definitions

(DEFN SEQJ (K)
  (IF (ZEROP K)
   NIL
   (CONS K (SEQJ (SUB1 K)))))

(DEFN SEQIJ (K)
  (IF (ZEROP K)
   NIL
   (APPEND2 (SEQIJ (SUB1 K)) (SEQJ K))))

(DEFN TAU (N) (SEQIJ N))

(DEFN PARI (K)
  (IF (ZEROP K)
   NIL
   (IF (EQUAL K 1)
     (LIST 1)
     (CONS K (PARI (SUB1 (SUB1 K))))))

(DEFN T1^- (I)
  (IF (ZEROP I)
   NIL
   (APPEND2 (T1^- (SUB1 I)) (LIST (PARI I))))))

(DEFN T2^- (I)
  (IF (ZEROP I)
   NIL
   (APPEND2 (LIST (PARI I)) (T2^- (SUB1 I))))))

(DEFN TAU^- (N)
  (APPEND2 (T1^- (SUB1 N)) (T2^- N)))

(DEFN UNRAVEL.PARI (K)
  (IF (ZEROP K)
   NIL
   (IF (EQUAL K 1)
     (LIST 1)
     (CONS K (LIST (PARI (SUB1 (SUB1 K)))))))))

(DEFN UNRAVEL.T2^- (I)
  (IF (ZEROP I)
   NIL
   (APPEND2 (UNRAVEL.PARI I)
     (UNRAVEL.T2^- (SUB1 I))))))
Trace Transformation Prerequisites

(DEFINE.LEMMA NUMBER.LISTP.PARI (REWITE)
  (NUMBER.LISTP (ALL.ATOMS (PARI K))))

(DEFINE.LEMMA MAX.PARI (REWITE)
  (IMPLIES (NUMBERP K)
    (EQUAL (MAX (ALL.ATOMS (PARI K))) K)))

(DEFINE.LEMMA IND.PARI (REWITE)
  (IMPLIES (LESSP 1 K)
    (HAVE.NO.NEIGHBOR (ALL.ATOMS K)
      (ALL.ATOMS (PARI (SUB1 (SUB1 K)))))))

(DEFINE.LEMMA NUMBER.LISTP.T2^- (REWITE)
  (NUMBER.LISTP (ALL.ATOMS (T2^- I))))

(DEFINE.LEMMA MAX.T2^- (REWITE)
  (IMPLIES (NUMBERP I)
    (EQUAL (MAX (ALL.ATOMS (T2^- I))) I)))

(DEFINE.LEMMA IND.T2^- (REWITE)
  (IMPLIES (LESSP 1 I)
    (HAVE.NO.NEIGHBOR (ALL.ATOMS I)
      (ALL.ATOMS (T2^- (SUB1 (SUB1 I)))))))

Trace Identities

(DEFINE.LEMMA T2^- SHIFT (REWITE)
  (IMPLIES (LESSP 2 N)
    (EQUAL (M.CS 'SEQ
      (UNRAVEL.PARI N)
      (M.CS 'SEQ (T2^- (SUB1 (SUB1 N)))) S))
    (CS N (M.CS 'SEQ (T2^- (SUB1 (SUB1 N)))) S))))

(DEFINE.LEMMA TAU^- SHIFT (REWITE)
  (IMPLIES (LESSP 0 N)
    (EQUAL (M.CS 'SEQ
      (T1^- N)
      (M.CS 'SEQ (T2^- (SUB1 N)) S))
    (M.CS 'SEQ
      (T1^- (SUB1 N))
      (M.CS 'SEQ (T2^- N) S))))
Trace Transformations

(Prove:Lemma CLOSE-SEQJ (Rewrite)
  (Implies (Numberp N)
    (Equal (CS (Add1 N) (M.CS 'SEQ (SEQJ N) S))
     (M.CS 'SEQ (SEQJ (ADD1 N)) S)))

(Prove:Lemma TAU-RAVEL1 (Rewrite)
  (Implies (Lessp 1 N)
    (Equal (M.CS 'SEQ (UNRAVEL.T2^-N) N) S)
     (M.CS 'SEQ
      (Append2 (T2^- (Sub1 (Sub1 N))) (SEQJ N))
     S)))
  ((Disable SEQJ PARI T2^- UNRAVEL.PARI)))

(Prove:Lemma TAU-RAVEL2 (Rewrite)
  (Implies (Lessp 0 N)
    (Equal (M.CS 'SEQ (T2^- N) S)
     (M.CS 'SEQ (UNRAVEL.T2^- N) S)))
  ((Disable SEQJ TAU-RAVEL1)))

(Defn RAVEL (N S)
  (If (Or (Zerop (Sub1 N))
     (Not (Numberp N)))
   T
   (RAVEL (Sub1 N)
     (M.CS 'SEQ (SEQJ N) S)))))

(Prove:Lemma TAU-MAIN (Rewrite)
  (Implies (Lessp 0 N)
    (Equal (M.CS 'SEQ (TAU^- N) N) S)
     (M.CS 'SEQ (TAU N) S)))
  ((Disable Close-SEQJ SEQJ T2^- UNRAVEL.T2^-)
   (Induct (RAVEL N S))))

/ 

Extension of TAU-MAIN to the Empty Trace / 

(Prove:Lemma TAU^-NIL (Rewrite)
  (Implies (Not (Lessp 0 N))
    (Equal (TAU^- N) NIL)))

(Prove:Lemma TAU-MAIN-COMPLETE (Rewrite)
  (Equal (M.CS 'SEQ (TAU^- N) N) S)
     (M.CS 'SEQ (TAU N) S))
  ((Disable TAU^-)))
Appendix B: Proof Outline of TAU's Transformation

Each transformation theorem expresses an identity of two M.CS function calls. We sketch each proof by presenting the major rewrites of the left M.CS call - these rewrites are named L1 to Lm, and the right M.CS call - these rewrites are named R1 to Rn. Lm and Rn yield matching calls.

TAU.RAVEL1

\[(M.CS \ 'SEQ \ (UNRAVEL.T2^N) \ S)\]

\[(L1) \ = \ (M.CS \ 'SEQ \ (UNRAVEL.PARI N) \ (M.CS \ 'SEQ \ (UNRAVEL.T2^N-1) \ S))\]

\[(L2) \ = \ (M.CS \ 'SEQ \ UNRAVEL.PARI N) \ (M.CS \ 'SEQ \ APPEND2 \ (UNRAVEL.T2^N-3) \ (SEQJ N-1) \ S))\]

\[(L3) \ = \ (M.CS \ 'SEQ \ (UNRAVEL.PARI N) \ (M.CS \ 'SEQ \ UNRAVEL.T2^N-3) \ (M.CS \ 'SEQ \ (SEQJ N-1) \ S)))\]

\[(L4) \ = \ (CS N \ (M.CS \ 'SEQ \ (T2^N-2) \ (M.CS \ 'SEQ \ (SEQJ N-1) \ S)))\]

\[(L5) \ = \ (M.CS \ 'SEQ \ (T2^N-2) \ (CS N \ (M.CS \ 'SEQ \ (SEQJ N-1) \ S)))\]

\[(L6) \ = \ (M.CS \ 'SEQ \ (T2^N-2) \ (M.CS \ 'SEQ \ (SEQJ N) \ S))\]

\[(M.CS \ 'SEQ \ APPEND2 \ (T2^N-2) \ (SEQJ N)) \ S)\]

\[(R1) \ = \ (M.CS \ 'SEQ \ (T2^N-2) \ (M.CS \ 'SEQ \ (SEQJ N) \ S))\]

(L1) by UNRAVEL.T2^N and M.CS.APPEND2
(L2) by induction hypothesis for N-1
(L3) by M.CS.APPEND2
(L4) by T2^N.SHIFT
(L5) by IND.T2^N, IS.IND.CS.IMPLIES.COMMUTATIVITY
(L6) by CLOSE.SEQJ
(R1) by M.CS.APPEND2
TAU.RAVEL2

(M.CS 'SEQ (T2~ N) S)

(L1)  = (M.CS 'PAR (PARI N)
       (M.CS 'SEQ (T2~ N-1) S))

(L2)  = (M.CS 'PAR (PARI N)
       (M.CS 'SEQ (UNRAVEL.T2~ N-1) S))

(L3)  = (M.CS 'PAR (CONS N (PARI N-2))
       (M.CS 'SEQ (UNRAVEL.T2~ N-1) S))

(L4)  = (M.CS 'SEQ (CONS N (LIST (PARI N-2)))
       (M.CS 'SEQ (UNRAVEL.T2~ N-1) S))

       (M.CS 'SEQ (UNRAVEL.T2~ N) S)

(R1)  = (M.CS 'SEQ (UNRAVEL.PARI N)
       (M.CS 'SEQ (UNRAVEL.T2~ N-1) S))

(R2)  = (M.CS 'SEQ (CONS(LIST (PARI N-2)))
       (M.CS 'SEQ (UNRAVEL.T2~ N-1) S))

(L1)  by T2~, M.CS, and M.CS.APPEND2
(L2)  by induction hypothesis for N-1
(L3)  by PARI
(L4)  by IND.PARI, G3i, and APPEND2
(R1)  by UNRAVEL.T2~ and M.CS.APPEND2
(R2)  by UNRAVEL.PARI
TAU.MAIN

(M.CS 'SEQ (TAU N) S)

(L1) = (M.CS 'SEQ (T1° N-1)
        (M.CS 'SEQ (T2° N) S))

(L2) = (M.CS 'SEQ (T1° N-1)
        (M.CS 'SEQ (UNRAVEL.T2° N) S))

(L3) = (M.CS 'SEQ (T1° N-1)
        (M.CS 'SEQ (T2° N-2)
        (M.CS 'SEQ (SEQJ N) S)))

(L4) = (M.CS 'SEQ (T1° N-2)
        (M.CS 'SEQ (T2° N-1)
        (M.CS 'SEQ (SEQJ N) S)))

(L5) = (M.CS 'SEQ (SEQIJ N-1)
        (M.CS 'SEQ (SEQJ N) S))

(R1) = (M.CS 'SEQ (SEQIJ N-1)
        (M.CS 'SEQ (SEQJ N) S))

(L1) by TAU° and M.CS.APPEND2
(L2) by TAU.RAVEL2
(L3) by TAU.RAVEL1 and M.CS.APPEND2
(L4) by TAU°.SHIFT
(L5) by induction hypothesis for N-1, TAU, and TAU°
(R1) by TAU and SEQIJ