# **Computing Queries from Derived Relations**\*

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Abstract. Assume that a set of derived relations are available in stored form. Given a query (or subquery), can it be computed from the derived relations and, if so, how? Variants of this problem arise in several areas of query processing. Relation fragments stored at a site in a distributed database system, database snapshots, and intermediate results obtained during the processing of a query are all examples of stored, derived relations. In this paper we give necessary and sufficient conditions for when a query is computable from a single derived relation. It is assumed that both the query and the derived relation are defined by PSJ-expressions, that is, relational algebra expressions involving only projections, selections, and joins, in any combination. The solution is constructive: not only does it tell whether the query is computable or not, but it also shows how to compute it.

## **1. INTRODUCTION**

Consider a database consisting of a number of (conceptual) relations  $R_1, R_2, ..., R_m$ , and assume that the extensions of the conceptual relations are not available in stored form. Instead we have available, in stored form, a set of *n* derived relations, defined by  $E_1, E_2, ..., E_n$  where each  $E_i$  is some expression in relational algebra. We are given a query  $E_0$ , that is, a relational algebra expression over  $R_1, R_2, ..., R_m$ . The problem is then the following: Can  $E_0$  be computed

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Permission to copy without fee all or part of this material is granted provided that the copies are not made or distributed for direct commercial advantage, the VLDB copyright notice and the title of the publication and its date appear, and notice is given that copying is by permission of the Very Large Data Base Endowment. To copy otherwise, or to republish, requires a fee and/or special permission from the Endowment. from the available, stored relations defined by  $E_1, E_2, ..., E_n$  and, if so, how?

This paper represents a first step towards the solution of this problem. We give necessary and sufficient conditions for deciding whether  $E_0$  can be computed from a single derived relation  $E_1$ , for the case when  $E_0$  and  $E_1$  are both *PSJ*-expressions. A relational algebra expression is a *PSJ*-expression if it involves only the operations project, select, join and Cartesian product. This may seem a very restricted class of expressions, but that is not the case. *PSJ*-queries (queries that can be represented by a *PSJ*-expression) are extremely common in relational database systems. Whatever the user query language, almost every query is either a *PSJ*-queries.

Different variants of this problem occur in several areas of query processing. The differences arise from different interpretations of stored, derived relations. In a distributed database context, derived relations can be interpreted as relation fragments stored at a site (or group of sites). This variant of the problem has been studied extensively, but normally under the assumption that each fragment is derived from a single relation using only selections and projections [CP84]. Database snapshots are another example of stored, derived relations [AL80].

The solution to this problem also has applications in "traditional" query optimization [MA83,UL82]. Here a derived relation can be interpreted as an intermediate result obtained in the process of computing a query. If some other part of the same query can be "easily" computed from intermediate results  $E_1, E_2, \dots, E_n$ , query processing may be speeded up. It may even be worthwhile retaining certain intermediate results, if the chance that they can be used in subsequent queries is high enough to warrant the extra storage space [FS82]. In this context the problem has been studied (usually) under the restriction that  $E_0$ must exactly match one of the expressions  $E_1, E_2, ..., E_n$ .

Our main motivation for studying this problem stems from a different area: physical database design for relational databases. In current relational systems

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the structure of the stored database is normally required to be in one-to-one correspondence with the conceptual schema. By this we mean that each conceptual relation exists as a separate stored relation (file). To speed up query processing some auxiliary access structures may be added: secondary indexes, join indexes, record linking, etc. [AS76,BL81,SW76]. This way of structuring the stored database is a simple and straightforward solution. However, it has the effect that the processing of a query almost always requires data to be collected from several physical files.

Instead of directly storing each conceptual relation, we suggest a more flexible approach where the physical database is structured as a set of stored, derived relations. The choice of stored relations should be guided by the (actual or anticipated) query load, so that frequently occurring queries can be processed rapidly. If it is advantageous to do so, some data may be redundantly stored in several physical relations. The structure of the stored database, and changes to the structure, should be completely transparent at the user level, and user queries and updates expressed solely in terms of conceptual relations. The system must be capable of automatically transforming a user query into an equivalent query against stored relations, and a user update to updates of (one or more) stored relations. To make the suggested approach viable the two fundamental problems of query transformation and update transformation must be solved. This paper is a first contribution towards the solution of the query transformation problem.

The idea of not necessarily having a one-to-one correspondence between conceptual relations and stored relations has so far received little attention (outside the area of distributed databases). Joining relations is often the most costly operation in query processing. Schkolnick and Sorensen [SS81] studied the effects of storing relations in prejoined form (called denormalization). Motivated by the availability of special hardware capable of rapid selection and projection, Babb proposed a "Joined Normal Form" [BA82]. It amounts to storing the whole database as one single relation, which is the join of all conceptual relations. Roussopoulos [RO82] originated the idea of view indexing. A view index represents a materialized user view, containing pointers to the tuples contained in the view. A view index is essentially an indirect version of a derived relation.

A classical problem related to the one studied here is that of query containment: given two queries defined by relational expressions  $E_0$  and  $E_1$ , is the result of  $E_0$  a subset of the result of  $E_1$ ? This problem, and the closely related problem of query equivalence, have been solved for certain classes of queries: conjunctive queries and tableau queries. For an overview and references, see the chapters on query optimization in [MA83] or [UL82]. Note, however, that the problem studied here is *not* equivalent to that of query containment. In addition to containment, we impose the further requirement that it must be possible to compute the result of  $E_0$  from the result of  $E_1$ .

# 2. NOTATION AND BASIC ASSUMPTIONS

A relation scheme is a list of attribute names. A database scheme is a set of relation schemes. An instance (or extension) of a relation is a set of mappings which maps a set of attribute names to values or, equivalently, a set of tuples. An instance (extension) of a database is a collection of instances of all relations in the database schema. When there is no risk of confusion we will use the shorter "relation" instead of "instance of a relation", and correspondingly for database.

Derived relations and queries are here expressed in relation algebra. The following notation will be used:

| $R \times S$          | Cartesian product of rela-<br>tions R and S.  |
|-----------------------|---|
| <i>R</i> { <i>C</i> } | Select all tuples from relation $R$ satisfying condition $C$ (a boolean expression).            |
| $R[A_1,,A_m]$         | <b>Project relation</b> $R$ onto attributes $A_1, \dots, A_m$ .                                 |
| R   C   S             | Join relations R and S<br>over condition C, that is,<br>$R \mid C \mid S = (R \times S) \{C\}.$ |
| α(C)                  | The set of all attributes appearing in condition C.   |

The logical connectives will be denoted by + for OR, juxtaposition (multiplication) for AND and  $\neg$  for NOT. A tuple *t* satisfies a condition *C*, if *C* evaluates to true when each attribute name in *C* is replaced by the corresponding attribute value from *t*. An instance of a relation satisfies *C* if all tuples in the relation satisfy *C*. Projections and selections are assumed to have higher precedence than binary operations.

A relational algebra expression has relation names as operands and relational algebra operators as operators. If each relation name is replaced by a corresponding instance the expression is evaluated and the result is a derived relation. The result of evaluating an expression E over an instance d of database  $D = \{R_1, R_2, ..., R_n\}$  will be denoted by V(E, d). An expression is considered to be defined over the full database scheme, even though the expression may only refer to a subset of the relations in the scheme.

We state the following without proof: Every (valid) expression constructed from an arbitrary number of Cartesian products, joins, selections and projections (*PSJ*-expression) can always be transformed to an equivalent expression in a standard form consisting of a Cartesian product, followed by a selection, followed by a final projection. One can easily see that this holds by considering the query tree corresponding to a *PSJ*-expression. The standard form is obtained by first pushing all projections to the root of the tree and thereafter all selection and join conditions.

From the above it follows that any PSJ-expression E can be written in the form  $E = \{R_1 \times R_2 \times \cdots \times R_k\} \{C\} [A_1, A_2, \dots, A_l]$ where  $R_1, R_2, \dots, R_k$  are relations, C is a selection condition and  $A_1, A_2, ..., A_l$  are the attributes of the final projection. We can therefore represent any PSJ-expression by a triple  $E = (\mathbf{A}, \mathbf{R}, C)$  where  $\mathbf{A} = \{A_1, A_2, \dots, A_l\}$  is called the attribute set,  $\mathbf{R} = \{R_1, R_2, ..., R_k\}$  the relation set or base, and C the selection condition.

**Example:** Consider the following expression defined over relations R(A, B, C, D) and S(E, F):

$$((R[A,B] \{A > 10\}) | B = F | (S\{E > 20\})) [A,F]$$
  
= (R × S) {(A > 10) (B = F) (E > 20)} [A,F]  
= ({A,F}, {R,S}, (A > 10) (B = F) (E > 20)).

The triple representation inherits the naming problem of relational algebra: if a relation R occurs more than once in the relation set, then a reference to one of its attributes in C or A is ambiguous. The ambiguity can be resolved by appropriate renaming of repeated relation. This problem will be discussed later, but until further notice we will assume that no relation occurs more than once in the relation set of an expression.

## 3. DERIVABILITY AND COVERAGE

A user query is expressed in terms of conceptual relations. Query transformation consists of transforming a user query into an equivalent query against stored relations. This process should be done automatically by the system and be completely transparent at the user level. Once an equivalent query has been constructed, the task of "optimizing" its evaluation still remains. This problem has been studied extensively and all the known techniques apply [MA83,UL82]. We are not at this stage concerned with constructing efficient queries; the only goal at this stage is to be able to correctly transform queries.

Both stored, derived relations and queries are represented by relational algebra expressions. Stated in its most general form, query transformation amounts finding to а relational algebra expression  $F(E_1, E_2, ..., E_n)$  which generates the same result as a given query  $E_0$ , for every instance of the database. Solving the problem in its general form appears extremely hard. We will reduce it to a more manageable level by imposing a certain restriction on the deriving expression F. The essence of the restriction is that "self-joins" are not allowed in F. A "self-join" is an expression of the type  $E_k | C | E_k$  or  $E_k \times E_k \{C\}$ . In our opinion this is not a severe restriction; "selfjoins" appear to be rare and of limited use.

Definition. (Derivability) Consider a set of relational algebra expressions  $\{E_0, E_1, \dots, E_n\}$  defined over a  $D = \{R_1, R_2, ..., R_m\},$  and let database schema  $\{E_{i_1}, E_{i_2}, \dots, E_{i_k}\}, k \le n$ , be a subset of  $\{E_1, E_2, \dots, E_n\}$ . We say that  $E_0$  is (algebraically) derivable from  $\{E_1, E_2, \dots, E_n\}$ , over the base  $E_{i_1}, E_{i_2}, \dots, E_{i_n}$ , if there exists a relational algebra expression F, containing no joins or Cartesian products, such that  $V(E_0,d) = V(F(E_{i_1} \times E_{i_2} \times \cdots \times E_{i_k}),d)$ for every instance d of the database D.  $\Box$ 

Note that the definition does not completely prohibit joins or Cartesian products; it only forbids "self-joins". There are no restrictions on the use of projections, selections and the standard set operations. An output tuple of F may be constructed by combining any tuple from  $E_j$  with any tuple from  $E_l$  provided that  $j \neq l$ . However, combining a tuple from  $E_j$  with another one from  $E_j$  is not allowed. Note also that F may not be unique. If the database contains redundant data, there may be many different ways of transforming the query.

If the relations defined by  $E_1, E_2, ..., E_n$  do not contain "sufficient" data to construct  $E_0$ , no deriving expression F can exist. If we can show that "sufficient" data exist, then we can proceed to the more difficult problem of constructing a deriving expression F. Saying that  $E_1, E_2, ..., E_n$  contains "sufficient" data vaguely means that "all the data values of  $E_0$  can be found in  $E_1, E_2, ..., E_n$ ". More precisely it means that any value (combination of values) of an attribute (set of attributes) found in a tuple in  $E_0$  can also be found, under the same attributes, by taking some combination of tuples from  $E_1, E_2, ..., E_n$ . Consider the following example:

**Example:**  $E_0$  is defined over  $R_1$  and  $R_2$ , and its attributes are  $[R_1.A, R_1.B, R_2.C]$ . Assume that a tuple (10, 5, -3) exists in  $E_0$  for some instance of  $R_1$  and

 $R_2$ . If the value 10 is not contained in any projection of  $E_1, E_2, ..., E_n$  over  $R_1.A$  for the same instance of  $R_1$ and  $R_2$ ,  $E_0$  cannot be derived from  $E_1, E_2, ..., E_n$ . The same holds if the tuple (10, 5) cannot be found in any projection over  $[R_1.A, R_1.B]$ .  $\Box$ 

This idea is formalized in the following two definitions.

**Definition.** Let  $r = (a_1, a_2, ..., a_m)$  and  $s = (b_1, b_2, ..., b_n), m \le n$ , be two tuples of relations  $R(A_1, A_2, ..., A_m)$  and  $S(B_1, B_2, ..., B_n)$ , respectively. Tuple r is said to be a **subtuple** of s if for every  $a_i$ ,  $1 \le i \le m$ , there exists an attribute  $j, 1 \le j \le n$ , such that  $a_i = b_j$  and  $A_i = B_j$ , and no two attributes in r map to the same attribute in s.  $\Box$ 

**Definition.** (Coverage) Let  $\{E_0, E_1, E_2, ..., E_n\}$  be a set of relational expressions defined over a common database schema  $D = \{R_1, R_2, ..., R_m\}$  and let  $\{E_{i_1}, E_{i_2}, ..., E_{i_k}\}$  be a subset of  $\{E_1, E_2, ..., E_n\}$ . We say that  $E_0$ , is **covered** by  $E_1, E_2, ..., E_n$ , over the base  $\{E_{i_1}, E_{i_2}, ..., E_{i_k}\}$ , if for every instance d of D and every tuple  $t_0 \in V(E_0, d)$  there exists a tuple  $t \in V(E_{i_1} \times E_{i_2} \times \cdots \times E_{i_k}, d)$  such that  $t_0$  is a subtuple of t.  $\Box$ 

It is easy to see that derivability implies coverage. Coverage is a necessary, but not sufficient, condition for derivability. Testing coverage is much easier than testing derivability as we will see. If the coverage test fails, there is no need to consider derivability.

# 4. TESTING THE VALIDITY OF BOOLEAN EXPRESSIONS

We will in the sequel frequently need an algorithm for testing whether a given boolean expression is valid, that is, always evaluates to true. Such an algorithm is developed in this section, for formulas containing no arithmetic expressions. It is based on the same idea as an algorithm given by Rosenkrantz and Hunt [RH80]. Our algorithm is faster but more restricted than theirs.

We will consider boolean expressions constructed from variables, constants, comparison operators (<, =, >), boolean connectives and parentheses. An **atomic conditions** is an expression of the form  $x \operatorname{cop} y$ where x is a variable, cop is one of '<', '=', or '>'and y is either a variable or a constant. If y is a constant, the condition is said to be a **simple condition**, otherwise a **connective condition**. Negation is not needed; any expression containing negation can be reduced to one without negation. Without loss of generality, all variables are assumed to be defined over some finite set of integer values. Variables correspond to attributes and in practice all attributes have a discrete, finite domain. Any discrete, finite domain can be mapped into a finite set of integer values.

Let  $C(x_1, x_2,..., x_n)$  be a boolean expression, as explained above, with variables  $x_1, x_2,..., x_n$ , that is, a formula in first order logic. A formula is valid if  $\forall x_1 \forall x_2 \cdots \forall x_n (C(x_1, x_2,..., x_n))$  holds, that is, if it evaluates to true for all possible values of  $x_1, x_2,..., x_n$ . Every  $x_i$ , i = 1, 2, ..., n is restricted to its domain, of course. When there is no risk of confusion we will use the shorter notation  $\forall x_1, x_2, ..., x_n(C)$ .

From elementary logic we know that  $\forall x_1, x_2 \cdots x_n(C) \Leftrightarrow \nexists x_1, x_2 \cdots x_n(\neg C)$ . Now expand  $\neg C$  into disjunctive form, that is, let  $\neg C$   $= B_1 + B_2 + \cdots + B_m$  where each  $B_i$  is of the form  $B_i = B_{i1}B_{i2} \cdots B_{ik_i}$  and each  $B_{ij}$  is an atomic condition. The quantifiers can be distributed over the  $B_i$ 's, and we obtain the following equivalences:

$$\forall x_1, x_2, \dots, x_n(C) \Leftrightarrow \nexists x_1, x_2, \dots, x_n(\neg C) \Leftrightarrow \nexists x_1, x_2, \dots, x_n(B_1 + B_2 + \dots + B_m) \Leftrightarrow (\nexists x_1, x_2, \dots, x_n(B_1))(\nexists x_1, x_2, \dots, x_n(B_2)) \dots (\nexists x_1, x_2, \dots, x_n(B_n)) .$$

To prove the validity of C it is sufficient to prove, for each  $B_i$  separately, that  $B_i$  is inconsistent (no value exists such that  $B_i$  evaluates to true). If, for any one of the  $B_i$ 's, we can find a value  $x'_1, x'_2, ..., x'_2$  such that  $B_i$  is true, then it immediately follows that C is not valid. Finding a value that satisfies all conditions in  $B_i$ , or proving that no such value exists, can be done by the algorithm explained below.

Let B be a conjunctive condition with variables  $x_1, x_2, ..., x_n$ . The first step of the algorithm is to define for each variable a permissible range which is consistent with its domain and the simple conditions of B. Consider a variable  $x_i$  and let its domain be  $\{L_{x_i}, L_{x_i} + 1, ..., H_{x_i}\}$ . Denote its permissible range by  $r_{x_i} = (a_{x_i}, b_{x_i})$ .

For each variable  $x_i$  of B we initialize its range to  $r_{x_i} = (L_{x_i}, H_{x_i})$ . Then we adjust these initial ranges by considering each simple condition of B in turn. Let w denote a simple condition involving  $x_i$ . The permissible range of  $x_i$  is then adjusted as follows:

(i) if  $w = (x_i > c)$  then  $r_{x_i} := (\max(a_{x_i}, c+1), b_{x_i})$ (ii) if  $w = (x_i < c)$  then  $r_{x_i} := (a_{x_i}, \min(b_{x_i}, c-1))$ (iii) if  $w = (x_i = c)$  then  $r_{x_i} = (c, c)$ .

At the end of this step we have for each  $x_i$  a permissible range which is consistent with its domain and the

simple conditions. If the range of any one variable  $x_i$  is empty then the condition *B* is (trivially) inconsistent. A range  $r_{x_i} = (a_{x_i}, b_{x_i})$  is empty if  $a_{x_i} > b_{x_i}$ .

Example: Consider the conjunctive condition

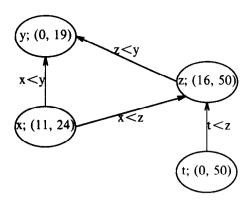
$$B' = (x > 10)(y < 20)(z > 15)(x < 25)(x < y)$$
$$(z < y)(x < z)(t < z)$$

where x, y, z are restricted to the domain  $\{0, 1, ..., 50\}$ . After the first step we have the permissible ranges  $r_x = (11, 24), \quad r_y = (0, 19), \quad r_z = (16, 50)$  and  $r_t = (0, 50)$ , so B' is not trivially inconsistent.  $\Box$ 

If the condition B is not trivially inconsistent we proceed to the next step, in which the connective conditions are taken into account. A directed graph is constructed where nodes represent variables and edges represent connective conditions of the form x > y. A node may represent several variables, as explained below. Each node also has an associated permissible range, which initially is the range obtained from the previous step.

We will denote a node by N(v; (a, b)) where v is the set of variables associated with the node and (a,b)is their permissible range. For each variable x in B a node  $N(x; (a_x, b_x))$  is created, where  $(a_x, b_x)$  is the permissible range obtained from the previous step. Next we merge nodes by using connective conditions of the form x = y. Whenever there is a condition x = y, we locate the node in whose variable set x occurs, and correspondingly for y. Denote these two nodes by  $N_x(v; (a, b))$  and  $N_v(u; (c, d))$ , respectively. modified  $N_{\mathbf{r}}$ to Node is  $N_x(v \cup u; (\max(a,c), \min(b,d)))$  and  $N_v$  is eliminated. This step terminates when all connective conditions with equality have been processed. If any node now has an empty range, condition B is inconsistent and the processing of B stops. Otherwise, there is still at least one value that satisfies all simple conditions and all connective conditions with equality: the value obtained by setting each variable equal to the lower bound of the permissible range of the node in which the variable occurs.

In the final step we take into account connective conditions of the form x > y. For each condition x > y we add an edge from the node in whose variable set y occurs to the node in whose variable set x occurs. For our example we obtain the following graph.



For a variable to have a value that satisfies all connective conditions, the value must be consistent with the values of all its predecessors in the graph. From the graph above we see that the lowest possible values are t = 0, x = 11,  $z = \max(16, 11+1, 0+1)$ = 16 and  $y = \max(0, 11+1, 16+1, 11+2, 0+2) = 17$ . This value satisfies all the conditions of B' and consequently we have found one value for which B' evaluates to true. However, if there had been an additional condition t > 20, then the minimum value for y would be  $y = 23 = \max(0, 16+1, 11+2, 21+2)$ . This is not a permissible value for y because the upper bound is 19, and consequently B' would be inconsistent.

Assume first that the graph obtained does not contain any cycles. Finding the minimum permissible value for each variable can then be done by processing the nodes in a certain order. Consider a node and its immediate predecessors in the graph. If the minimum permissible value for each immediate predecessor has been determined, then the minimum permissible value for (all the variables of) the node is 1 plus the highest of the minimum permissible values of all its immediate predecessors. Any lower value will violate at least one of the conditions. Whenever there are no cycles in the graph, we can easily find a processing sequence such that no node is processed before all its predecessors have been processed. First find all nodes that have no immediate predecessors and mark them. Then repeatedly select any node having only marked immediate predecessors, adjust the lower bound of its range and mark it. If at any stage during this process the range of a node t becomes empty, condition B is inconsistent (because there is no value for the variables of tthat satisfies all the conditions). If, when all nodes have been processed, all permissible ranges are nonempty, then there exists at least one combination of values that satisfies all the conditions: the one obtained by setting each variable equal to the minimum value in its permissible range.

Now consider the situation when there are one or more cycles in the graph. Then there is at least one node t which is its own predecessor. Because all the predecessors of a node must be processed before a node can be processed, node t will never be processed. This will in the algorithm lead to a situation where there are still unmarked nodes in the graph, but they all have at least one unmarked (immediate) predecessor. This situation is easily detected. A cycle in the graph can arise only when there is a subset of connecconditions in tive B such that  $(x_{i_1} > x_{i_2})(x_{i_2} > x_{i_3}) \cdots (x_{i_k} > x_{i_1})$ . This set of conditions can never be satisfied and consequently B is inconsistent.

The above process for testing the inconsistency of a conjunctive boolean expression is summarized in the algorithm INCONSISTENT below. It is possible to strictly prove its correctness but we will not include the proof here. However, we think that its correctness is fairly obvious from the way it was constructed.

The validity of a universally quantified boolean expression can then be proved, or disproved, by the simple algorithm VALID-C which follows.

**procedure** VALID-C(C : boolean expression): boolean

 $\{C \text{ is a boolean expression as explained}\}$ in the beginning of this section}

convert  $\neg C$  into disjunctive form, that is,  $\neg C = B_1 + B_2 + \cdots + B_n$ 

```
for each B_i do
    if not INCONSISTENT (B_i)
     then return (false)
   od:
   return (true);
end {VALID-C}
```

# procedure INCONSISTENT (B: boolean expression): boolean

{B must be of the form  $B = C_1 C_2 \cdots C_m$  where each  $C_i$  is an atomic condition. Each variable x in B must have a finite range  $(L_x, H_x)$ .

G: directed graph with nodes having the format N(v;(a,b)) where v is a set of variable names and (a,b) is the permissible range of the variables;

# begin

```
for each variable x in B do
   insert a new node N_x(\{x\}; (L_x, H_x)) into G;
od;
```

for each simple condition  $(x \circ p c)$  in B do find node  $N_x(\{x\}; (a_x, b_x));$ case op of  $<: (a_x, b_x) := (a_x, \min(b_x, c-1));$  $=:(a_x,b_x):=(c,c):$ >:  $(a_x, b_x)$ : =  $(\max(a_x, c+1), b_x)$ ; end; if  $a_x > b_x$  then return (true); od:

for each connective condition (x = y) in B do find nodes  $N_x(v; (a, b))$  and  $N_y(u; (c, d))$ such that  $x \in v$  and  $y \in u$ ;  $v := v \cup u$ :  $(a,b):=(\max(a,c),\min(b,d));$ delete  $N_{\nu}$  from G; if a > b then return (true):

od:

for each connective condition (x > y) in B do find nodes  $N_x(v; (a,b))$  and  $N_y(u; (c,d))$ such that  $x \in v$  and  $y \in u$ ; insert into G an edge from  $N_{\nu}$  to  $N_{\nu}$ ;

od;

while G is not empty do find any node  $N_p(v; (a_n, b_n))$ without incoming edges; if none exists then return (true); {detected a cycle} for each edge s starting from  $N_p$  do let  $N_s(u,(a_s,b_s))$  be the end node of s;  $(a_s, b_s) := (\max(a_s, a_p + 1), b_s)$ if  $a_s > b_s$  then return (true); od delete node  $N_p$  from G; od;

return (false); end {INCONSISTENT};

Testing the validity of a boolean expression is equivalent to the satisfiability problem for boolean expressions. It is well-known that testing satisfiability may, in the worst case, take exponential time. However, we do not consider this to be a very serious problem. Selection expressions are normally short, and the exponential explosion occurs only for expressions involving several "not equals" [RH80].

#### 5. TESTING COVERAGE

We are now ready to consider the question whether a *PSJ*-expression  $E_0$  is covered by another *PSJ*-expression  $E_1$ . Coverage means that every tuple obtained by evaluating  $E_0$  must occur (as subtuples) in the relation obtained by evaluating  $E_1$ , and that this must hold for every instance of the database. Before stating the main theorem we need the notion of an extended attribute set.

**Definition.** Let C be a boolean expression with variables  $x_1, x_2, ..., x_n, y_1, y_2, ..., y_m$ . The variable  $y_i$ ,  $1 \le i \le m$ , is said to be **uniquely determined** by  $x_1, x_2, ..., x_n$ , with respect to C, if the following holds:

$$\forall x_{1},...,x_{n},y_{1},...,y_{m},y_{1}',...,y_{m}'(C(x_{1},...,x_{n},y_{1},...,y_{m}))$$

$$C(x_{1},...,x_{n},y_{1}',...,y_{m}') \Rightarrow y_{i} = y_{i}'). \square$$

Let C' stand for  $C(x_1,...,x_n,y_1',...,y_m')$ .

The above condition is equivalent to  $\neg((y_i > y'_i + y_i < y'_i)CC')$  which can be tested using the algorithm VALID-C. The attribute set of a derived relation can be extended to include all variables uniquely determined by those already in the attribute set.

**Definition.** Let E = (A, R, C) be a *PSJ*-expression and let **B** be the set of all attributes uniquely determined by the attributes of **A**. Then  $A \cup B$ , denoted by A+, is called the **extended attribute set** of E.  $\Box$ 

The set **B** in the definition above must obviously be a subset of  $\alpha(C)$ . An attribute not mentioned in C cannot be uniquely determined by the attributes of **A**. The importance of the notion of an extended attribute set stems from the fact that given a tuple from E we can correctly derive the missing values of all the variables in the set **B**. The procedure is explained below.

Let  $A = \{x_1, x_2, ..., x_n\}$  and  $\alpha(C) - A$ =  $\{y_1, y_2, ..., y_m\}$ . Assume that  $y_1, y_2, ..., y_k$ ,  $k \le m$ , have been shown to be uniquely determined by  $x_1, x_2, ..., x_n$ . We are given a tuple  $t = (x_1^0, x_2^0, ..., x_n^0)$ from the relation defined by E. The values of the variables  $y_1, ..., y_m$  are not known, but they must have been such that the tuple t, before the projection, satisfied the condition C. In other words, the following must hold:  $\exists y_1, y_2, ..., y_m(C(x_1^0, ..., x_n^0, y_1, y_2, ..., y_m))$ .

This is equivalent to  $\neg(\forall y_1, y_2, ..., y_m(\neg C))$ . The algorithm VALID-C is now applied to  $\forall y_1, ..., y_m(\neg C)$ , modifying it to return the values found. Denote the values returned by  $y_1^0, y_2^0, ..., y_m^0$ . Such a value combination always exists, because otherwise t would not have satisfied C in the first place. The values for the

variables  $y_1, y_2, ..., y_k$  are guaranteed to be unique, because each variable  $y_1, y_2, ..., y_k$  is uniquely determined by  $x_1, x_2, ..., x_n$ . In other words, we have correctly reconstructed the missing values of  $y_1, y_2, ..., y_k$ . The following examples illustrates the discussion.

| Example:  | Let        |
|---|------------|
| $E = (\{A\}, R, (B=A)((A>7)(C=5) + (A<5)(C=10))$    |            |
| where $R = (A, B, C, D)$ . One can easily show that | t <b>B</b> |
| and $C$ (but not $D$ ) are uniquely determined by   | Α.         |
| Now consider the following instance of $R$ and      | the        |
| result when $E$ is applied to that instance:        |            |

| <b>R</b> : | A  | B  | С  | D | <i>E</i> : | A  |
|------------|----|----|----|---|------------|----|
|            | 10 | 10 | 5  | 4 |            | 10 |
|            | 10 | 11 | 5  | 6 |            | 4  |
|            | 11 | 11 | 7  | 4 |            |    |
|            | 4  | 4  | 10 | 6 |            |    |
|            | 3  | 3  | 14 | 6 |            |    |
|            |    |    |    |   |            |    |

Each tuple appearing in the result of E must have satisfied the selection condition. For the first tuple in E we can conclude that B = 10 (because of the condition B = A) and that C = 5 (because A > 7). In the same way, we find for the second tuple that B = 4 and C = 10.  $\Box$ 

**Lemma.** Let  $E = (\mathbf{A}, \mathbf{R}, C)$  be a PSJ-expression. Given a tuple from E, the value of an attribute y, is guaranteed to be uniquely reconstructible if and only if  $y \in \mathbf{A} + . \square$ 

Sketch of proof: The procedure above for reconstructing missing values shows that the condition  $y \in \mathbf{A} +$  is sufficient. To show that it is also necessary, assume that  $y \notin A +$ . Then we can find two tuples  $t_0$  and  $t_1$  over the base of E (the Cartesian product of all relations in **R**) such that they agree on all attributes in A but differ at least on the attribute y. This follows directly from the definition of "uniquely determined" above. The database instance obtained by projecting  $t_0$  and  $t_1$  onto the attributes of each relation in **R** is a valid instance. When  $t_0$  and  $t_1$ are projected onto the attributes of A they will map to the same tuple, because they agree on all attributes of A. Given only this tuple from E (and the condition C) the value of y cannot be uniquely determined. There are at least two different values of y, which when combined with the given tuple satisfy C.  $\Box$ 

Given a tuple in E, we know the values of all attributes in A and we can correctly determine the values of all other attributes in A+. The above lemma shows that A+ is the maximal set of attributes whose values can be determined from a tuple in E. The following theorem gives necessary and sufficient conditions on coverage. Note the requirement that all instances of the relations involved must be non-empty.

**Theorem 1.** Consider two PSJ-expressions  $E_0 = (\mathbf{A}_0, \mathbf{R}_0, C_0)$  and  $E_1 = (\mathbf{A}_1, \mathbf{R}_1, C_1)$  defined over a database where every relation in  $\mathbf{R}_0 \cup \mathbf{R}_1$  is nonempty. Let  $\alpha(C_0) \cup \alpha(C_1) = \{x_1, x_2, \dots, x_n\}$ . Then  $E_0$  is covered by  $E_1$  if and only if  $\mathbf{A}_0 \subseteq \mathbf{A}_1 +$  and  $\forall x_1, x_2, \dots, x_n (C_0(x_1, x_2, \dots, x_n) \Rightarrow C_1(x_1, x_2, \dots, x_n))$ .  $\Box$ 

Sketch of Proof: First extend the relation set of  $E_0$  and  $E_1$  to  $\mathbf{R}_0 \cup \mathbf{R}_1$  so that they are defined over a common base (Cartesian product). Adding a new relation S to the relation set of an expression does not change the result of the expression (as long as the new relation always contains at least one tuple). A tuple t in the base before adding S will give rise to a number of "copies" in the extended base, where the copies differ only in the values of the attributes of S. If tsatisfied a given selection condition, then all its "copies" will satisfy the condition as well. If t did not satisfy the condition, then none of the "copies" will satisfy it. The added attributes have no effect on the selection. The final projection will reduce all "copies" of t to a single tuple, exactly the same one as the projection of t.

Assume that  $A_0 \subseteq A_1 +$  and that the implication holds. The fact that the implication holds, guarantees that each tuple t in the common base satisfying  $C_0$ , will also satisfy  $C_1$ . Because  $A_0 \subseteq A_1 +$ , the projection of t onto  $A_0$  is a subtuple of the projection of t onto  $A_1 +$ . Consequently  $E_0$  is covered by  $E_1$ .

It is obvious that the condition  $A_0 \subseteq A_1 + is$ necessary for coverage. Now assume that  $E_0$  is covered by  $E_1$  and that  $A_0 \subseteq A_1 +$ , but that the implication does not hold. Then there must exist a tuple t such that it satisfies  $C_0$  but not  $C_1$ . Now consider the database instance where each relation  $S_i \in \mathbf{R}_0 \cup \mathbf{R}_1$  consist of the single tuple obtained by projecting t onto the attributes of  $S_i$ . When  $E_0$  is evaluated over this instance, the result will be one tuple: the projection of t onto  $A_0$ . However, when evaluating  $E_1$  over the same database instance the result will be the empty set. Hence  $E_0$  is not covered by  $E_1$  for this instance, contradicting the assumption that  $E_0$  is covered by  $E_1$  for every instance of the database.  $\Box$ 

By the above theorem testing coverage is reduced to two simpler tests. Note that it is not always necessary to construct  $A_1 + .$  If  $A_0 \subseteq A_1$  the first condition is trivially satisfied. If this is not the case, then we must test whether every variable in  $A_0 - A_1$  is uniquely determined by the attributes in  $A_1$ . Let x be a variable in  $A_0 - A_1$ . A first simple test is to check whether x is in  $\alpha(C_1)$ . If x is not mentioned in  $C_1$ , then it cannot be uniquely determined by those in  $A_1$ .

Note that Theorem 1 can be used to test equivalence of relational expressions for a wider class of expressions than those considered in [AS79] and [SY81].

#### 6. TESTING DERIVABILITY

The fact that an expression  $E_0$  is covered by a derived relation  $E_1$  does not automatically guarantee derivability. All the necessary tuples are guaranteed to exist in  $E_1$ , but it may be impossible to select out the ones belonging to  $E_0$ , as illustrated by the following example.

**Example:** Consider a database consisting of the single relation S(A,B,C). Let  $E_0 = (\{A\},\{S\},(B>10)(C<20))$  and  $E_1 = (\{A,C\},\{S\},B>5)$ . It is easy to see that  $E_0$  is covered by  $E_1$  because every tuple satisfying (B>10)(C<20) will automatically satisfy B>5. However, to compute  $E_0$  from the derived relation  $E_1$ , attribute B is needed (to further qualify the tuples) and B is not present in  $E_1$  after the projection. Hence  $E_0$  cannot be derived from  $E_1$ .  $\Box$ 

Consider  $E_0 = (\mathbf{A}_0, \mathbf{R}_0, C_0)$  and  $E_1 = (\mathbf{A}_1, \mathbf{R}_1, C_1)$ and assume that  $E_0$  is covered by  $E_1$ . If  $\alpha(C_0) \subseteq \mathbf{A}_1 +$ , that is, if all the attributes needed to evaluate the condition  $C_0$  are present in  $\mathbf{A}_1 +$ , it is obvious that  $E_0$  can be derived from  $E_1$ . However, even when some attributes in  $\alpha(C_0)$  are missing from  $\mathbf{A}_1$ + it may still be possible to derive  $E_0$  from  $E_1$ . The missing attributes may not be needed. In the example above that would be the case if the condition in  $E_0$  were B > 5 instead of B > 10. To handle this type of situation we introduce the notion of essential and nonessential variables.

**Definition.** Consider a boolean expression C with variables  $x_1, x_2, ..., x_n$ . The variable  $x_k$  is said to be **nonessential** if the following holds

$$\forall x_1, ..., x_k, ..., x_n, x_k'(C(x_1, ..., x_k, ..., k_n)) = C(x_1, ..., x_k', ..., x_n)).$$

Otherwise  $x_k$  is an essential variable of C.  $\Box$ 

A nonessential variable can be eliminated from the condition simply by replacing it with any value from its domain. This will in no way change the value of the expression. However, for derivability we need to know whether a variable in a condition is essential given that another condition holds. This is covered by the next definition.

**Definition.** Let  $C_0$  and  $C_1$  be boolean expressions with variables  $x_1, x_2, ..., x_n$ . The variable  $x_k$  is said to be computationally nonessential in  $C_0$  with respect to  $C_1$  if the following condition holds:

$$\begin{aligned} &\forall x_1, \dots, x_k, \dots, x_n, x_k' \\ & (C_1(x_1, \dots, x_k, \dots, x_n) C_1(x_1, \dots, x_k', \dots, x_n) \\ &\Rightarrow C_0(x_1, \dots, x_k, \dots, x_n) = C_0(x_1, \dots, x_k', \dots, x_n)) . \end{aligned}$$

The basic idea of this definition is that the exact value of  $x_k$  does not matter when evaluating  $C_0$ . It can be replaced by any value  $x_k'$  as long as the value  $x_k'$  is such that  $C_1$  is satisfied. The algorithm VALID-C in section 3 can be used to test whether the condition holds. Let  $C_1'$  stand for  $C_1(x_1,...,x_k',...,x_n)$  and similarly for  $C_0'$ . Then the condition is equivalent to

$$\forall x_1, \dots, x_k, \dots, x_n, x_k'$$
  
( $\neg (C_1 \ C_1' (C_0 + C_0') (\neg C_0 + \neg C_0'))) ,$ 

to which the algorithm can be applied directly.

Both definitions above can trivially be extended to the case of several nonessential variables by interpreting  $x_k$  (and  $x_{k'}$ ) as sets of variables. Note also that a variable not occurring in  $C_0$  is (trivially) nonessential.

If a variable (or set of variables) y of  $C_0$  is computationally nonessential with respect to  $C_1$ , we can correctly evaluate  $C_0$  without knowing the exact value of y. However, y cannot be assigned an arbitrary value in its domain. Consider two PSJ-expressions  $E_0 = (\mathbf{A}_0, \mathbf{R}_0, C_0)$  and  $E_1 = (\mathbf{A}_1, \mathbf{R}_1, C_1)$  where  $E_0$  is covered by  $E_1$ . Let  $\alpha(C_0) = \{x_1, x_2, \dots, x_n, y\}$  where  $\{x_1, \dots, x_n\} \subseteq \mathbf{A}_1 + \text{ and } y \notin \mathbf{A}_1 + \dots$  Assume that we have shown that y is computationally nonessential with respect to  $C_1$ . Then y must be one of the variables in  $C_1$ . Now consider an arbitrary but fixed tuple  $t_1 = (x_1', x_2', ..., x_n', y', ...)$  from  $E_1$  before the projection onto  $A_1$ . After the projection y disappears and its exact value in  $t_1$  is not known. However, we know that it was such that  $t_1$  satisfied  $C_1$ . If we somehow can find a surrogate value y'' such that the tuple  $t_2$  $= (x_1', x_2', ..., x_n', y'', ...)$  satisfies  $C_1$ , then from the fact that y is computationally nonessential it follows that  $C_0(x_1', x_2', ..., x_n', y') = C_0(x_1', x_2', ..., x_n', y'').$  In other words, using the surrogate value y'' when evaluating  $C_0$  gives exactly the same result as would have been obtained using the correct, but unknown, value y'.

**Example:** Let R(A, B, C, D),  $E_0 = (\{A, B\}, \{R\}, (A < 50)(B > 10)(C < 40)(D < C)(D > B))$ and  $E_1$   $= (\{A, B, C\}, \{R\}, (B > 10)(C < 40)(D < C)(D > B) + (B < 5)).$   $E_1$  covers  $E_0$  and it is easy to show that D is computationally nonessential in  $C_0$ . Now consider the following instance of the relation defined by  $E_1$ :

| 1: | A   | B  | С  |
|----|-----|----|----|
|    | 100 | 30 | 35 |
|    | 20  | 15 | 25 |
|    | 20  | 4  | 0  |

Ε

For the first tuple we can deduce that the missing value of D must have been in the range 30 < D < 35. We can set D to any value in this range, D = 31, for example. Using D = 31 and the given values for the other attributes, we find that the first tuple does not satisfy the condition in  $E_0$ . For the second tuple the range is 15 < D < 25. Setting D = 16 we find that the second tuple satisfies the selection condition of  $E_0$ . For the third tuple, D can be assigned any value in its range, D = 0, for example. (The tuple satisfies the condition B < 5 independently of the value of D.) With D = 0 (or any other value) the tuple does not satisfy the condition in  $E_0$ . The final result of  $E_0$  is hence the single tuple (20, 15).  $\Box$ 

Surrogate values for computationally nonessential variables can easily be found using the algorithm given in section 3. Let  $A_1 = \{x_1, x_2, \dots, x_n\}$  and let  $s = (x_1', ..., x_n')$  be a tuple in the relation defined by  $E_1$ . This tuple must then have satisfied  $C_1$ . Let  $\alpha(C_1) - \mathbf{A}_1 = \{y_1, y_2, \dots, y_k\}$ , that is, the variables of  $C_1$ not retained after the projection. All computationally nonessential variables of  $C_0$  are guaranteed to be in the set  $\{y_1, \dots, y_k\}$ . Finding surrogate values for  $y_1, y_2, \dots, y_k$  to be used with the given tuple proving  $s = (x_1', x_2', \dots, x_n')$ amounts to  $\exists y_1, y_2, \dots, y_k (C_1(x_1', x_2', \dots, x_n', y_1, y_2, \dots, y_k)).$  Proving this equivalent to disproving is  $\forall y_1, y_2, ..., y_k (\neg C_1(x_1', ..., x_n', y_1, ..., y_k))$  which can be done by applying the algorithm of section 3 (modifying it to return the value found). Surrogate values  $y_1', y_2', \dots, y_k'$  will always exist, because otherwise s could not have satisfied  $C_1$ . It may be necessary to compute new surrogate values for every tuple, as we saw in the example above. However, when given a new tuple, it may be worthwhile trying to use previously computed surrogate values. If the attribute values of the new tuple combined with the 'old' surrogate values satisfy  $C_1$ , then there is no need to recompute the surrogates.

The following theorem summarizes the main result of this paper:

**Theorem 2.** Let  $E_0 = (\mathbf{A}_0, \mathbf{R}_0, C_0)$  and  $E_1 = (\mathbf{A}_1, \mathbf{R}_1, C_1)$  be PSJ-expressions and assume that  $E_0$  is covered by  $E_1$ . Then  $E_0$  is derivable from  $E_1$  if and only if all variables in  $\alpha(C_0) - \mathbf{A}_1 + are$  computationally nonessential with respect to  $C_1$ .  $\Box$ 

**Sketch of Proof:** If the condition holds then  $E_0$ can be computed in the way discussed above, so the condition is clearly sufficient. To show that it is necessary, assume that  $\alpha(C_0) - A_1 + is$  nonempty, that  $y \in \alpha(C_0) - A_1 +$  is a computationally essential variable, and that  $E_0$  is derivable from  $E_1$ . Then there must exist a relational algebra expression F, not involving the missing attribute y, which correctly computes  $E_0$  from  $E_1$  for every instance of the database. Denote by S the base of  $E_1$ , that is, the Cartesian product of all relations in  $\mathbf{R}_1$ . Because y is computationally essential, we can construct two tuples  $t_1$  and  $t_2$ over S which differ only in the value of y and which both satisfy  $C_1$  but only one,  $t_1$  say, satisfies  $C_0$ . Now consider the following two database instances:  $d_1$ obtained by projecting  $t_1$  onto the relations in  $\mathbf{R}_1$  and  $d_2$  by similarly projecting  $t_2$ .  $E_0$  evaluated over  $t_1$ must then result in one tuple, while  $E_0$  evaluated over  $t_2$  must be empty. However, evaluating F over  $t_1$ gives exactly the same result as evaluating F over  $t_2$ , because they differ only in the value of y and the expression F does not involve the attribute y. Hence, F gives an incorrect result for one of the two database instances, contradicting the assumption that  $E_0$  is derivable from  $E_1$ .  $\Box$ 

Note that the proof does not impose any other restrictions on the deriving expression F than that it does not involve attribute y. This implies that the theorem holds also for expressions that involve "self-joins". In other words, if  $E_0$  cannot be derived from  $E_1$  then it cannot be derived from  $E_1 \times E_1$  or  $E_1 \times E_1 \times E_1, \ldots$  either.

#### 7. DISCUSSION

In this section we summarize the main assumptions made and briefly discuss the effect of relaxing them.

One of the main restrictions was not to allow "self-joins". A "self-join" in a *PSJ*-expression is indicated by the same relation name occurring more than once in the relation set when the expression is converted into triple format. This restriction can be slightly relaxed. Assume that a relation R occurs k, k > 1, times in the relation set of  $E_0$  and that the k occurrences have been renamed  $R^1, R^2, \dots, R^k$ . If the same relation occurs in  $E_1$  with at least the same mul-

tiplicity,  $E_0$  may be derivable from  $E_1$ . Assume that there are  $l, l \ge k$ , occurrences of R in  $E_1$  and that they have been renamed  $S^1, S^2, ..., S^l$ . If we can find a mapping of the names  $R^1, R^2, ..., R^k$  onto a subset of  $S^1, S^2, ..., S^l$  such that  $E_0$  is derivable from  $E_1$  under the mapping, the problem is solved. (A mapping associates each name  $R^1, R^2, ..., R^k$  with some name  $S^1, S^2, ..., S^l$ .) There does not seem to be any better way of solving this than trying all l(l-1)...(l-k+1) possible mappings.

The algorithm in section 4 works only for a restricted class of boolean expressions: only those where atomic condition are restricted to a comparison of two variables or a comparison of a variable with a constant. Note, however, that none of the results in sections 5 and 6 depend directly on this algorithm. Any class of boolean expressions can be handled provided that we have an algorithm for testing the validity (or satisfiability) of expressions of that class. The core of such an algorithm is an algorithm for testing whether a set of inequalities and/or equalities can all be simultaneously satisfied. Whether such an algorithm exists or not, and its complexity, depend completely on the type of expressions (functions) allowed in the (in)equalities and the domain of the variables. If linear expressions with variables ranging over the real numbers are allowed, the problem is equivalent to findingng a feasible solution to a linear programming problem.

The reader may have noticed that so far we have not mentioned keys, functional dependencies, inclusion dependencies and so on. What effect will they have on derivability? The only constraints taken into account in this paper are domain constraints: any combination of attribute values drawn from their respective domains represents a tuple that potentially may occur in an instance of the relation in guestion. The only effect of functional dependencies, multivalued dependencies, inclusion dependencies, etc. is to impose constraints on what tuples can occur simultaneously in the database, thereby reducing the set of permitted database instances. Clearly the conditions given for derivability are still sufficient, but they may not be necessary. The proof that the stated conditions are necessary for derivability relies on a certain database instance, consisting of two tuples, to produce a contradiction. In the presence of additional constraints, this may not be a permissible database instance.

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