

DECOMPOSABILITY AND ITS ROLE IN PARALLEL LOGIC-PROGRAM EVALUATION

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Decomposability and Its Role in Parallel Logic-Program Evaluation

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ABSTRACT

This paper is concerned with the issue of parallel evaluation of logic programs. We define the concept of *program decomposability*, which means that the load of evaluation can be partitioned among a number of processors, without a need for communication among them. This in turn results in a very significant speed-up of the evaluation process. Some programs are decomposable, whereas others are not. We completely syntactically characterize three classes of single rule programs with respect to decomposability: nonrecursive, simple linear, and simple chain programs. We also establish two sufficient conditions for decomposability.

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1. Introduction

We propose a new method of evaluating logic programs in parallel. The method is suitable for sharing the computation load among an arbitrary number of processors, which either have common memory or communicate by message passing. This makes it applicable to a large class of hardware architectures. Let us demonstrate the method using the classical example of the program computing the transitive closure of a graph. The arcs of the graph are given by the tuples of a database relation A . The program is written in DATALOG (see [MW]):

$$\begin{aligned} T(x,y) &:- T(x,z), A(z,y) \\ T(x,y) &:- A(x,y). \end{aligned}$$

If the relation A is replicated at two different processors, $p1$ and $p2$, we can partition the work of computing (the relation for) the predicate T as follows. The above program, with the arithmetic predicate $even(x)$ appended to the body of the second rule, is assigned to processor $p1$. In other words, $p1$ executes the program:

$$\begin{aligned} T(x,y) &:- T(x,z), A(z,y) \\ T(x,y) &:- A(x,y), even(x). \end{aligned}$$

On the other hand, processor $p2$ executes the program:

$$\begin{aligned} T(x,y) &:- T(x,z), A(z,y) \\ T(x,y) &:- A(x,y), odd(x). \end{aligned}$$

The effect of the evaluation of $p1$ is that it computes the tuples (x,y) of the transitive closure, in which x is even. Similarly, $p2$ computes those tuples in which x is odd. For example, if the input graph is $1 \rightarrow 2 \rightarrow 3 \rightarrow 4 \rightarrow 5$, then $p1$ computes the set of output tuples $\{ (1,2), (1,3), (1,4), (3,4), (3,5) \}$, and $p2$ computes the set $\{ (2,3), (2,4), (2,5), (4,5) \}$.

A moment of reflection will reveal that the above partitioning of the work has several nice properties. First, no processor computes a tuple which is also computed by the other processor, thus there is no work-duplication in this sense. Second, if the relation computed by each processor is output to the same device, or stored in the same file, the result is always the complete transitive closure, regardless of the input graph. Third, no communication between the two processors is required during the computation. Fourth, the work-partitioning does not require complicated program transformations, only adding evaluable predicates to the body of some rules of the original program.

Assume that the whole relation for T has to be evaluated, and $p1$ and $p2$ start at the same time and execute their programs in parallel. Assume further that at the same time a single processor, using the original program, starts the evaluation of T . It is quite intuitive that, for an "average" (large enough) graph, the partitioned evaluation of T will complete much sooner than the single-processor evaluation. Furthermore, note that the evaluation can be divided among k processors, for any $k \geq 2$. The only difference from the above example is that processor p_i executes a copy of the program with the predicate $i \bmod k(x)$ added to the nonrecursive rule. The exact time-speedup achieved by the work-partitioning scheme depends on many parameters outside the scope of this paper, however, here we are interested in a qualitative issue.

We postulate that in general, a work-partitioning scheme with the properties enumerated above, is very desirable. If it can be applied to the evaluation of a predicate in a program, then

we say that the predicate is decomposable. Not every predicate is decomposable. Even for the same problem of computing the transitive closure, we will prove that the predicate T in the program:

$$\begin{aligned} T(x,y):- T(x,z),T(z,y) \\ T(x,y):- A(x,y) \end{aligned}$$

is not decomposable. The proof of this fact will be given in section 6. This indicates that decomposability is a syntactic rather than semantic property. We feel that it is both practically and theoretically important to first formally define decomposability, and then characterize the decomposable predicates.

In this paper we completely characterize three subclasses of single rule programs (sirups) with respect to decomposability: nonrecursive, simple linear, and simple chain programs. Sirups were first studied as a syntactically restricted class of programs by Cosmadakis and Kanellakis ([CK]). They have only one output predicate, therefore we interchangeably use the term decomposability of a predicate or of a program. We also provide two sufficient conditions for any sirup to be decomposable. Simple linear programs and simple chain programs are important subclasses of sirups from the practical point of view.

This work is related to the general subject of parallel evaluation of logic programs. The subject has recently emerged as a very important and active area of research ([AP], [K], [U], [UV]). Most existing research is concerned with membership in the complexity class NC. This class is a mathematical tool for analyzing parallel algorithms in general. Here we show that for analyzing parallel evaluation of logic programs, a different tool can be used. Loosely speaking, if a logic program is in NC it does not guarantee that it has all the nice properties of a decomposable predicate. In particular, the processors executing an NC type algorithm usually have to communicate extensively, and therefore communication is assumed to take place through common memory. Also, a speedup for such an algorithm is not guaranteed unless the number of processors is polynomial in the size of the input.

The remainder of the paper is organized as follows. In section 2 we introduce the necessary definitions and notations used throughout the paper. In section 3 we prove that any nonrecursive sirup is decomposable. In section 4 we provide two sufficient conditions for a general sirup to be decomposable, and in section 5 we show that one of these conditions, called pivoting, is also necessary for decomposability of a simple linear sirup. In section 6 it is proven that a simple chain program is decomposable if and only if it is regular. In section 7 we discuss future work.

2. Preliminaries

In this section we present the basic definitions and terminology that will be used throughout this paper.

2.1. Program Structure

An *atom* is a predicate symbol with a constant or a variable in each argument position. We assume that the constants are the natural numbers. An *R-atom* is an atom having R as the predicate symbol. A *rule* consists of an atom, Q , designated as the *head*, and a conjunction of one or more atoms, denoted Q^1, \dots, Q^k , designated as the *body*. Such a rule is denoted

$Q :- Q^1, \dots, Q^k$, which should be read "Q if Q^1 and Q^2 , and, ..., and Q^k ." A rule or an atom is an *entity*. If an entity has a constant in each argument position, then it is a *ground* entity. For a predicate symbol, R , a finite set of R -ground-atoms is a *relation* for R .

A DATALOG program P , or a program for short, is a finite set of rules whose predicate symbols are divided into two disjoint subsets: the *base* predicates, and the *derived* predicates. The base predicates are distinguished by the fact that they do not appear in any head of a rule. An *input* to P is a relation for each base predicate. An *output* of P is a relation for each derived predicate of P . A *substitution* applied to an entity, or a sequence of entities, is the replacement of each variable in the entity by a variable or a constant. It is denoted $x_1/y_1, x_2/y_2, \dots, x_n/y_n$ indicating that x_i is replaced by y_i . A substitution is *ground* if the replacement of each variable is by a constant. A ground substitution applied to a rule is an *instantiation* of the rule.

A *database* for P is a relation for each predicate of P . The output of P , given an input I , is the set of relations for the derived predicates in the database, obtained by the following procedure, called *bottom up evaluation* :

- (1) Start with an initial database consisting of the relations of I .
- (2) If there is an instantiation of a rule of P such that all the ground atoms in the body are in the database generated so far, and the one in the head is not, then: add to the database the ground atom in the head of the instantiated rule, and reexecute (2).
- (3) Stop.

This procedure is guaranteed to terminate, and produce a finite output for any given P and I ([VEK]). The output is unique, in the sense that any order in which *bottom up evaluation* adds the atoms to the database will produce the same output. For simplicity we assume that the rules of a program are *range restricted*, i.e., every variable in the head of a rule also appears in the body of that rule. Furthermore, we assume that the rules do not have constants, and each query is to evaluate a whole relation for a predicate.

A predicate Q in a program P *derives* a predicate R , if Q occurs in the body of a rule whose head is a R -atom. Q is *recursive* if (Q, Q) is in the nonreflexive transitive closure of the "derives" relation. A program is recursive if it has a recursive predicate. A rule is recursive if the predicate in its head transitively derives some predicate in its body.

A *single rule program* (see [CK]) is a DATALOG program having a single derived predicate, denoted S in our paper, and consisting of:

1. A nonrecursive rule,

$$S(x_1, \dots, x_n) :- B(x_1, \dots, x_n)$$

where the x_i 's are distinct variables.

2. One other, possibly recursive, rule in which the predicate symbol B does not appear.

2.2. Restricted Versions of Programs

An *evaluable predicate* is an arithmetic predicate (see [BR]). Examples of evaluable predicates are sum, greater than, modulo, etc. A rule re is a *restricted version* of some rule r , if r and re have exactly the same variables, and r can be obtained by omitting zero or more evaluable

predicates from the body of re . In other words, re is r with some evaluable predicates added to the body, and the arguments of these evaluable predicates are variables of r . For example, if r is:

$$S(x,y,z):-S(w,x,y), A(w,z)$$

then one possible re rule is:

$$S(x,y,z):-S(w,x,y), A(w,z), x-y=5.$$

A program P_i is a *restricted version* of program P if each one of its rules is a restricted version of some rule of P . Note that P_i may have more than one restricted version of a rule r of P . To continue the above example, if P has the rule r , then P_i may have the rule re as well as the rule re' :

$$S(x,y,z):-S(w,x,y), A(w,z), x-y=6.$$

Throughout this paper, only restricted versions of a program may have evaluable predicates.

The input of a program with evaluable predicates, i.e. a restricted version, is defined as before. The output is also defined as before, except that step (2) of the procedure bottom-up-evaluation also verifies that the substitution satisfies the evaluable predicates in the ground rule; only then the atom in the head is added to the database and step (2) is reexecuted. For example, for the rule re' above, the substitution $x/14,y/8$ satisfies the evaluable predicate $x-y=6$, whereas the substitution $x/13,y/9$ does not do so.

3. Decomposability

In this section we first define and discuss the key notion of decomposability, then prove that a nonrecursive sirup is decomposable. Let P be a program, let P_1, \dots, P_r be restricted copies of P , and let T be a derived predicate of P . We denote by T_i the relation output by P_i for T . (Observe that this is a somewhat unconventional notation, since the relation name is different than the predicate name).

We say that predicate T is *decomposable* in P with respect to P_1, \dots, P_r if the following two conditions hold:

1. For each input I to P, P_1, \dots, P_r
 - i. $\bigcup_i T_i \supseteq T$ (completeness).
 - ii. T_i and T_j are disjoint for each $i \neq j$; furthermore, if some derived predicate Q transitively derives T in P , then Q_i and Q_j are disjoint (lack-of-duplication).
2. For some input I to P_1, \dots, P_r , each T_i is nonempty (nontriviality).

The above definition is central to this paper, and we shall discuss it next.

Requirement 1.i states that no output is lost by evaluating the relation for T in each P_i rather than the relation for T in P ; the fact that no additional output is generated is implied by the fact that each P_i is a restricted version of P . Requirement 1.ii states that in the process of evaluating T , each new ground atom (or intermediate result) is computed by a unique processor. Assume that, along the lines suggested in [BR section 4], we measure the cost of evaluating the relation T , in terms of the number of new ground atoms generated in the evaluation process. Then, loosely speaking, requirement 1 says the following. For every input (i.e. set of base

relations replicated at each processor), the evaluation by r processors is equivalent, in terms of the output produced and the total evaluation cost, to the single-processor evaluation.

The strength of requirement 1 enables the relaxed form of requirement 2. It is enough that for "some" inputs each T_i is nonempty, since for those inputs the evaluation cost incurred by each processor is smaller than that of a single processor executing the program P . Then the evaluation of T completes sooner in the distributed case. In other words, since there is nothing to lose by distributing the computation, it is enough that we gain only in some cases to make the scheme worthwhile. However, for the decomposable predicates that we discuss in this paper, nontriviality holds for more than an isolated case input.

For instance, in the transitive closure example nontriviality holds for any input graph in which arcs exit both, even and odd nodes. Specifically, for the class of predicates that we prove decomposable in this paper, decomposability is shown using the *odd-even* predicates alone. This has two implications. First, the work performed by each processor for an arbitrary input, is roughly equal (e.g. for an arbitrary graph, the number of odd and even nodes is roughly equal). In these cases we expect the distributed evaluation to be faster than the single-processor evaluation, by a factor which is close to two, i.e. the number processors. Second, note that the *odd* and *even* predicates are a special case of the $i \bmod r$ predicates, for $r=2$. When we show that T is decomposable in P with respect to P_1 and P_2 , then it should be easy for the readers to convince themselves that for any r , there are restricted copies P_1, \dots, P_r such that T is decomposable in P with respect to P_1, \dots, P_r . This means that the work can be divided among any number of processors. For instance, in the transitive closure example, in order to do so processor i evaluates T_i where:

$$P_i. \quad T(x,y):-T(x,z),A(z,y). \\ T(x,y):-A(x,y), x=i \bmod r.$$

These facts stress the robustness of the decomposability definition.

We say that predicate T is *decomposable* in P if it is decomposable with respect to some restricted copies P_1, \dots, P_r for $r > 1$.

Theorem 1: If a sirup P is nonrecursive, then its derived predicate is decomposable.

Proof: Assume that P is:

$$S(x_1, \dots, x_n):-Q^1(\dots), \dots, Q^k(\dots) \\ S(x_1, \dots, x_n):-B(x_1, \dots, x_n)$$

where B and each Q^i are base predicates. Consider the following restricted copies of P :

$$P_1. \quad S(x_1, \dots, x_n):-Q^1(\dots), \dots, Q^k(\dots), \text{even}(x_1) \\ S(x_1, \dots, x_n):-B(x_1, \dots, x_n), \text{even}(x_1) \\ P_2. \quad S(x_1, \dots, x_n):-Q^1(\dots), \dots, Q^k(\dots), \text{odd}(x_1) \\ S(x_1, \dots, x_n):-B(x_1, \dots, x_n), \text{odd}(x_1).$$

It is easy to see that S is decomposable in P with respect to P_1 and P_2 . \square

4. Sufficient Conditions for Decomposability

In this section we provide two sufficient conditions for decomposability of a general sirup. The first one is motivated by the next example, which also merits attention for the following reason. From the preceding discussion one might suspect that our notion of decomposability is equivalent to "naive" propagation of variable bindings (see introduction of [BKBR]). The latter notion means simply substituting a constant for a variable in some rules. The constant is usually taken from a query. For example, in order to find all the arcs exiting the node 2 in the transitive closure of a graph, the constant can be naively propagated into the program as follows:

$$\begin{aligned} T(2,y):- T(2,z),A(z,y) \\ T(2,y):- A(2,y). \end{aligned}$$

It is quite clear that if a sirup is amenable to naive propagation of variable bindings, then it is decomposable. However, the reverse is not true. For example, consider the program:

$$\begin{aligned} S(x,y):- S(y,x) \\ S(x,y):- A(x,y). \end{aligned}$$

which outputs an arc in both directions for every arc of an input graph. It is easy to see that a binding cannot be naively propagated into this program, but the sirup is decomposable; one restricted copy has the nonrecursive rule:

$$S(x,y):- A(x,y),\text{even}(x+y).$$

and the other:

$$S(x,y):- A(x,y),\text{odd}(x+y).$$

Note that appending to the body of the nonrecursive rule the predicates $\text{odd}, \text{even}(x*y)$, or any other commutative function of x and y , works as well. Our first sufficient condition for decomposability, introduced below, is based on the preceding observation.

Let R be a set of atoms, each of which has a variable in each argument position. The set R is *pivoting* if there is a subset d of argument positions, such that in the positions of d :

1. The same variables appear (possibly in a different order) in all atoms of R , and
2. Each variable appears the same number of times in all atoms of R .

A member of d is called a *pivot*. Note that a variable that appears in a pivot may or may not appear in a nonpivot position of the same atom.

The recursive rule of a sirup is *pivoting* if all the occurrences of the recursive predicate in the rule constitute a pivoting set. For example, the rule

$$S(w,x,x,y,z) :- S(u,y,x,x,w), S(v,x,y,x,w), A(u,v,z)$$

is pivoting, with argument positions 2, 3 and 4 of S being the pivots.

Theorem 2: If the recursive rule of a sirup is pivoting, then the sirup is decomposable.

Proof: Assume that argument positions i_1, \dots, i_k of S are the pivots. Consider restricted copy P_1 of P which has the same recursive rule as P , and a nonrecursive rule

$$S(x_1, \dots, x_n) :- B(x_1, \dots, x_n), \text{even}(x_{i_1} + x_{i_2} + \dots + x_{i_k}).$$

Restricted copy P_2 of P is the same, except that the nonrecursive rule is

$$S(x_1, \dots, x_n) :- B(x_1, \dots, x_n), \text{odd}(x_{i_1} + x_{i_2} + \dots + x_{i_k}).$$

Assume that for input I , the ground atom $a=S(c_1, \dots, c_n)$ is in the relation S output by P . Assume further, without loss of generality, that $c=c_{i_1} + \dots + c_{i_k}$ is even. Denote by t the necessary and sufficient number of iterations of step (2) of bottom-up-evaluation for adding a to the database, in evaluating P . It is easy to see by induction on t , that t iterations are necessary and sufficient to add a to S_1 . It is also easy to see that a is not in S_2 , and that nontriviality holds. \square

Theorem 2 can be extended to general DATALOG programs, not necessarily sirups, provided that they do not have repeated variables in the heads of rules. A rule in such a program is pivoting, if all its derived-predicate-atoms (in the head and the body) constitute a pivoting set. A program is pivoting if each one of its rules is pivoting, with the same argument positions being the pivots in all the rules. For example, the program

$$\begin{aligned} S(x,y,z) &:- R(y,x,w), A(w,z) \\ R(x,y,z) &:- R(x,y,w), B(w,z) \\ R(x,y,z) &:- C(x,y,z) \end{aligned}$$

is pivoting, with positions 1 and 2 being the pivots. A predicate in such a program is decomposable if the rules which derive the predicate constitute a pivoting program. For example, predicate S in the program above is decomposable (add $\text{odd-even}(x+y)$ to the body of the third rule).

The condition of theorem 2 is not necessary for decomposability. For example, the sirup

$$\begin{aligned} S(x,x) &:- S(y,y), A(x,y) \\ S(x,y) &:- B(x,y) \end{aligned}$$

is obviously not pivoting, but it is decomposable. Again, $\text{odd-even}(x+y)$ is added to the body of the nonrecursive rule. The intuition indicates that in this example the computation load for an arbitrary input is not evenly divided between the processors executing the two restricted versions of the program (because only the processor executing the copy with the *even* evaluable predicate can output an atom as a result of instantiation of the recursive rule). The example is unique (throughout the paper) in this respect. Expectedly, the last example motivates our next sufficient condition for decomposability of a sirup. It is defined as follows. Assume that R is a set of atoms with each atom having the same predicate symbol, Q , and a variable in each argument position. The set R is *repeating* if there are at least two argument positions of Q , i and j , such that the same variable appears in position i and position j , and this is true for each member of R (note that the variable of one member of R may be different than the variable of another). The recursive rule of a sirup is *repeating* if all the occurrences of the recursive predicate in the rule constitute a repeating set. For example, the rule

$$S(x,z,x) :- S(z,z,z), S(x,x,x)$$

is repeating because of argument positions 1 and 3.

Theorem 3: If the recursive rule of a sirup is repeating, then the sirup is decomposable.

Proof: Very similar to the proof of theorem 2 thus omitted. The only difference between the proofs is that $\text{odd-even}(xi + xj)$ replaces $\text{odd-even}(xi_1 + \dots + xi_k)$, where i and j are the positions of the repeated variable. \square

Obviously, the condition of theorem 3 is not necessary for decomposability either.

5. Simple Linear Sirups

In this section we completely characterize the class of simple linear sirups with respect to decomposability. A sirup is *linear* if it is recursive, and in the body of the recursive rule there is exactly one occurrence of the recursive predicate. A linear sirup is *simple* if it does not have repeated variables in an occurrence of the recursive predicate.

The characterization of simple linear sirups with respect to decomposability is done by proving that the sufficient condition of theorem 2 is also necessary. We assume that the recursive rule is:

$$S(x_1, \dots, x_n) :- S(Y_1, \dots, Y_n), A_1(\dots), \dots, A_k(\dots)$$

where the A_i 's are base predicates. Observe the notation used in this section to distinguish between two types of variables. The ones starting with a lowercase letter are logic program variables, or variables for short, as before. The ones starting with an upper case letter (e.g. Y_1), are *metalinguistic-variables*. They denote program variables. For example, Y_1 may denote the variable x_n .

If the predicate $S(x_1, \dots, x_n)$ in a (not necessarily linear) sirup P is decomposable with respect to P_1, \dots, P_r , then we define the *home-site* of a sequence of n constants, $\bar{c} = c_1, \dots, c_n$. It is the S_i to which the output atom $S(\bar{c})$ belongs, if each P_i is given the input consisting of a unique atom, $B(\bar{c})$. Note that the home-site of a sequence is unique (lack-of-duplication), every sequence of n constants has a home-site (completeness), and each S_i , $1 \leq i \leq r$, has a sequence of constants for which S_i is the home-site. Let $\bar{c} = c_1, \dots, c_n$ and $\bar{d} = d_1, \dots, d_n$ be two sequences of constants. The ordered pair of ground atoms $\langle S(\bar{d}), S(\bar{c}) \rangle$ is a *one-step-derivation* if there is an instantiation of the recursive rule of P , in which the first atom is in the head and the second is in the body.

Lemma 1: If the derived predicate, S , of a simple linear sirup, P , is decomposable, and there are two sequences of constants $\bar{d} = d_1, \dots, d_n$ and $\bar{c} = c_1, \dots, c_n$ such that $\langle S(\bar{d}), S(\bar{c}) \rangle$ is a one-step-derivation, then the home-site of \bar{d} and \bar{c} is identical.

Proof: Let the instantiation of the recursive rule which results in the one-step-derivation be:

$$S(\bar{d}) :- S(\bar{c}), A_1(\bar{a}_1), \dots, A_k(\bar{a}_k)$$

Consider the input:

$$I_1 = \{A_1(\bar{a}_1), \dots, A_k(\bar{a}_k), B(\bar{c})\}.$$

Assume that P is decomposable with respect to P_1, \dots, P_r . Note that $S(\bar{c})$ must be in the output of some restricted version of P . Assume that $S(\bar{c})$ is in S_j . By completeness, $S(\bar{d})$ is also in the output of some restricted version. This output must be of P_j , for the following reason. In the input I_1 there is only one B -ground-atom, therefore the output of any restricted version other than P_j is empty. Now consider the input $I_2 = I_1 \cup \{B(\bar{d})\}$. By lack of duplication, for the input I_2 the ground atom $S(\bar{d})$ is still in S_j . Therefore, for the input consisting of the single atom $B(\bar{d})$, the output ground atom $S(\bar{d})$ must be in S_j . \square

Let P be a simple linear sirup, having the recursive predicate denoted S , and the recursive rule denoted r . Let us define the sequence of S -atoms *Distinct-Vars* as follows. The first member, m_0 , is $S(x_1, \dots, x_n)$, where the x_i 's are variables. Subsequently, member m_i is defined as the head of the recursive rule, r' , obtained by applying to r a substitution which satisfies the following two conditions:

1. Each one of the variables in the S -atom in the body of r is replaced by another variable, such that m_{i-1} appears in the body of r' .
2. Each one of the other variables in r is replaced by a distinct variable that does not appear in m_1, m_2, \dots, m_{i-1} .

For example, consider the recursive rule: $S(x_1, x_2, x_3) :- S(x_4, x_1, x_2), A(x_4, x_3)$. Then the following is a prefix of the sequence *Distinct-Vars*: $S(x_1, x_2, x_3)$, $S(x_2, x_3, x_4)$, $S(x_3, x_4, x_5)$, and $S(x_4, x_5, x_6)$.

We shall prove that S is not decomposable, if the sequence *Distinct-Vars* has a member in which none of the variables is one of the x_i 's; then we shall prove that if so, then the recursive rule of P is not pivoting.

By definition of *Distinct-Vars*, we immediately obtain the following.

Lemma 2: Assume that $S(Y_1, \dots, Y_n)$ and $S(Z_1, \dots, Z_n)$ are two consecutive members of *Distinct-Vars*. Furthermore, assume that there is a ground substitution ρ of the program variables in the sequence $S(Y_1, \dots, Y_n)$, $S(Z_1, \dots, Z_n)$, resulting in the sequence of ground atoms $S(c_1, \dots, c_n)$, $S(d_1, \dots, d_n)$. Then the pair $\langle S(d_1, \dots, d_n), S(c_1, \dots, c_n) \rangle$ is a one-step-derivation.

Lemma 3: Assume that m_i is a member of *Distinct-Vars*, such that no variable in the set x_1, \dots, x_n appears in m_i . Then P is not decomposable.

Proof: Assume that P is decomposable with respect to P_1, \dots, P_r , and let c_1, \dots, c_n and d_1, \dots, d_n be two arbitrary sequences of constants. We will show that both have the same home-site, contradicting nontriviality. Consider the sequence of atoms $s: m_1 = S(x_1, \dots, x_n), \dots, m_i = S(Z_1, \dots, Z_n)$. The substitution $\rho = x_1/c_1, \dots, x_n/c_n, Z_1/d_1, \dots, Z_n/d_n$ is valid for any values of $c_1, \dots, c_n, d_1, \dots, d_n$, since $\{x_1, \dots, x_n, Z_1, \dots, Z_n\}$ is a set of distinct variables. Let ρ' be an extension of ρ to a ground substitution of the sequence s . The sequence of ground atoms $s\rho'$ has the property that any two consecutive atoms in it constitute a one step derivation (by Lemma 2). Therefore, by Lemma 1, the constant-sequences c_1, \dots, c_n and d_1, \dots, d_n have the same home-site. \square

Lemma 4: If the recursive rule of P is not pivoting, then there is a member, m_i , of *Distinct-Vars*, such that no variable in the set $\{x_1, \dots, x_n\}$ appears in m_i .

Proof: Construct a graph, G , in which the nodes are the argument positions of S , and there is an additional node called "new". There is an edge from p to q if the same variable appears in position p in the occurrence of S in the body of the recursive rule, and in position q in the occurrence of S in the head. If in position q in the occurrence of S in the head there is a variable which does not appear in the occurrence of S in the body, then draw an arc from "new" to q . Every node except "new" has exactly one entering arc and one exiting arc, because there are no repeated variables in an occurrence of the recursive predicate. It is easy to see that if G has a

cycle, then P is pivoting, with the nodes of the cycle being the pivots. Since P is not pivoting, G is acyclic, and we conclude that there must be a path from "new" to every other node in G . Assume that the shortest path from "new" to some other node, p , is of length k . It can be shown by induction on k , that position p of m_k will have a variable which is not in the set $\{x_1, \dots, x_n\}$. By definition of *Distinct-Vars* the lemma follows. \square

Theorem 4: A simple linear sirup is decomposable if and only if its recursive rule is pivoting.

Proof: (if) Special case of Theorem 2.

(only if) Immediate from Lemmas 3 and 4. \square

6. Simple Chain Programs

A simple chain program is a recursive sirup in which:

- (a) All the predicates are binary.
- (b) The argument positions in the left hand side of the recursive rule have distinct variables, and these variables appear in the first argument position of the first atom in the body, and in the last argument position of the last atom, respectively.
- (c) All the argument positions in the body of the recursive rule have distinct variables, except that the first argument position of the second atom has the same variable as the last argument position of the first atom, the first argument position of the third atom has the same variable as the last argument position of the second atom, etc.

For example, the following is a simple chain program:

$$S(x,y):- A(x,z_1),S(z_1,z_2),S(z_2,z_3),C(z_3,z_4),D(z_4,y)$$

$$S(x,y):- B(x,y)$$

where the A,B,C,D are base relations. A simple chain program is *regular* if in its recursive rule there is one occurrence of the predicate S and this occurrence is the first or the last in the body of the recursive rule. Note that a simple chain program is pivoting if and only if it is regular.

Theorem 5: A simple chain program P is decomposable if and only if it is regular.

Proof: (if) Immediate, based on Theorem 2.

(only if) Assume that P is not regular, and is decomposable with respect to restricted copies P_1, P_2, \dots, P_r of P , for $r > 1$. Denote the recursive rule of P by:

$$S(x,y):- Q^1(x,z_1), \dots, Q^t(z_{t-1},y)$$

where some of the Q^i 's are S 's, and $t > 1$. Using the usual notation, the nonrecursive rule is:

$$S(x,y):- B(x,y).$$

By nontriviality there are two sequences of constants, j_1, k_1 and j_2, k_2 with home sites S_1 and S_2 respectively. Since the recursive rule of P is not regular, there are two cases to analyze:

Case 1: There is a subsequence in the body of the recursive rule, of the following form:

$$Q^{i-1}(z_{i-2},z_{i-1}),S(z_{i-1},z_i),Q^{i+1}(z_i,z_{i+1}).$$

Let I^2 consist of the set of ground atoms:

$$Q^1(c_1, c_2), Q^2(c_2, c_3), \dots, Q^{i-1}(c_{i-1}, j_1), B(j_1, k_1), Q^{i+1}(k_1, c_{i+1}), Q^{i+2}(c_{i+1}, c_{i+2}), \dots, Q^t(c_{t-1}, c_t)$$

where:

- 1) Each predicate $S(m, n)$ in the list is a notation for $B(m, n)$.
- 2) Each pair of different c 's represents different constants.
- 3) None of the c 's is in the set $\{j_1, k_1, j_2, k_2\}$.

For the input I^2 , the ground atom $S(c_1, c_t)$ is in the output S of P . By completeness, for this input, $S(c_1, c_t)$ is in some S_j . We will show that $S(c_1, c_t)$ is in S_1 . Assume otherwise, i.e. $S(c_1, c_t)$ is in S_b for $b \neq 1$. The atom $B(c_1, c_t)$ is not in I^2 because $t > 1$, therefore $S(c_1, c_t)$ must be added to the database by instantiating the recursive rule of P_b in step 2 of bottom-up-evaluation. However, to generate an atom using the recursive rule of P_b , requires a 'chain' of atoms of length t . But $I^2 - B(j_1, k_1)$ does not contain such a chain, since it only contains $t-1$ atoms, and has no cycles (by the choice of constants).

Therefore, for input I^2 to P_1, \dots, P_r , $S(c_1, c_t)$ is in S_1 . Now consider the input I^3 , which is defined identically to I^2 , except that the constants j_1, k_1 are replaced by j_2, k_2 respectively. Similar arguments as before will reveal that $S(c_1, c_t)$ is in S_2 . The proof of this case is completed by noticing that for the input $I^2 \cup I^3$, the ground atom $S(c_1, c_t)$ is in both, S_1 and S_2 , contradicting lack-of-duplication.

Case 2: The body of the recursive rule of P is of the following form:

$$S(x, z), S(z, y).$$

Consider the input I^4 consisting of the ground atom $B(j_3, k_3)$, where j_3 and k_3 are distinct, and none of them is in the set $\{j_1, j_2, k_1, k_2\}$. Assume without loss of generality that the home site of j_3, k_3 is S_d , for $d \neq 1$ (otherwise the analysis below can be carried out by replacing j_1, k_1 by j_2, k_2 respectively).

Subcase 2.1: Assume that $j_1 \neq k_1$. Let input $I^5 = \{B(j_1, k_1), B(k_1, j_3), B(j_3, k_3)\}$. This input relation can be regarded as a graph consisting of a path, therefore $S(j_1, j_3)$ and $S(k_1, k_3)$ are in S . Assume that the home site of k_1, j_3 is S_i for $i \neq 1$. But then it is easy to see that $S(j_1, j_3)$ is not in any S_i ; contradicting completeness. If the home site of k_1, j_3 is S_1 , then it is easy to see that for input I^5 the atom $S(k_1, k_3)$ is not in any S_i ; again contradicting completeness.

Subcase 2.2: Assume that $j_1 = k_1$. In other words, the home site of j_1, j_1 , is S_1 . Let input $I^6 = \{B(j_1, j_3), B(j_3, k_3), B(k_3, j_1)\}$. This input relation can be regarded as a graph consisting of a cycle, therefore $S(j_1, j_1)$ is in S . Since $S(j_3, k_3)$ is in S_d only, for the input I^6 , $S(j_1, j_1)$ cannot be in any S_i other than S_d . But then, for the input $I^6 \cup \{B(j_1, j_1)\}$ the ground atom $S(j_1, j_1)$ is in both, S_1 and S_d ; this contradicts lack of duplication. \square

Note that theorem 5, combined with the results in [AP, UV], indicate that in the class of simple chain programs, the subclass of programs in NC properly contains the subclass of decomposable programs. The reason for this is that, clearly, every regular chain program is in NC, and the program:

$$S(x, y):- A(x, z_1), S(z_1, z_2), A(z_2, z_3), S(z_3, y)$$

$$S(x,y):-B(x,y)$$

is one of the programs in NC that is not decomposable. Recent results from [CW] indicate that, outside the class of simple chain programs, there are decomposable programs that are P-complete.

7. Future Work

We shall continue the work on decomposability in several directions. One of them is to extend the characterization of decomposable predicates to other sirups first, e.g. typed (see [K]), and then to general logic programs. Another direction is to determine whether decomposition implies that the work can be evenly divided among the processors, as we have seen that can be done using the *mod* predicate. For this purpose a notion of *fair* decomposition should be defined. Another topic which merits attention is minimizing communication when evaluating nondecomposable predicates in a distributed environment. We feel that the work on decomposability should also be helpful in this area. More specifically, observe that the method proposed in this paper to partition the load in evaluating decomposable predicates, can be applied to non-decomposable ones as well; however in that case communication among the processors is necessary. The question is, how does the amount of necessary communication compare in different partitioning schemes. Finally, we shall mention that we intend to study the relationship between the class of decomposable programs and the programs in the complexity class NC. Also, our notion of program-decomposability may be related to algebraic-operator decomposition, discussed in [IW], and to clausal decomposition, discussed in [LM] (although both papers, in contrast to ours, do not require disjointness of the output sets, and not provide a syntactic characterization of programs). We intend to investigate these possible relationships as well.

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